

## ***Initial Site Characterization***

Radford High School Track and Field Excavation  
4361 Salt Lake Boulevard  
(Portion of TMK No: [1] 9-9-002: Parcel 023)  
Honolulu, Oahu, Hawaii

March 21, 2014  
Project No. 17012-012148.00/Task 48

*Prepared for:*

**STATE OF HAWAII DEPARTMENT OF EDUCATION**  
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*Move Forward with Confidence*

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## CONTENTS

| <u>Section</u>  | <u>Page</u> |
|---|-------------|
| List of Acronyms.....   | iv          |
| Executive Summary.....  | v           |
| <b>1.0 INTRODUCTION AND PURPOSE.....</b>                          | <b>1</b>    |
| <b>2.0 BACKGROUND.....</b>  | <b>1</b>    |
| 2.1 SITE DESCRIPTION.....   | 1           |
| 2.1.1 Soil Conditions.....  | 1           |
| 2.1.2 Groundwater Conditions.....                                 | 1           |
| 2.2 HISTORICAL SITE USE.....                                      | 2           |
| 2.3 RECENT CONSTRUCTION ACTIVITIES.....                           | 2           |
| <b>3.0 APPLICABLE ACTION LEVELS.....</b>                          | <b>2</b>    |
| <b>4.0 FIELD ACTIVITIES.....</b>                                  | <b>2</b>    |
| 4.1 INITIAL RESPONSE.....   | 2           |
| 4.2 INVESTIGATION PLANNING.....                                   | 4           |
| 4.3 SOIL SAMPLING ACTIVITIES.....                                 | 4           |
| 4.4 FIELD REPLICATE SAMPLING.....                                 | 5           |
| 4.5 CHAIN-OF-CUSTODY.....   | 6           |
| 4.6 DECONTAMINATION PROCEDURES.....                               | 6           |
| 4.7 INVESTIGATION DERIVED WASTE.....                              | 6           |
| <b>5.0 LABORATORY ANALYTICAL RESULTS.....</b>                     | <b>6</b>    |
| 5.1 INITIAL INVESTIGATION SOIL ANALYTICAL RESULTS.....            | 7           |
| 5.1.1 TPH.....  | 7           |
| 5.1.2 SVOCs.....  | 7           |
| 5.1.3 RCRA 8 Metals.....  | 7           |
| 5.1.4 PCBs.....   | 7           |
| 5.1.5 Organochlorine Pesticides.....                              | 7           |
| 5.1.6 Dioxin.....   | 8           |
| 5.2 FOLLOW-UP INVESTIGATION SOIL ANALYTICAL RESULTS.....          | 9           |
| 5.2.1 TPH.....  | 9           |
| 5.2.2 SVOCs.....  | 9           |
| 5.2.3 RCRA 8 Metals.....  | 9           |
| 5.2.4 PCBs.....   | 9           |
| 5.2.5 Organochlorine Pesticides.....                              | 10          |
| 5.2.6 Dioxin.....   | 10          |
| 5.3 LABORATORY QUALITY CONTROL.....                               | 11          |
| 5.4 FIELD QUALITY CONTROL.....                                    | 11          |
| 5.5 PRELIMINARY ENVIRONMENTAL HAZARD EVALUATION - FIELD AREA..... | 13          |
| <b>6.0 SUMMARY AND CONCLUSION.....</b>                            | <b>13</b>   |
| <b>7.0 LIMITATIONS.....</b>                                       | <b>16</b>   |



## **CONTENTS (Continued)**

### **References**

### **Figures**

- 1 Site Location Map
- 2 Initial Sample Area
- 3 Decision Unit Area Results Exceeding HDOH Tier 1 EALs

### **Tables**

- 1 Analytical Results for Initial Soil Sample
- 2 Dioxin Toxic Equivalency (TEQ) Calculations for Initial Soil Sample
- 3 Analytical Results for Soil Samples from Decision Unit Areas
- 4 Analytical Results for Replicate Samples
- 5 Statistical Calculations for Replicate Samples

### **Photographs**

### **Appendices**

- A Laboratory Analytical Reports and Chain of Custody Forms (Test America)
- B Laboratory Analytical Reports and Chain of Custody Forms (NVL)
- C Laboratory Analytical Reports and Chain of Custody Forms (ESN)
- D Asbestos Disposal Manifest



## List of Acronyms

|                |   |
|----------------|---|
| ACM            | Asbestos Containing Material                            |
| bgs            | below ground surface                                    |
| Bureau Veritas | Bureau Veritas North America, Inc.                      |
| COPC           | Constituents of Potential Concern                       |
| DOE            | Department of Education                                 |
| DRO            | Diesel range organics                                   |
| DU             | Decision Unit   |
| EALs           | Environmental Action Levels                             |
| EHE            | Environmental Hazard Evaluation                         |
| EHMP           | Environmental Hazard Management Plan                    |
| EPA            | Environmental Protection Agency                         |
| ESN            | Environmental Services Network Pacific, Inc.            |
| HDOH           | State of Hawaii Department of Health                    |
| HEER           | Hazard Evaluation and Emergency Response                |
| HQ             | Hazard Quotient   |
| IDW            | Investigation Derived Waste                             |
| mg/kg          | milligrams per kilogram                                 |
| mg/L           | milligram per liter                                     |
| ng/kg          | nanograms per kilogram                                  |
| NVLAP          | National Voluntary Laboratory Accreditation Program     |
| PCB            | Polychlorinated Biphenyls                               |
| PLM            | Polarized Light Microscopy                              |
| PPE            | Personal Protective Equipment                           |
| RCRA           | Resource Conservation Recovery Act                      |
| RL             | Reporting Limit   |
| RPD            | Relative Percent Difference                             |
| RRO            | Residual Range Organics                                 |
| RSD            | Relative Standard Deviation                             |
| SOP            | Standard Operating Procedure                            |
| SPLP           | Synthetic Precipitation Leaching Procedure              |
| SVOCs          | Semi-Volatile Organic Compounds                         |
| QA/QC          | Quality Assurance/Quality Control                       |
| TCDD           | tetrachlorodibenzo-p-dioxin                             |
| TEF            | Toxicity Equivalence Factor                             |
| TEQ            | Toxicity Equivalent                                     |
| TGM            | Technical Guidance Manual                               |
| TMK            | Tax Map Key   |
| TPH            | Total Petroleum Hydrocarbons                            |
| TPH-DRO        | Total Petroleum Hydrocarbons as Diesel Range Organics   |
| TPH-RRO        | Total Petroleum Hydrocarbons as Residual Range Organics |
| UIC            | Underground Injection Control                           |
| USGS           | United States Geological Survey                         |
| WHO            | World Health Organization                               |





## Executive Summary

The State of Hawaii Department of Education (DOE) retained Bureau Veritas North America, Inc. (Bureau Veritas) to collect and analyze soil samples and suspect asbestos materials from the track and field area at Radford High School. This report summarizes the data collected at the property described as being located at 4361 Salt Lake Boulevard, Tax Map Key (TMK) Number (1) 9-9-002: Parcel 023 (hereinafter referred to as the "site").

On December 19, 2013 buried debris was encountered during excavation activities at the track and field area of the site. A topographic map from 1953 depicts the site and adjoining areas as a naval reservation located within Makalapa Crater. The topographic maps from 1959 through 2013 showed the site as a built-up area; labeled "Radford High School". The encountered buried debris is assumed to be associated with the site's former use as a military reservation, which may have included a former military dumpsite.

The purpose of this investigation was to assess the presence and magnitude of the soil contamination within the track construction area and associated soil stockpiles.

On December 20, 2013, after the discovery of buried debris, Bureau Veritas responded onsite to collect one multi-increment soil sample and six suspect Asbestos Containing Material (ACM) samples. Two types of suspect ACM, an off-white fibrous material with debris and an off-white/brown fibrous material with debris, were observed. Three samples of each type of suspect ACM were collected and sent to NVL Laboratories, Inc., a National Voluntary Laboratory Accreditation Program (NVLAP) laboratory located in Seattle, Washington. The samples were analyzed for asbestos content utilizing Polarized Light Microscopy (PLM) analysis. The three samples of the first type of suspect ACM (off-white fibrous material with debris) tested positive for asbestos with 51% to 55% Chrysotile. The three samples of the second type of suspect ACM (off-white/brown fibrous material with debris) did not test positive for asbestos. After the samples were collected, MEI personnel covered the buried debris with geofabric. The asbestos was subsequently bagged and stored in a locked storage area. On February 11, 2014, Unitek arrived onsite and double bagged five bags of ACM, which were subsequently properly disposed.

The multi-increment soil sample was a representative sample collected from one decision unit (RHS-01), that was established on the northeastern portion of the site, where buried debris was uncovered. The sample was collected to assess the soil in the track area surrounding the buried debris. The soil sample was analyzed for Total Petroleum Hydrocarbons as Diesel Range Organics (TPH-DRO), Total Petroleum Hydrocarbons as Residual Range Organics (TPH-RRO), Semi-Volatile Organic Compounds (SVOCs), Resource Conservation and Recovery Act (RCRA) 8 Metals, polychlorinated biphenyls (PCBs), Organochlorine Pesticides, and dioxin. Analysis of the soil sample indicated that the soil concentrations of arsenic, cadmium, lead and dioxin exceeded their respective Hawaii Department of Health (HDOH) Tier 1 Environmental Action Levels (EALs). Further sampling and analysis were recommended to delineate the extent of the soil contamination.

A scope of work was prepared, submitted, and approved by the State of Hawaii Department of Health (HDOH) prior to conducting a follow-up investigation. On February 4 and 5, 2014, a follow-up soil investigation was conducted by Bureau Veritas. To facilitate the additional investigation, the site was divided into the following three areas of concern: (1) the excavated track area, (2) the soil stockpiles, and (3) the grassy/original field area. Each of the areas of concern was further divided into DUs based on visual contamination and size. A total of 29 DUs (RHS-DU-1 through RHS-DU-29) were established across the site to facilitate the collection of multi-increment soil samples from the surface soil layer in each DU.



One multi-increment surface soil sample was collected from each DU for a total of 29 samples. Three sets of replicate samples were also collected from three DUs selected to represent each of the three areas of concern. The samples were submitted to Environmental Services Network Pacific (ESN) Analytical Laboratory located in Honolulu, Hawaii for analytical testing and were analyzed for the following:

- TPH-DRO using Environmental Protection Agency (EPA) Method 8015B.
- TPH-RRO using EPA Method 8015B
- SVOCs using EPA Method 8270C
- RCRA 8 Metals using EPA Methods 6010B and 7471
- PCBs using EPA Method 8082
- Organochlorine Pesticides using EPA Method 8081A
- Dioxin using EPA Method 8290

TPH-DRO and TPH-RRO were not detected in the 29 soil samples at concentrations greater than the laboratory reporting limits.

One SVOC analyte, benzo(a)pyrene, was detected at concentrations that equaled or exceeded the HDOH Tier 1 EAL. In three of the 29 samples, benzo(a)pyrene concentrations that ranged between 0.15 and 0.43 milligrams per kilogram (mg/kg) were equal to or exceeded the HDOH Tier 1 EAL for benzo(a)pyrene of 0.15 mg/kg.

Five of the eight RCRA metals, arsenic, barium, cadmium, lead, and mercury, were detected at concentrations that exceeded their respective HDOH Tier 1 EALs. Arsenic was detected in seven of the 29 samples at concentrations ranging from 24 to 35 mg/kg, which were equal to or exceeded the HDOH Tier 1 EAL for arsenic of 24 mg/kg. Barium was detected in one of the samples at a concentration of 1,200 mg/kg, which exceeded the HDOH Tier 1 EAL for barium of 1,000 mg/kg. Cadmium was detected in one of the samples at a concentration of 14 mg/kg, which was equal to the HDOH Tier 1 EAL for cadmium of 14 mg/kg. Lead was detected in 15 of the samples at concentrations ranging from 240 to 6,200 mg/kg, which exceeded the HDOH Tier 1 EAL for lead of 200 mg/kg. Mercury was detected in 22 of the samples at concentrations ranging from 4.8 to 200 mg/kg, which exceeded the HDOH Tier 1 EAL for mercury of 4.7 mg/kg.

Organochlorine pesticides were not detected in the 29 samples at concentrations greater than the laboratory reporting limits.

One PCB, Aroclor 1260, was detected in five of the 29 samples at concentrations that exceeded the HDOH Tier 1 EAL. The concentrations of the PCB Aroclor 1260 ranged from 1.19 to 3.27 mg/kg, which exceeded the HDOH Tier 1 EAL for PCBs of 1.1 mg/kg.

Dioxin was detected in six of the 29 soil samples with Toxicity Equivalent (TEQ) concentrations above the HDOH Tier 1 EAL of 240 nanograms per kilogram (ng/kg) ranging from 290 ng/kg to 710 ng/kg. The concentrations in these six soil samples, RHS-DU-1, RHS-DU-2, RHS-DU-9, RHS-DU-13, RHS-DU-18, and RHS-DU-19, indicated soils within those DUs are categorized as HDOH Soil Management Category C. According to HDOH guidance (2012), Soil Management Category C soils should be restricted to commercial/industrial land use with a formal restriction to the deed against sensitive land uses (e.g., residential, schools, etc.) in the absence of significant institutional and engineered control and HDOH approval. In addition, an Environmental Hazard Management Plan (EHMP) is required if soil is left on site for long-term management.



It should be noted that a replicate sample collected from one DU (RHS-DU-1) contained a Dioxin TEQ concentration of 2,900 ng/kg indicating soil from that DU may fall into the HDOH Soil Management Category D. HDOH guidance (2012) indicates that Soil Management Category D soils require remedial action under any land use scenario in order to reduce potential exposure and adverse health risks.

A preliminary Environmental Hazard Evaluation (EHE) was prepared for the field area of the site because of the potential for exposures to student athletes competing on the grassy/original field area. Based on the analytical laboratory results of the multi-increment surface soil samples collected from the four field area DUs, one contaminant, mercury, was detected in one of the four DUs, RHS-DU-29, at a concentration exceeding the HDOH Tier 1 EALs. The surface soil sample collected from RHS-DU-29 contained a mercury concentration of 5.1 mg/kg, which exceeded the HDOH Tier 1 EAL for mercury of 4.7 mg/kg. The surface soil samples from the remaining three field area DUs did not contain COPC concentrations exceeding the HDOH Tier 1 EALs.

The HDOH Tier 1 EAL for mercury is the direct contact EAL for residential sites, which is conservatively based on a noncarcinogen Hazard Quotient (HQ) of 0.2 to be protective of cumulative health risks potentially posed by sites with multiple noncarcinogens (up to a total of five). However, soil in RHS-DU-29 contained only one noncarcinogen at a concentration exceeding its HDOH Tier 1 EAL. As such, the direct contact EAL calculated using a HQ of 1.0 of 23 mg/kg may be used (see Table I-1 of the Technical Guidance Manual [TGM], HDOH, 2012). The mercury concentration of 5.1 mg/kg measured in RHS-DU-29 does not exceed the direct contact HDOH EAL for residential sites for mercury calculated using a HQ of 1.0 of 23 mg/kg. In other words, mercury surface soil concentrations measured in RHS-DU-29 are not expected to pose a risk for noncancer health effects assuming the strictest exposure parameters (i.e., residential).

Based on the laboratory analytical results, Bureau Veritas recommends the following:

- The stockpile DUs (RHS-DU-14 through RHD-DU-25) should be profiled and properly disposed.
- Conduct an additional site characterization investigation to further delineate the horizontal extent of contamination and the vertical extent of contamination.
- After the additional site characterization is conducted, an EHE should be prepared to assess the potential environmental hazards posed by the soil contamination.
- Develop remedial alternatives based on the results of the EHE.



## **1.0 INTRODUCTION AND PURPOSE**

The State of Hawaii Department of Education (DOE) retained Bureau Veritas North America, Inc. (Bureau Veritas) to collect and analyze soil samples and suspect asbestos materials from the track and field area of Radford High School. This report summarizes the data collected at the property described as being located at 4361 Salt Lake Boulevard, Tax Map Key (TMK) Number (1) 9-9-002: Parcel 023 (hereinafter referred to as the "site"). A site location map is presented as Figure 1, located behind the Figures tab.

On December 19, 2013 buried debris was encountered during excavation activities of the track and field area of the site. A topographic map from 1953 depicts the site and adjoining areas as a naval reservation located within Makalapa Crater. The topographic maps from 1959 through 2013 showed the site as a built-up area; labeled "Radford High School". The encountered buried debris is assumed to be associated with the site's former use as a military reservation, which may have included a former military dumpsite.

The purpose of this investigation was to assess the presence and magnitude of the soil contamination within the track construction area and associated soil stockpiles.

## **2.0 BACKGROUND**

### **2.1 SITE DESCRIPTION**

The site is located on the northwestern corner of a portion of TMK Number: (1) 9-9-002: Parcel 023 in a residential area of Honolulu, Hawaii. The site consists of approximately 25 acres of land located approximately one mile from Pearl Harbor. The site was developed as Radford High School by the State of Hawaii in 1957. The initial sampling area is presented as Figure 2.

#### **2.1.1 Soil Conditions**

The U.S. Department of Agriculture Soil Conservation Service (Foote, et al., 1972) classifies the soil within the area of the site as Makalapa clay, 2 to 6 percent slopes. This series consists of well-drained soils on uplands on the island of Oahu, near Salt Lake Crater, Diamond Head, and the Mokapu Peninsula. These soils formed in volcanic tuff. They are gently sloping to moderately steep. Elevations range from 20 to 200 feet. The annual rainfall amounts to 20 to 35 inches.

#### **2.1.2 Groundwater Conditions**

Bureau Veritas reviewed the Aquifer Identification and Classification for Oahu: Groundwater Protection Strategy for Hawaii (Mink, J.F. and L.S. Lau, 1990), published by the Water Resources Research Center at the University of Hawaii, for information on groundwater conditions below the subject property. The report describes the upper and lower aquifers below the site as part of the Waimalu aquifer system of the Pearl Harbor sector, on the island of Oahu.

The upper aquifer is an unconfined basal aquifer of the sedimentary type, occurring in non-volcanic lithology. Its status is described as an irreplaceable, low (250-1,000 milligrams per liter [mg/L] Chloride) water supply. It is considered a currently used ecologically important water source, and has a high vulnerability to contamination.



The lower aquifer is a confined basal aquifer of the dike type, occurring in dike compartments. Its status is described as an irreplaceable, low (250-1,000 mg/L Chloride) water supply, considered a currently used ecologically important water source, and has a moderate vulnerability to contamination.

The site is located above the State of Hawaii Department of Health (HDOH) defined Underground Injection Control (UIC) line. Areas above the UIC line denote potential underground drinking water aquifers. Areas below the UIC line generally denote aquifers that are unsuitable for drinking water purposes. Consequently, the aquifers above the site are considered drinking water aquifers.

The subsurface conditions under the site are interpreted from available data and may vary. Estimated groundwater flow direction is based on topography and nearby water features unless otherwise noted. Topography is not always a reliable basis for predicting groundwater flow direction. The local groundwater gradient under the site may be influenced naturally by tidal influences, zones of higher or lower permeability, or artificially by nearby pumping or recharge, and may deviate from the regional trend.

## **2.2 HISTORICAL SITE USE**

On December 19, 2013 buried debris was encountered during excavation activities of the track and field area of the site. A topographic map from 1953 depicts the site and adjoining areas as a naval reservation located within Makalapa Crater. The topographic maps from 1959 through 2013 showed the site as a built-up area labeled "Radford High School". The encountered buried debris is assumed to be associated with the site's former use as a military reservation, which may have included a former military dumpsite.

## **2.3 RECENT CONSTRUCTION ACTIVITIES**

Buried debris was encountered during recent construction activities involving refurbishing of the track and field area. The track and the end zone areas have been excavated, as well as the long-jump area and drain line within the field. In addition, the soil stockpiles are present onsite associated with the construction activities.

## **3.0 APPLICABLE ACTION LEVELS**

The applicable action levels for this site were established using the HDOH guidance document entitled "Evaluation of Environmental Hazards at Sites with Contaminated Soil and Groundwater" (HDOH, 2012), which is published by the HDOH Hazard Evaluation and Emergency Response (HEER) Office.

The site is located above the HDOH defined UIC line. Consequently, the aquifers below the site are considered drinking water aquifers. Furthermore, the site lies greater than 150 meters from a surface water body. Therefore, the analytical results were compared to the HDOH Tier 1 Environmental Action Levels (EALs) where groundwater is a current or potential source of drinking water and the nearest surface water body is greater than 150 meters from the site (HDOH, 2012; Table A).

## **4.0 FIELD ACTIVITIES**

### **4.1 INITIAL RESPONSE**

On December 20, 2013, Bureau Veritas responded onsite to collect one soil sample and six suspect asbestos containing material (ACM) samples at the site because buried debris had been found by MEI Corporation while excavating the northeastern portion of Radford High Schools track and field.



## **Asbestos**

Two types of suspect ACM, an off-white fibrous material with debris and an off-white/brown fibrous material with debris, were observed. Three samples of each type of suspect ACM were collected and sent to NVL Laboratories, Inc., a National Voluntary Laboratory Accreditation Program (NVLAP) laboratory located in Seattle, Washington. The samples were analyzed for asbestos content utilizing Polarized Light Microscopy (PLM) analysis. The three samples of the first type of suspect ACM (off-white fibrous material with debris) tested positive for asbestos with 51% to 55% Chrysotile. The three samples of the second type of suspect ACM (off-white/brown fibrous material with debris) did not test positive for asbestos. In summary, based on the laboratory results, one buried material (off-white fibrous material with debris) was confirmed as ACM (containing greater than [ $>$ ] 1% asbestos).

The two types of suspect ACM are presented in the table below. Laboratory results for the suspect ACM samples collected during Bureau Veritas' site visit are presented in Appendix B.

**Summary of Suspect ACM Sample Collection Information**

| <b>Material Description</b>                              | <b>Material Location</b>                     | <b>Concentration</b>  |
|--|--|-----------------------|
| Off-white fibrous material with debris (3 Samples)       | Northeastern corner of track and field area. | 51% to 55% Chrysotile |
| Off-white/brown fibrous material with debris (3 Samples) | Northeastern corner of track and field area. | None Detected         |

After the samples were collected, MEI personnel covered the buried debris with geofabric. The asbestos was subsequently bagged and stored in a locked storage area. On February 11, 2014, Unitek arrived onsite and double bagged five bags of ACM, which were subsequently properly disposed. The waste disposal manifest is presented in Appendix D.

## **Soil**

Also on December 20, 2013, Bureau Veritas collected one multi-increment soil sample, comprised of 30 sub-samples during potholing activities along the northeastern corner of the track and field area. The multi-increment soil sample was a representative sample collected from one decision unit (DU), RHS-01, that was established on the northeastern portion of the site, where buried debris was uncovered. The boundary of RHS-01 is shown on Figure 2, located behind the *Figures* tab. The sample was collected to assess the soil in the surrounding track area. The soil sample was analyzed for the following Chemicals of Potential Concern (COPC): Total Petroleum Hydrocarbons as Diesel Range Organics (TPH-DRO), TPH as Residual Range Organics (TPH-RRO), Semi-Volatile Organic Compounds (SVOCs), Resource Conservation and Recovery Act (RCRA) 8 Metals, polychlorinated biphenyls (PCBs), Organochlorine Pesticides, and dioxins.

Analysis of the soil sample reported that the concentrations of arsenic, cadmium, lead and dioxin exceeded their respective HDOH Tier 1 EALs. Further sampling and analysis were recommended to assess the extent of the soil contamination within the track construction area.





## 4.2 INVESTIGATION PLANNING

Following HDOH guidance (HDOH, 2013), a multi-increment soil sampling approach was selected for this site investigation. The HDOH defines a DU as a portion of a site where a decision is to be made regarding the extent and magnitude of contaminants identified within the unit, as well as the potential environmental hazards posed by the contaminants. A scope of work was prepared, submitted, and approved by HDOH prior to conducting a follow-up investigation.

To facilitate the additional investigation, the site was divided into the following three main areas of concern: (1) the excavated track area, (2) the soil stockpiles, and (3) the grassy/original field area. Each of the areas of concern was further divided into DUs based on visual contamination and size. A total of 29 DUs (RHS-DU-1 through RHS-DU-29) were established across the site to facilitate the collection of multi-increment soil samples from the surface soil layers in each DU. The boundaries of each DU are shown on Figure 3, located behind the *Figures* tab.

## 4.3 SOIL SAMPLING ACTIVITIES

On February 4 and 5, 2014, Bureau Veritas collected multi-increment surface soil samples from the 29 DUs established at the site. Soil samples were collected using a soil sampling probe that was advanced approximately six inches into the soil. Each increment consisted of approximately one to two ounces of soil. A total of 50 increments were collected using a systematic random approach across each DU. The increments were placed in a new, plastic Ziploc™ plastic bag. The soil coring device was decontaminated using Alconox™ cleaner and double rinsed with distilled water prior to collecting the sample.

The sample container was labeled and placed into a cooler with wet-ice to begin the preservation process during delivery to the laboratory. The sample was logged on a chain-of-custody form, which accompanied the sample to the laboratory. Three sets of replicate soil samples were also collected, one set from each area of concern, and analyzed (see Section 4.4).

A summary of the primary soil samples is presented in the following table:

**Summary of Primary Soil Samples**

| Sample ID | DU   | Date/Time Sample Collected | Sample Interval (bgs) | Description of Sample Location |
|-----------|------|----------------------------|-----------------------|--------------------------------|
| RHS-DU-1  | DU1  | 2-4-14/1525                | 0 to 0.5 feet         | Eastern portion of the track   |
| RHS-DU-2  | DU2  | 2-4-14/1500                | 0 to 0.5 feet         | Eastern portion of the track   |
| RHS-DU-3  | DU3  | 2-4-14/1403                | 0 to 0.5 feet         | Eastern portion of the track   |
| RHS-DU-4  | DU4  | 2-4-14/1145                | 0 to 0.5 feet         | Southern portion of the track  |
| RHS-DU-5  | DU5  | 2-4-14/1220                | 0 to 0.5 feet         | Southern portion of the track  |
| RHS-DU-6  | DU6  | 2-4-14/1402                | 0 to 0.5 feet         | Southern portion of the track  |
| RHS-DU-7  | DU7  | 2-4-14/1131                | 0 to 0.5 feet         | Southern portion of the track  |
| RHS-DU-8  | DU8  | 2-4-14/1424                | 0 to 0.5 feet         | Western portion of the track   |
| RHS-DU-9  | DU9  | 2-4-14/1455                | 0 to 0.5 feet         | Western portion of the track   |
| RHS-DU-10 | DU10 | 2-5-14/0915                | 0 to 0.5 feet         | Northern portion of the track  |
| RHS-DU-11 | DU11 | 2-5-14/0900                | 0 to 0.5 feet         | Northern portion of the track  |



| Sample ID | DU   | Date/Time Sample Collected | Sample Interval (bgs) | Description of Sample Location   |
|-----------|------|----------------------------|-----------------------|----------------------------------|
| RHS-DU-12 | DU12 | 2-5-14/0945                | 0 to 0.5 feet         | Northern portion of the track    |
| RHS-DU-13 | DU13 | 2-5-14/0945                | 0 to 0.5 feet         | Northern portion of the track    |
| RHS-DU-14 | DU14 | 2-4-14/1252                | 0 to 0.5 feet         | Northern soil stockpile          |
| RHS-DU-15 | DU15 | 2-4-14/1345                | 0 to 0.5 feet         | Northern soil stockpile          |
| RHS-DU-16 | DU16 | 2-4-14/1411                | 0 to 0.5 feet         | Northern soil stockpile          |
| RHS-DU-17 | DU17 | 2-4-14/1436                | 0 to 0.5 feet         | North of end zone soil stockpile |
| RHS-DU-18 | DU18 | 2-4-14/1300                | 0 to 0.5 feet         | North of end zone soil stockpile |
| RHS-DU-19 | DU19 | 2-4-14/1422                | 0 to 0.5 feet         | North of end zone soil stockpile |
| RHS-DU-20 | DU20 | 2-4-14/1146                | 0 to 0.5 feet         | End zone soil stockpile          |
| RHS-DU-21 | DU21 | 2-4-14/1226                | 0 to 0.5 feet         | End zone soil stockpile          |
| RHS-DU-22 | DU22 | 2-4-14/1015                | 0 to 0.5 feet         | Western soil stockpile           |
| RHS-DU-23 | DU23 | 2-4-14/1023                | 0 to 0.5 feet         | Western soil stockpile           |
| RHS-DU-24 | DU24 | 2-4-14/1050                | 0 to 0.5 feet         | Southern soil stockpile          |
| RHS-DU-25 | DU25 | 2-4-14/1115                | 0 to 0.5 feet         | Southern soil stockpile          |
| RHS-DU-26 | DU26 | 2-4-14/1020                | 0 to 0.5 feet         | Northern portion of the field    |
| RHS-DU-27 | DU27 | 2-4-14/1000                | 0 to 0.5 feet         | Eastern portion of the field     |
| RHS-DU-28 | DU28 | 2-4-14/1007                | 0 to 0.5 feet         | Western portion of the field     |
| RHS-DU-29 | DU29 | 2-4-14/0945                | 0 to 0.5 feet         | Southern portion of the field    |

#### 4.4 FIELD REPLICATE SAMPLING

Field Quality Assurance/Quality Control (QA/QC) was performed through the collection of three sets of replicate samples, each consisting of the primary sample and an associated duplicate sample and triplicate sample. The replicate samples were collected from the upper six inches of DU1, DU17 and DU26. A summary of the primary/replicate soil samples is presented in the following table:

**Summary of Primary/Replicate Soil Samples**

| Sample ID   | Sample Type | DU   | Date/Time Sample Collected | Sample Interval (bgs) | Description of Sample Location   |
|-------------|-------------|------|----------------------------|-----------------------|----------------------------------|
| RHS-DU-1    | primary     | DU1  | 2-4-14/1525                | 0 to 0.5 feet         | Eastern portion of the track     |
| RHS-DU-1.2  | duplicate   | DU1  | 2-5-14/0933                | 0 to 0.5 feet         | Eastern portion of the track     |
| RHS-DU-1.3  | triplicate  | DU1  | 2-5-14/1030                | 0 to 0.5 feet         | Eastern portion of the track     |
| RHS-DU-17   | primary     | DU17 | 2-4-14/1436                | 0 to 0.5 feet         | North of end zone soil stockpile |
| RHS-DU-17.2 | duplicate   | DU17 | 2-4-14/1449                | 0 to 0.5 feet         | North of end zone soil stockpile |
| RHS-DU-17.3 | triplicate  | DU17 | 2-4-14/1502                | 0 to 0.5 feet         | North of end zone soil stockpile |
| RHS-DU-26   | primary     | DU26 | 2-4-14/1020                | 0 to 0.5 feet         | Northern portion of the field    |
| RHS-DU-26.2 | duplicate   | DU26 | 2-4-14/1059                | 0 to 0.5 feet         | Northern portion of the field    |
| RHS-DU-26.3 | triplicate  | DU26 | 2-4-14/1145                | 0 to 0.5 feet         | Northern portion of the field    |





Soil analytical results for the primary and replicate samples are summarized in Table 4, located behind the *Tables* tab.

The data from the replicate samples were used to calculate important statistical measures, including Relative Percent Difference (RPD) and Relative Standard Deviation (RSD). A detailed description of the calculations performed using data from the replicate samples is presented in Section 5.4.

#### **4.5 CHAIN-OF-CUSTODY**

Each soil sample was logged on a chain-of-custody form upon collection. The chain-of-custody forms accompanied the samples from the field to the laboratory. Whenever the samples changed hands, the custodian of the samples signed the “relinquished by” section and the receiving person signed the “received by” section. The soil samples were delivered by hand under standard chain-of-custody procedures to TestAmerica and Environmental Services Network (ESN) Pacific Inc.

#### **4.6 DECONTAMINATION PROCEDURES**

New disposable gloves were worn by the sampler during the soil sample collection activities. The gloves were disposed of following the collection of a soil sample and new gloves were donned prior to the collection of the following soil sample. Equipment, such as the soil sampling probes and trowels, were decontaminated between sample collections as follows:

- Washed with Alconox™
- Double rinsed with distilled water
- Air dried

#### **4.7 INVESTIGATION DERIVED WASTE**

The field activities generated different types of investigation derived waste (IDW), including used personal protective equipment (PPE) and disposable equipment (i.e. nitrile gloves, paper towels, etc.).

Used PPE and disposable equipment was sealed in a plastic bag and placed in a municipal refuse dumpster. The used PPE and equipment were not considered hazardous and were sent to a municipal landfill.

### **5.0 LABORATORY ANALYTICAL RESULTS**

The soil samples collected as part of the initial response/investigation were submitted to TestAmerica located in Aiea, Hawaii. The samples collected during the follow-up investigation were submitted to ESN Analytical Laboratories located in Honolulu, Hawaii for analytical testing. The soil samples were submitted to the lab in a cooler containing wet ice. Copies of the Analytical Reports and Chain-of-Custody documents are presented in Appendices A through C. Tables 1 through 5, located behind the *Tables* tab, present the laboratory analytical results for the samples.

The following analyses were performed on the soil samples collected from the site:

- TPH-DRO using Environmental Protection Agency (EPA) Method 8015B.
- TPH-RRO using EPA Method 8015B
- SVOCs using EPA Method 8270C
- RCRA 8 Metals using EPA Methods 6010B and 7471



- PCBs using EPA Method 8082
- Organochlorine Pesticides using EPA Method 8081A
- Dioxin using EPA Method 8290

## 5.1 INITIAL INVESTIGATION SOIL ANALYTICAL RESULTS

During the initial investigation, one multi-incremental soil sample was collected from one DU (RHS-01). Table 1 summarizes the analytical results of the soil sample and compares the results to the HDOH Tier 1 EALs. The soil sample was obtained from 0 to 0.5 feet below ground surface (bgs).

### 5.1.1 TPH

TPH-DRO was detected at a concentration of 24 milligrams per kilogram (mg/kg), which was below the HDOH Tier 1 EAL for TPH-DRO of 100 mg/kg. TPH-RRO was detected at a concentration of 23 mg/kg, which was below the HDOH Tier 1 EAL for TPH-RRO of 500 mg/kg.

### 5.1.2 SVOCs

Four SVOC analytes, 2-methylnaphthalene, benzo[b]fluoranthene, indeno[1,2,3-cd]pyrene, and naphthalene, were detected, but at concentrations less than their respective HDOH Tier 1 EALs. The remaining SVOC analytes were not detected at concentrations greater than the laboratory reporting limits.

### 5.1.3 RCRA 8 Metals

Three of the eight RCRA metals, arsenic, cadmium, and lead, were detected at concentrations greater than their respective HDOH Tier 1 EALs. Arsenic was detected at a concentration of 43 mg/kg, which was above the HDOH Tier 1 EAL for arsenic of 24 mg/kg. Cadmium was detected at a concentration of 26 mg/kg, which was above the HDOH Tier 1 EAL for cadmium of 14 mg/kg. Lead was detected at a concentration of 5,300 mg/kg, which was above the HDOH Tier 1 EAL for lead of 200 mg/kg.

Four of the eight RCRA metals, barium, chromium, silver, and mercury, were detected, but at concentrations below their respective HDOH Tier 1 EALs. Barium was detected at a concentration of 940 mg/kg, which was below the HDOH Tier 1 EAL for barium of 1,000 mg/kg. Chromium was detected at a concentration of 210 mg/kg, which was below the HDOH Tier 1 EAL for chromium of 1,100 mg/kg. Silver was detected at a concentration of 14 mg/kg, which was below the HDOH Tier 1 EAL for silver of 78 mg/kg. Mercury was detected at a concentration of 1.1 mg/kg, which was below the HDOH Tier 1 EAL for mercury of 4.7 mg/kg.

Selenium was not detected at a concentration greater than the laboratory reporting limit.

### 5.1.4 PCBs

The PCB Aroclor 1260 was detected at a concentration of 0.094 mg/kg, which was below the HDOH Tier 1 EAL for PCBs of 1.1 mg/kg.

### 5.1.5 Organochlorine Pesticides

Five organochlorine pesticide analytes, DDE, DDT, dieldrin, endrin, and endrin aldehyde, were detected, but at concentrations less than their respective HDOH Tier 1 EALs. The remaining organochlorine pesticide analytes were not detected at concentrations greater than the laboratory reporting limits.



### 5.1.6 Dioxin

The term dioxin is used to refer to a family of chlorinated compounds with similar chemical structures and mechanisms of toxicity, referred to as “congeners.” Because the various dioxins/furans congeners are not equally toxic, concentrations are reported relative to 2, 3, 7, 8-tetrachlorodibenzo-p-dioxin (2, 3, 7, 8-TCDD) referred to as the Dioxin Toxicity Equivalent (TEQ) concentration.

For the calculations of the Dioxin TEQ concentrations, the 2005 World Health Organization (WHO) toxicity equivalence factors (TEFs) were used (Van den Berg, et. al, 2006). The Dioxin TEQ calculations are presented in Table 2, located behind the *Tables* tab.

The HDOH published the document titled “Evaluation of Environmental Hazards at Sites with Contaminated Soil and Groundwater,” dated Fall 2011 (HDOH, 2012). The technical memorandum indicates that the Tier 1EAL for Dioxin TEQ is 240 ng/kg.

The HDOH document also provides a Summary of TEQ Dioxin Soil Action Levels and associated management categories. An abbreviated version of the Summary is shown in the Table below.

**Summary of TEQ Dioxin Soil Action Levels and Associated Management Categories**

| Soil Management Category | TEQ                               | Recommended Action  |
|--------------------------|-----------------------------------|---|
| A                        | Less than 20 ng/kg                | <b>Background.</b> Within range of expected background conditions in non-agricultural and non-industrial areas. No further action required and no restrictions on land use.   |
| B                        | Between 20 ng/kg and 240 ng/kg    | <b>Minimally Impacted.</b> Exceeds expected background conditions but within range anticipated for agricultural fields. Potential health risks considered insignificant. Include Category B soil in remedial actions for more heavily contaminated spill areas as practicable in order to reduce exposure (e.g., outer margins of pesticide mixing areas).  |
| C                        | Between 240 ng/kg and 1,500 ng/kg | <b>Moderately impacted.</b> Typical of incinerator ash, burn pits, wood treatment operations that used pentachlorophenol, and the margins of heavily impacted, pesticide mixing areas associated with former sugarcane operations that used pentachlorophenol. Restriction to commercial/industrial land use required with a formal restriction to the deed against sensitive land uses (e.g., residential, schools, etc.) in the absence of significant institutional and engineered control and HDOH approval. Environmental Hazard Management Plan (EHMP) required if soil left on site for long-term management. This includes controls to ensure no off-site dispersion (e.g., dust or surface runoff) or inadvertent excavation and reuse at properties with sensitive land uses. |
| D                        | Greater than 1,500 ng/kg          | <b>Heavily Impacted.</b> Typical of former pesticide mixing areas that used pentachlorophenol. Remedial actions required under any land use scenario in order to reduce potential exposure. Potentially adverse health risks under both sensitive and commercial/industrial land use scenarios in the absence of significant institutional and/or engineered controls.  |

Sample RHS-01 had a TEQ dioxin concentration of 2,019 ng/kg, which exceeded the HDOH Tier 1 EAL for both unrestricted use (240 ng/kg) and commercial/industrial use (1,500 ng/kg). Following HDOH guidance (2012), the calculated TEQ Dioxin soil concentration is within the HDOH TEQ Dioxin Soil Management Category D (Heavily Impacted). The Guidance indicates that remedial actions are required for Category D soils under any land use scenario in order to reduce potential exposure.



## 5.2 FOLLOW-UP INVESTIGATION SOIL ANALYTICAL RESULTS

A total of 29 primary, multi-increment, surface soil samples were collected (i.e., one soil sample from each DU) during the follow-up investigation. Replicate multi-increment, surface soil samples were also collected from three DUs (i.e., one DU representing each of the three areas of concern: the track, the field, and the soil stockpile areas). Samples were obtained from the upper 6-inches of soil. Table 3, summarizes the analytical results of the soil samples and compares the results to the HDOH Tier 1 EALs.

### 5.2.1 TPH

TPH-DRO and TPH-RRO were not detected in the 29 soil samples at concentrations greater than the laboratory reporting limits.

### 5.2.2 SVOCs

One SVOC analyte, benzo(a)pyrene, was detected at concentrations that equaled or exceeded the HDOH Tier 1 EAL. In three of the 29 samples, benzo(a)pyrene concentrations that ranged between 0.15 and 0.43 mg/kg were equal to or exceeded the HDOH Tier 1 EAL for benzo(a)pyrene of 0.15 mg/kg.

Seven SVOC analytes, benzo(a)anthracene, benzo(b)fluoranthene, benzo(k)fluoranthene, chrysene, fluoranthene, phenanthrene, and pyrene, were detected in one or more of the 29 samples, but at concentrations below their respective HDOH Tier 1 EALs.

The remaining SVOCs analytes were not detected at concentrations greater than the laboratory reporting limits.

### 5.2.3 RCRA 8 Metals

Five of the eight RCRA metals, arsenic, barium, cadmium, lead, and mercury, were detected at concentrations that exceeded their respective HDOH Tier 1 EALs. Arsenic was detected in seven of the 29 samples at concentrations ranging from 24 to 35 mg/kg, which were equal to or exceeded the HDOH Tier 1 EAL for arsenic of 24 mg/kg. Barium was detected in one of the samples at a concentration of 1,200 mg/kg, which exceeded the HDOH Tier 1 EAL for barium of 1,000 mg/kg. Cadmium was detected in one of the samples at a concentration of 14 mg/kg, which was equal to the HDOH Tier 1 EAL for cadmium of 14 mg/kg. Lead was detected in 15 of the samples at concentrations ranging from 240 to 6,200 mg/kg, which exceeded the HDOH Tier 1 EAL for lead of 200 mg/kg. Mercury was detected in 22 of the samples at concentrations ranging from 4.8 to 200 mg/kg, which exceeded the HDOH Tier 1 EAL for mercury of 4.7 mg/kg.

Chromium was detected in all of the 29 samples, but at concentrations below the HDOH Tier 1 EAL for chromium of 1,100 mg/kg. Selenium and silver were not detected in the 29 samples at concentrations greater than the laboratory reporting limits.

### 5.2.4 PCBs

One PCB, Aroclor 1260, was detected in five of the 29 samples at concentrations that exceeded the HDOH Tier 1 EAL. The concentrations of the PCB Aroclor 1260 ranged from 1.19 to 3.27 mg/kg, which exceeded the HDOH Tier 1 EAL for PCBs of 1.1 mg/kg. The remaining PCB Aroclors were not detected at concentrations greater than the laboratory reporting limits.



### 5.2.5 Organochlorine Pesticides

Seven of the organochlorine pesticides analytes, alpha-chlordane, gamma-chlordane, chlordane (technical), dieldrin, endosulfan II, DDE, and DDT, were detected in one or more of the soil samples, but at concentrations less than their respective HDOH Tier 1 EALs. The remaining organochlorine pesticides were not detected at concentrations greater than the laboratory reporting limits.

### 5.2.6 Dioxin

The Dioxin TEQ calculations were performed by the analytical laboratory and are presented within the laboratory analytical reports (see Appendix C). The analytical laboratory calculated the Dioxin TEQs following three different approaches (i.e., a lower bound, a medium bound, and an upper bound). The medium bound approach used one-half the laboratory detection limit as a proxy concentration for the congeners that were not detected. The medium bound approach followed the approach recommended by the HDOH (HDOH, 2012). As such, the medium bound Dioxin TEQ concentrations were used for comparison to the HDOH action levels (see Table 3).

Dioxin was detected in 6 of the 29 soil samples with TEQ concentrations above the HDOH TEQ Dioxin SAL of 240 ng/kg ranging from 290 ng/kg to 710 ng/kg. Dioxin TEQ concentrations in the remaining 23 samples did not exceed the HDOH Tier EAL of 240 ng/kg. The concentrations in these six soil samples, RHS-DU-1, RHS-DU-2, RHS-DU-9, RHS-DU-13, RHS-DU-18, and RHS-DU-19, indicated soils within those DUs are categorized as HDOH Soil Management Category C. According to HDOH guidance (2012), Soil Management Category C soils should be restricted to commercial/industrial land use with a formal restriction to the deed against sensitive land uses (e.g., residential, schools, etc.) in the absence of significant institutional and engineered control and HDOH approval. In addition, an Environmental Hazard Management Plan (EHMP) is required if soil is left on site for long-term management.

It should be noted that a replicate sample collected from one DU (RHS-DU-1) contained a Dioxin TEQ concentration of 2,900 ng/kg indicating soil from that DU may fall into the HDOH Soil Management Category D. HDOH guidance (2012) indicates that Soil Management Category D soils require remedial action under any land use scenario in order to reduce potential exposure and adverse health risks.

Six of the samples, RHS-DU-20, RHS-DU-21, RHS-DU-23, RHS-DU-24, RHS-DU-25, and RHS-DU-28, had Dioxin TEQ concentrations less than 20 ng/kg, indicating those six DUs had Dioxin TEQ concentrations that are within the range of expected background conditions in non-agricultural and non-industrial areas (HDOH Soil Management Category A). No further actions are required and no restrictions on land use are necessary for HDOH Soil Management Category A soils.

The remaining 17 samples had Dioxin TEQ concentrations greater than 20 ng/kg, but less than 240 ng/kg indicating those 17 DUs had Dioxin TEQ concentrations that exceeded expected background conditions but are within the range anticipated for agricultural fields (HDOH Soil Management Category B): RHS-DU-3, RHS-DU-4, RHS-DU-5, RHS-DU-6, RHS-DU-7, RHS-DU-8, RHS-DU-10, RHS-DU-11, RHS-DU-12, RHS-DU-14, RHS-DU-15, RHS-DU-16, RHS-DU-17, RHS-DU-22, RHS-DU-26, RHS-DU-27, and RHS-DU-29. Potential health risks are considered insignificant for HDOH Soil Management Category B soils.



### 5.3 LABORATORY QUALITY CONTROL

The soil sample collected during the initial investigation was submitted to TestAmerica Analytical Laboratory in Aiea, Oahu, Hawaii for laboratory analysis. The soil samples collected as part of the follow-up investigation were submitted to ESN Analytical Laboratory in Honolulu, Oahu, Hawaii for laboratory analysis. Analytical data were generated following EPA methods (SW-846 protocols), laboratory standard operating procedures (SOP), and QA/QC guidelines for sample analysis. Common laboratory QC checks include the use of Method Blank, Matrix Spike and Matrix Spike Duplicate, and Laboratory Control and Laboratory Control Duplicate samples.

Adequate reporting levels of the chemicals of concern are dependent on the sample matrix, naturally occurring background concentrations, and laboratory instrumentation. For some of the constituents, the laboratory reporting limits were greater than the HDOH Tier 1 EALs. As stated in the HDOH guidance document (HDOH, 2012), laboratory reporting limits of chemicals were not directly considered in development of the HDOH EALs, and therefore, some standards are not able to be achieved through standard laboratory analysis. In these cases, it is generally acceptable to consider the reporting limit in place of the screening level.

### 5.4 FIELD QUALITY CONTROL

Field QA/QC was performed through the collection of three sets of replicate samples consisting of the primary sample and an associated duplicate and triplicate sample. The table below lists sample identification numbers for the primary samples and the associated duplicate and triplicate samples.

**Duplicate and Triplicate Sample Identification**

| Primary Sample Identification Number | Associated Duplicate (Dup) and Triplicate (Trip) Sample Identification Number |
|--------------------------------------|---|
| RHS-DU-1                             | RHS-DU-1.2 and RHS-DU-1.3   |
| RHS-DU-17                            | RHS-DU-17.2 and RHS-DU-17.3   |
| RHS-DU-26                            | RHS-DU-26.2 and RHS-DU-26.3   |

The statistical summaries were calculated for all chemical constituents with detected concentrations in the primary, duplicate, and triplicate samples and are presented in Table 5. For some analytes, the statistical calculations could not be performed because the analyte was not detected above laboratory reporting limits in one or more of the samples (i.e., primary, duplicate, or triplicate sample). In these cases, the RPD and RSD calculations were not performed.

#### Relative Percent Difference

The RPD, expressed as a percent, is a measure of precision between two sample values (the primary sample and the duplicate sample as well as the primary sample and the triplicate sample).

The RPD is calculated as the positive difference between two measurements (primary and duplicate; primary and triplicate) divided by the average of the two measured values and multiplied by 100. Typically, if the RPD is less than or equal to 35%, then the quality of the data is acceptable. The goal for this investigation was 35% or less.





## Standard Deviation and Relative Standard Deviation

The standard deviation is a statistical measure of the scatter, or variability, of several sample values around their average. The lower the standard deviation, the lower the variability of the sample values observed in the data.

The RSD, expressed as a percent, is a measure of precision between several sample values (the primary, duplicate, and triplicate samples). The RSD differs from the RPD in that it measures the precision between several sample values versus just two sample values. The RSD is calculated as the standard deviation divided by the mean (average). The RSD is useful for comparing the uncertainty between different measurements. Typically, if the RSD is less than or equal to 35%, then the quality of the data is acceptable. The goal for this investigation was 35% or less.

### RHS-DU-1

The RPDs and RSDs for the duplicate and triplicate samples of sample RHS-DU-1 exhibited moderate to significant variability. The calculated RPDs comparing the primary and the duplicate sample for dioxin, arsenic, barium, cadmium, chromium, and lead met the goal of less than 35%. The calculated RPD comparing the primary and duplicate sample for mercury was greater than 35%.

The calculated RPDs comparing the primary and the triplicate sample for arsenic, barium and chromium met the goal of less than 35%. The calculated RPDs comparing the primary and the triplicate sample for dioxin, cadmium, lead and mercury were greater than 35%.

The calculated RSDs for arsenic, barium, and chromium met the goal of less than 35%. The calculated RSDs for dioxin, cadmium, lead, and mercury were greater than 35%.

### RHS-DU-17

The RPDs and RSDs for the duplicate and triplicate samples of sample RHS-DU-17 exhibited moderate to significant variability. The calculated RPDs comparing the primary and the duplicate sample for dioxin, arsenic, barium, cadmium, chromium, mercury, chrysene, fluoranthene, pyrene, Aroclor 1260 (PCB), and DDT met the goal of less than 35%. The calculated RPD comparing the primary and duplicate sample for lead was greater than 35%.

The calculated RPDs comparing the primary and the triplicate sample for dioxin, arsenic, barium, cadmium, chromium, mercury, Aroclor 1260 (PCB), and DDT met the goal of less than 35%. The calculated RPDs comparing the primary and the triplicate sample for lead, chrysene, fluoranthene, and pyrene were greater than 35%.

The calculated RSDs for dioxin, arsenic, barium, cadmium, chromium, lead, mercury, Aroclor 1260 (PCB), and DDT met the goal of less than 35%. The calculated RSDs for chrysene, fluoranthene, and pyrene were greater than 35%.

### RHS-DU-26

The RPDs and RSDs for the duplicate and triplicate samples of sample RHS-DU-26 exhibited moderate to significant variability. The calculated RPDs comparing the primary and the duplicate sample for chromium and mercury met the goal of less than 35%. The calculated RPDs comparing the primary and duplicate sample for dioxin, barium, and lead were greater than 35%.

The calculated RPDs comparing the primary and the triplicate sample for barium, chromium, lead, and mercury met the goal of less than 35%. The calculated RPD comparing the primary and the triplicate sample for dioxin was greater than 35%.



The calculated RSDs for dioxin, arsenic, barium, chromium, lead, and mercury met the goal of less than 35%.

### Summary

The RPDs and RSDs that exceeded the goal of 35% are likely the result of heterogeneous distribution of the contaminants in the soil column and to some degree variability in the sample collection process. Alternatives to reducing the variability may be applicable, such as reducing the area of each DU investigated. Overall, the data quality appears to be sufficient for decision making purposes.

## **5.5 PRELIMINARY ENVIRONMENTAL HAZARD EVALUATION - FIELD AREA**

A preliminary Environmental Hazard Evaluation (EHE) was prepared for the field area of the site because of the potential for exposures to student athletes competing on the grassy/original field area. The field area was divided into four DUs (i.e., RHS-DU-26, RHS-DU-27, RHS-DU-28, and RHS-DU-29) as shown in Figure 3. Based on the analytical laboratory results of the multi-increment surface soil samples collected from the four field area DUs, one contaminant, mercury, was detected in one of the four DUs, RHS-DU-29, at a concentration exceeding the HDOH Tier 1 EALs. The surface soil sample collected from RHS-DU-29 contained a mercury concentration of 5.1 mg/kg, which exceeded the HDOH Tier 1 EAL for mercury of 4.7 mg/kg. The surface soil samples from the remaining three field area DUs did not contain COPC concentrations exceeding the HDOH Tier 1 EALs.

The HDOH Tier 1 EAL for mercury is the direct contact EAL for residential sites, which is conservatively based on a noncarcinogen Hazard Quotient (HQ) of 0.2 to be protective of cumulative health risks potentially posed by sites with multiple noncarcinogens (up to a total of five). However, soil in RHS-DU-29 contained only one noncarcinogen at a concentration exceeding its HDOH Tier 1 EAL. As such, the direct contact EAL calculated using a HQ of 1.0 of 23 mg/kg may be used (see Table I-1 of the Technical Guidance Manual [TGM], HDOH, 2012). The mercury concentration of 5.1 mg/kg measured in RHS-DU-29 does not exceed the direct contact HDOH EAL for residential sites for mercury calculated using a HQ of 1.0 of 23 mg/kg. In other words, mercury surface soil concentrations measured in RHS-DU-29 are not expected to pose a risk for noncancer health effects assuming the strictest exposure parameters (i.e., residential).

## **6.0 SUMMARY AND CONCLUSION**

Buried debris was discovered during excavation activities of the track and field areas, which is assumed to be associated with the site's former use as a military reservation and possibly a military dumpsite. The purpose of this investigation was to assess the presence and magnitude of the soil contamination within the track construction area and associated soil stockpiles.

On December 20, 2013, after the discovery of buried debris, Bureau Veritas responded onsite to collect one multi-increment soil sample and six suspect ACM samples. Two types of suspect ACM, an off-white fibrous material with debris and an off-white/brown fibrous material with debris, were observed. Three samples of each type of suspect ACM were collected and sent to NVL Laboratories, Inc., a NVLAP laboratory located in Seattle, Washington. The samples were analyzed for asbestos content utilizing PLM analysis. The three samples of the first type of suspect ACM (off-white fibrous material with debris) tested positive for asbestos with 51% to 55% Chrysotile. The three samples of the second type of suspect ACM (off-white/brown fibrous material with debris) did not test positive for asbestos. After the samples were collected, MEI personnel covered the buried debris with geofabric. The asbestos was subsequently





bagged and stored in a locked storage area. On February 11, 2014, Unitek arrived onsite and double bagged five bags of ACM, which were subsequently properly disposed.

The multi-increment soil sample was a representative sample collected from one decision unit (RHS-01), that was established on the northeastern portion of the site, where buried material was uncovered. The sample was collected to assess the soil in the track area surrounding the buried debris. The soil sample was analyzed for TPH-DRO, TPH-RRO, SVOCs, RCRA 8 Metals, PCBs, Organochlorine Pesticides, and dioxin. Analysis of the soil sample indicated that the soil concentrations of arsenic, cadmium, lead and dioxin exceeded their respective HDOH Tier 1 EALs. Further sampling and analysis was recommended to delineate the extent of the soil contamination.

A scope of work was prepared, submitted, and approved by HDOH prior to conducting a follow-up investigation. On February 4 and 5, 2014, a follow-up soil investigation was conducted by Bureau Veritas. To facilitate the additional investigation, the site was divided into the following three areas of concern: (1) the excavated track area, (2) the soil stockpiles, and (3) the grassy/original field area. Each of the areas of concern was further divided into DUs based on visual contamination and size. A total of 29 DUs (RHS-DU-1 through RHS-DU-29) were established across the site to facilitate the collection of multi-increment soil samples from the surface soil layer in each DU.

One multi-increment surface soil sample was collected from each DU for a total of 29 samples. Three sets of replicate samples were also collected from three DUs selected to represent each of the three areas of concern. The samples were submitted to ESN Analytical Laboratory located in Honolulu, Hawaii for analytical testing and were analyzed for the following:

- TPH-DRO using EPA Method 8015B.
- TPH-RRO using EPA Method 8015B
- SVOCs using EPA Method 8270C
- RCRA 8 Metals using EPA Methods 6010B and 7471
- PCBs using EPA Method 8082
- Organochlorine Pesticides using EPA Method 8081A
- Dioxin using EPA Method 8290

TPH-DRO and TPH-RRO were not detected in the 29 soil samples at concentrations greater than the laboratory reporting limits.

One SVOC analyte, benzo(a)pyrene, was detected at concentrations that equaled or exceeded the HDOH Tier 1 EAL. In three of the 29 samples, benzo(a)pyrene concentrations that ranged between 0.15 and 0.43 mg/kg were equal to or exceeded the HDOH Tier 1 EAL for benzo(a)pyrene of 0.15 mg/kg.

Five of the eight RCRA metals, arsenic, barium, cadmium, lead, and mercury, were detected at concentrations that exceeded their respective HDOH Tier 1 EALs. Arsenic was detected in seven of the 29 samples at concentrations ranging from 24 to 35 mg/kg, which were equal to or exceeded the HDOH Tier 1 EAL for arsenic of 24 mg/kg. Barium was detected in one of the samples at a concentration of 1,200 mg/kg, which exceeded the HDOH Tier 1 EAL for barium of 1,000 mg/kg. Cadmium was detected in one of the samples at a concentration of 14 mg/kg, which was equal to the HDOH Tier 1 EAL for cadmium of 14 mg/kg. Lead was detected in 15 of the samples at concentrations ranging from 240 to 6,200 mg/kg, which exceeded the HDOH Tier 1 EAL for lead of 200 mg/kg. Mercury was detected in 22 of the samples at concentrations ranging from 4.8 to 200 mg/kg, which exceeded the HDOH Tier 1 EAL for mercury of 4.7 mg/kg.



Organochlorine pesticides were not detected in the 29 samples at concentrations greater than the laboratory reporting limits.

One PCB, Aroclor 1260, was detected in five of the 29 samples at concentrations that exceeded the HDOH Tier 1 EAL. The concentrations of the PCB Aroclor 1260 ranged from 1.19 to 3.27 mg/kg, which exceeded the HDOH Tier 1 EAL for PCBs of 1.1 mg/kg.

Dioxin was detected in six of the 29 soil samples with TEQ concentrations above the HDOH Tier 1 EAL of 240 ng/kg ranging from 290 ng/kg to 710 ng/kg. The concentrations in these six soil samples, RHS-DU-1, RHS-DU-2, RHS-DU-9, RHS-DU-13, RHS-DU-18, and RHS-DU-19, indicated soils within those DUs are categorized as HDOH Soil Management Category C. According to HDOH guidance (2012), Soil Management Category C soils should be restricted to commercial/industrial land use with a formal restriction to the deed against sensitive land uses (e.g., residential, schools, etc.) in the absence of significant institutional and engineered control and HDOH approval. In addition, an Environmental Hazard Management Plan (EHMP) is required if soil is left on site for long-term management.

It should be noted that a replicate sample collected from one DU (RHS-DU-1) contained a Dioxin TEQ concentration of 2,900 ng/kg indicating soil from that DU may fall into the HDOH Soil Management Category D. HDOH guidance (2012) indicates that Soil Management Category D soils require remedial action under any land use scenario in order to reduce potential exposure and adverse health risks.

A preliminary Environmental Hazard Evaluation (EHE) was prepared for the field area of the site because of the potential for exposures to student athletes competing on the grassy/original field area. Based on the analytical laboratory results of the multi-increment surface soil samples collected from the four field area DUs, one contaminant, mercury, was detected in one of the four DUs, RHS-DU-29, at a concentration exceeding the HDOH Tier 1 EALs. The surface soil sample collected from RHS-DU-29 contained a mercury concentration of 5.1 mg/kg, which exceeded the HDOH Tier 1 EAL for mercury of 4.7 mg/kg. The surface soil samples from the remaining three field area DUs did not contain COPC concentrations exceeding the HDOH Tier 1 EALs.

The HDOH Tier 1 EAL for mercury is the direct contact EAL for residential sites, which is conservatively based on a noncarcinogen Hazard Quotient (HQ) of 0.2 to be protective of cumulative health risks potentially posed by sites with multiple noncarcinogens (up to a total of five). However, soil in RHS-DU-29 contained only one noncarcinogen at a concentration exceeding its HDOH Tier 1 EAL. As such, the direct contact EAL calculated using a HQ of 1.0 of 23 mg/kg may be used (see Table I-1 of the Technical Guidance Manual [TGM], HDOH, 2012). The mercury concentration of 5.1 mg/kg measured in RHS-DU-29 does not exceed the direct contact HDOH EAL for residential sites for mercury calculated using a HQ of 1.0 of 23 mg/kg. In other words, mercury surface soil concentrations measured in RHS-DU-29 are not expected to pose a risk for noncancer health effects assuming the strictest exposure parameters (i.e., residential).

Based on the laboratory analytical results, Bureau Veritas recommends the following:

- The stockpile DUs (RHS-DU-14 through RHD-DU-25) should be profiled and properly disposed.
- Conduct an additional site characterization investigation to further delineate the horizontal extent of contamination and the vertical extent of contamination.



- After the additional site characterization is conducted, an EHE should be prepared to assess the potential environmental hazards posed by the soil contamination.
- Develop remedial alternatives based on the results of the EHE.

## 7.0 LIMITATIONS

This report is for the exclusive use of the State of Hawaii Department of Education and no other party shall have any right to rely on any service provided by Bureau Veritas without prior written consent. The information and opinions expressed in this report are given in response to a limited assignment and should be considered and implemented only in light of that assignment. The services provided by Bureau Veritas in completing this project were consistent with normal standards of the profession. No other warranty, expressed or implied, is made. Bureau Veritas will not distribute or publish this report without consent except as required by law or court order.

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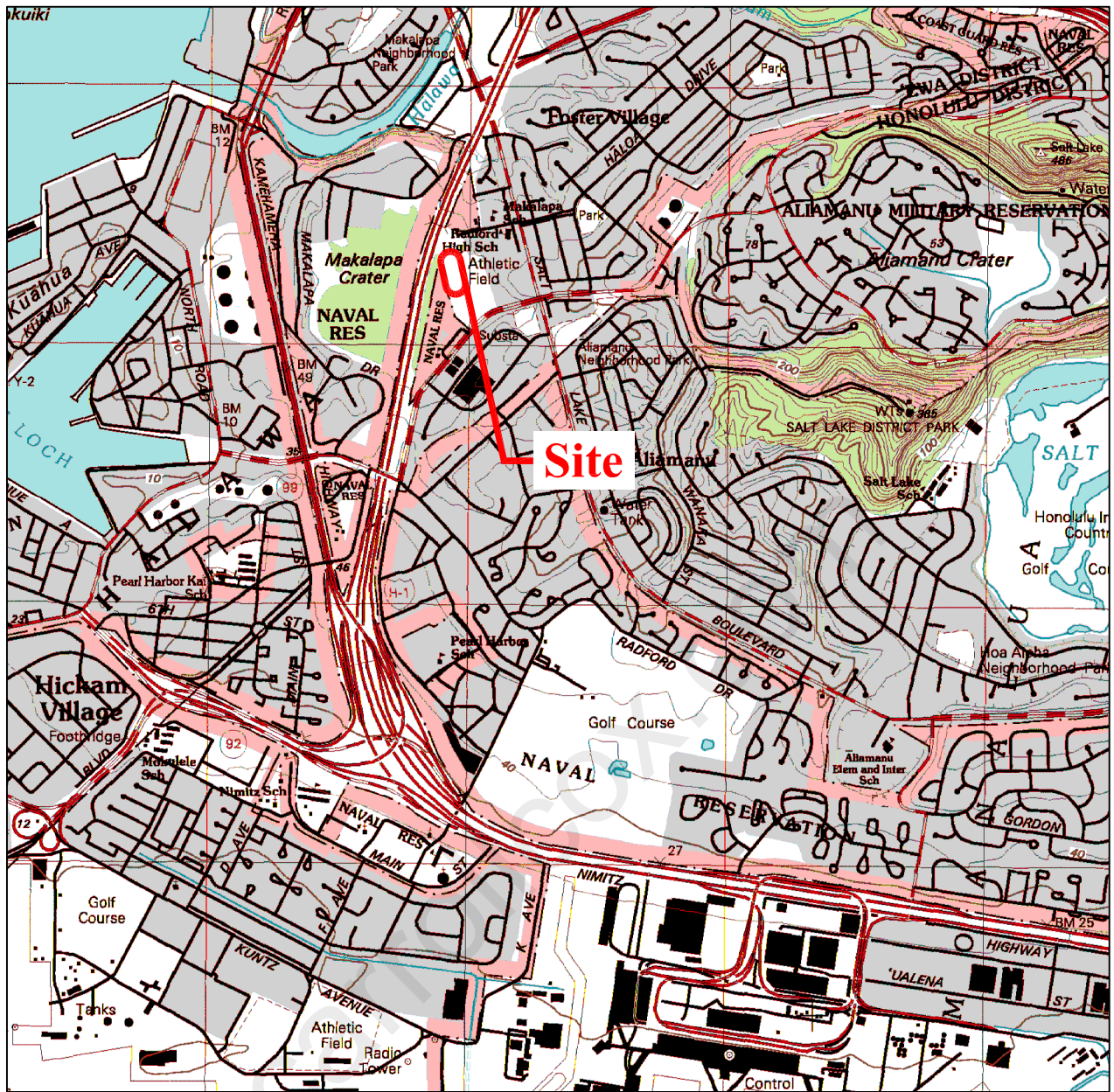
Van den Berg, M., Birnbaum, L.S., Denison, M., De Vito, M., Farland, W., Feeley, M., Fiedler, H., Hakansson, H., Hanberg, A., Haws, L., Rose, M., Safe, S., Schrenk, D., Tohyama, C., Tritscher, A., Tuomisto, J., Tysklind, M., Walker, N., Peterson, R.E.; 2006. Review: The 2005 World Health Organization Reevaluation of Human and Mammalian Toxic Equivalency Factors for Dioxins and Dioxin-Like Compounds. *Toxicological Sciences* 93(2), 223-241.



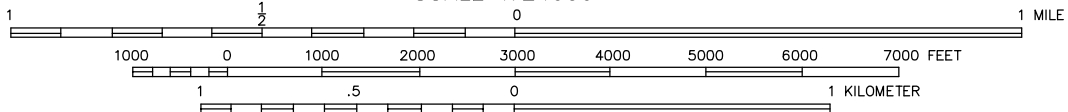
## FIGURES

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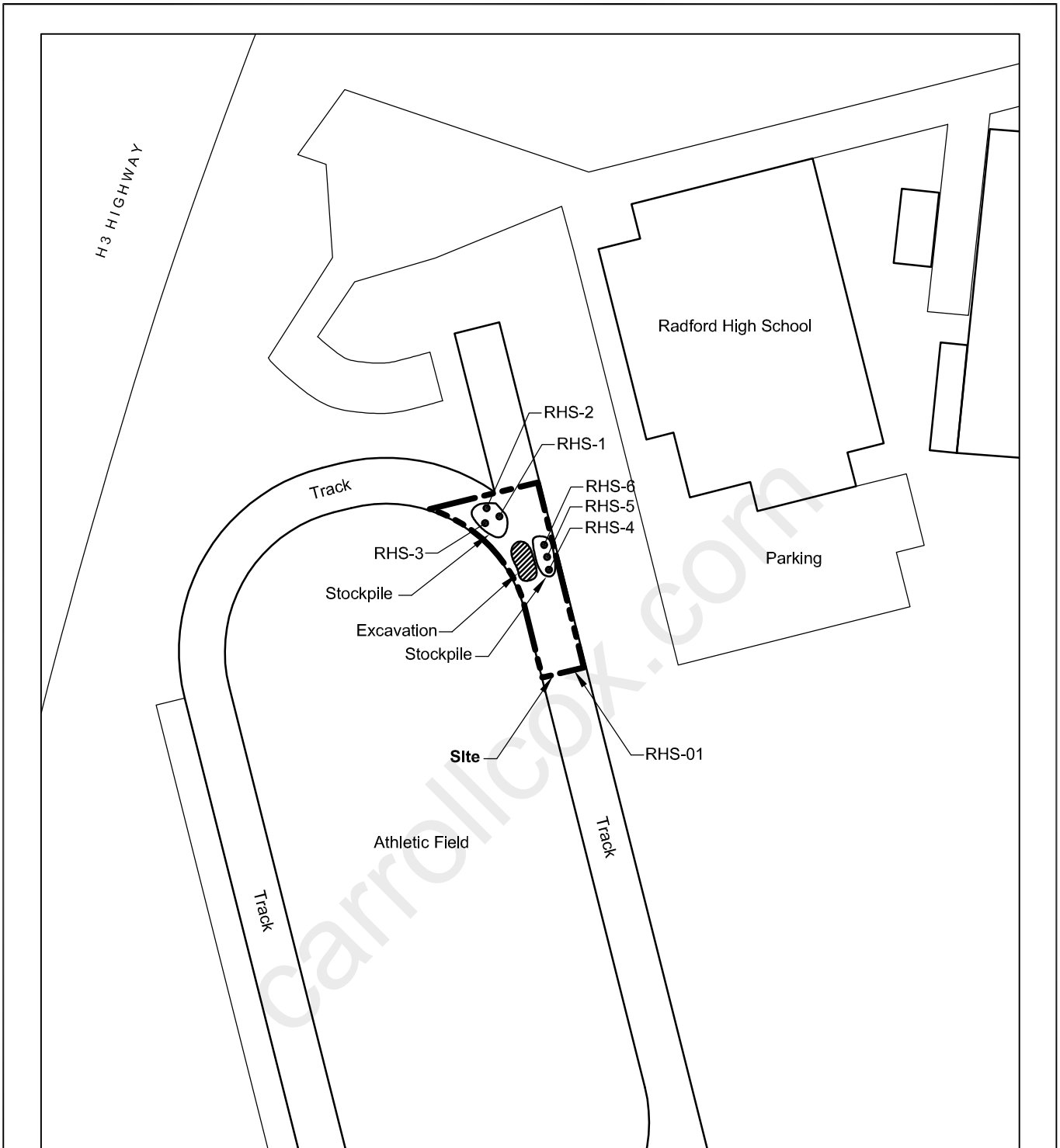



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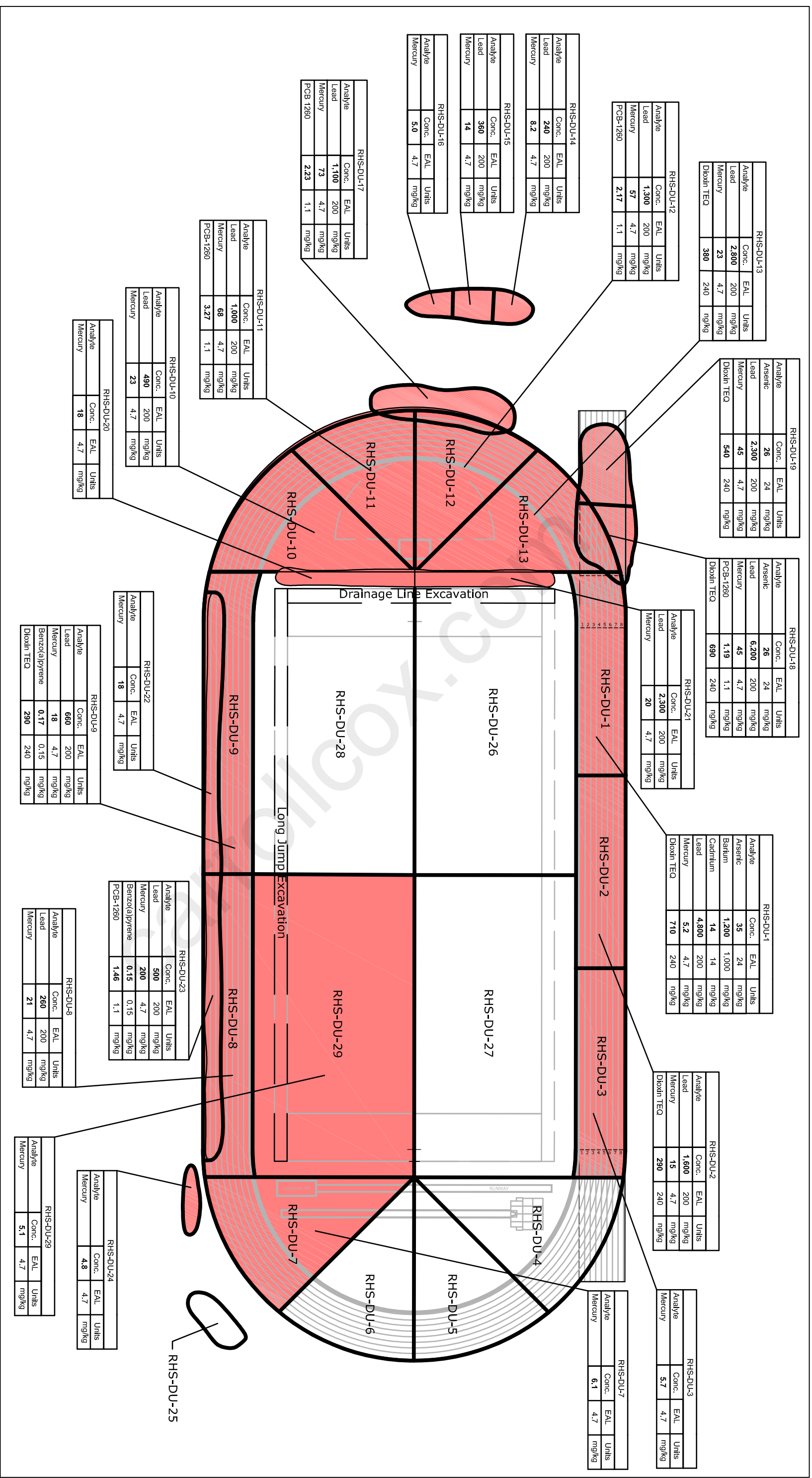
Portion of 7.5-minute Series (Topographic) Maps  
 United States Department of Interior  
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 1998

|  |                                 |  |                               |
|--|---------------------------------|--|-------------------------------|
| <br><b>BUREAU VERITAS</b> | Project No.:<br>17012-012148.00 | Title:<br><b>Site Location Map</b>   | <b>FIGURE</b><br><br><b>1</b> |
|  | Date:<br>3/20/14                | Location:<br>Radford High School<br>4361 Salt Lake Blvd.<br>Honolulu, Oahu, Hawaii 96818 |                               |
|  | Revised By:<br>JC               | Client:<br>State of Hawaii Department of Education                                       |                               |
|  | Checked By:<br>KG               |  |                               |

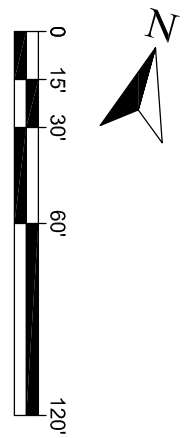


|  |                                 |  |                               |
|--|---------------------------------|--|-------------------------------|
| <br><b>BUREAU VERITAS</b> | Project No.:<br>17012-012148.00 | Title:<br><b>Initial Sample Area</b>   | <b>FIGURE</b><br><br><b>2</b> |
|  | Date:<br>3/20/14                | Location:<br>Radford High School<br>4361 Salt Lake Blvd.<br>Honolulu, Oahu, Hawaii 96818 |                               |
|  | Revised By:<br>JC               | Client:<br>State of Hawaii Department of Education                                       |                               |
|  | Checked By:<br>KG               |  |                               |





**Legend:**  
 Analyte Detected Greater than the HDOH Unrestricted EALs or the HDOH 2010 TEQ Dioxin SALs



Project No.: 17012-012148.00  
 Date: 3/20/14  
 Revised By: JC  
 Checked By: KG  
 Decision Unit Area Results Exceeding HDOH Tier 1 EALS  
 Radford High School  
 4361 Salt Lake Blvd.  
 Honolulu, Oahu, Hawaii 96818  
 State of Hawaii Department of Education

FIGURE 3





## TABLES

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**Table 1**  
**Analytical Results for Initial Soil Sample**  
**State of Hawaii Department of Education**  
**Radford High School Track and Field Investigation**  
**Honolulu, Oahu, Hawaii**

**Project No: 17012-012148.00 / Task 048**

| Sample ID:  | <b>RHS-01</b> | HDOH   |
|---|---------------|--------|
| Date Sampled:   | 12/20/2013    | Tier 1 |
| Analytes  | Units:        | EAL    |
|   | mg/kg         |        |
| <b>Total Petroleum Hydrocarbons (TPH) / EPA Method 8015B</b>      |               |        |
| TPH-Diesel Range Organics   | 24            | 100    |
| TPH-Residual Range Organics                                       | 23            | 500    |
| <b>Total Metals / EPA Method 6010B/7471</b>                       |               |        |
| Arsenic   | <b>43*</b>    | 24     |
| Barium  | 940           | 1,000  |
| Cadmium   | <b>26*</b>    | 14     |
| Chromium  | 210           | 1,100  |
| Lead  | <b>5,300*</b> | 200    |
| Selenium  | ND< 10        | 78     |
| Silver  | 14            | 78     |
| Mercury   | 1.1           | 4.7    |
| <b>Semi-volatile Organic Compounds (SVOCs) / EPA Method 8270C</b> |               |        |
| 1,2,4-Trichlorobenzene  | ND< 0.69      | 0.098  |
| 1,2-Dichlorobenzene   | ND< 0.69      | 0.75   |
| 1,3-Dichlorobenzene   | ND< 0.69      | 0.57   |
| 1,4-Dichlorobenzene   | ND< 0.69      | 0.047  |
| 1-Methylnaphthalene   | ND< 0.73      | 1.8    |
| 2,4,5-Trichlorophenol   | ND< 0.69      | 29     |
| 2,4,6-Trichlorophenol   | ND< 0.69      | 1.8    |
| 2,4-Dichlorophenol  | ND< 0.69      | 0.025  |
| 2,4-Dimethylphenol  | ND< 0.69      | 9.9    |
| 2,4-Dinitrophenol   | ND< 1.4       | 5.6    |
| 2,4-Dinitrotoluene  | ND< 0.69      | 0.021  |
| 2,6-Dinitrotoluene  | ND< 0.69      | 3.6    |
| 2-Chloronaphthalene   | ND< 0.69      | NS     |
| 2-Chlorophenol  | ND< 0.69      | 0.0092 |
| 2-Methylnaphthalene   | 0.25          | 4.1    |
| 2-Methylphenol  | ND< 0.69      | NS     |
| 2-Nitroaniline  | ND< 0.69      | NS     |
| 2-Nitrophenol   | ND< 0.69      | NS     |
| 3,3'-Dichlorobenzidine  | ND< 1.7       | 0.079  |
| 3 & 4 Methylphenol  | ND< 0.69      | NS     |
| 3-Nitroaniline  | ND< 0.69      | NS     |
| 4,6-Dinitro-2-methylphenol  | ND< 0.88      | NS     |
| 4-Bromophenyl phenyl ether  | ND< 0.69      | NS     |

Notes on page 3



Table 1 (continued)

| Total Metals / EPA Method 6010B/7471 (Continued) |          |          |
|--|----------|----------|
| 4-Chloro-3-methylphenol                          | ND< 0.69 | NS       |
| 4-Chloroaniline                                  | ND< 0.69 | 0.0063   |
| 4-Chlorophenyl phenyl ether                      | ND< 0.69 | NS       |
| 4-Nitroaniline                                   | ND< 1.7  | NS       |
| 4-Nitrophenol                                    | ND< 1.7  | NS       |
| Acenaphthene                                     | ND< 0.69 | 120      |
| Acenaphthylene                                   | ND< 0.69 | 100      |
| Anthracene                                       | ND< 0.69 | 4.3      |
| Benzo[a]anthracene                               | ND< 0.69 | 1.5      |
| Benzo[a]pyrene                                   | ND< 0.69 | 0.15     |
| Benzo[b]fluoranthene                             | 0.15     | 1.5      |
| Benzo[g,h,i]perylene                             | ND< 0.69 | 35       |
| Benzo[k]fluoranthene                             | ND< 0.69 | 15       |
| Benzoic acid                                     | ND< 1.7  | NS       |
| Benzyl alcohol                                   | ND< 0.69 | NS       |
| Bis(2-chloroisopropyl)ether                      | ND< 0.69 | NS       |
| Bis(2-chloroethoxy)methane                       | ND< 0.69 | NS       |
| Bis(2-chloroethyl)ether                          | ND< 0.69 | 0.000064 |
| Bis(2-ethylhexyl) phthalate                      | ND< 0.69 | 35       |
| Butyl benzyl phthalate                           | ND< 0.69 | NS       |
| Carbazole  | ND< 0.69 | NS       |
| Chrysene   | ND< 0.69 | 30       |
| Dibenzo(a,h)anthracene                           | ND< 0.88 | 0.15     |
| Dibenzofuran                                     | ND< 0.69 | NS       |
| Diethyl phthalate                                | ND< 0.69 | 16       |
| Dimethyl phthalate                               | ND< 0.69 | 22       |
| Di-n-butyl phthalate                             | ND< 0.69 | NS       |
| Di-n-octyl phthalate                             | ND< 0.69 | NS       |
| Fluoranthene                                     | ND< 0.69 | 460      |
| Fluorene   | ND< 0.69 | 100      |
| Hexachlorobenzene                                | ND< 0.69 | 0.30     |
| Hexachlorobutadiene                              | ND< 0.69 | 0.18     |
| Hexachlorocyclopentadiene                        | ND< 1.7  | NS       |
| Hexachloroethane                                 | ND< 0.69 | 0.27     |
| Indeno[1,2,3-cd]pyrene                           | 0.49     | 1.5      |
| Isophorone                                       | ND< 0.69 | 0.77     |
| Naphthalene                                      | 2.7      | 4.4      |
| Nitrobenzene                                     | ND< 0.69 | 0.0046   |
| N-Nitrosodi-n-propylamine                        | ND< 0.52 | NS       |
| N-Nitrosodiphenylamine                           | ND< 0.69 | NS       |
| Pentachlorophenol                                | ND< 1.7  | 0.82     |
| Phenanthrene                                     | ND< 0.69 | 440      |
| Phenol   | ND< 0.69 | 0.16     |
| Pyrene   | ND< 0.69 | 44       |

Notes on page 3



Table 1 (continued)

| <b>Polychlorinated Biphenyls (PCBs) / EPA Method 8082</b> |                    |         |
|---|--------------------|---------|
| PCB-1016  | ND< 0.026          | 1.1     |
| PCB-1221  | ND< 0.026          | 1.1     |
| PCB-1232  | ND< 0.026          | 1.1     |
| PCB-1242  | ND< 0.026          | 1.1     |
| PCB-1248  | ND< 0.026          | 1.1     |
| PCB-1254  | ND< 0.026          | 1.1     |
| PCB-1260  | 0.094              | 1.1     |
| <b>Organochlorine Pesticides / EPA Method 8081A</b>       |                    |         |
| p,p'-DDD  | ND< 0.0026         | 2.0     |
| p,p'-DDE  | 0.0046             | 1.4     |
| p,p'-DDT  | 0.0045             | 1.7     |
| Aldrin  | ND< 0.0026         | 0.92    |
| alpha-BHC   | ND< 0.0026         | NS      |
| beta-BHC  | ND< 0.0026         | NS      |
| delta-BHC   | ND< 0.0052         | NS      |
| Chlordane (technical)                                     | ND< 0.026          | 16      |
| Dieldrin  | 0.0035             | 1.5     |
| Endosulfan I  | ND< 0.0026         | 18      |
| Endosulfan II   | ND< 0.0026         | NS      |
| Endosulfan sulfate  | ND< 0.0052         | NS      |
| Endrin  | 0.00086            | 3.7     |
| Endrin aldehyde   | 0.0035             | NS      |
| Endrin ketone   | ND< 0.0026         | NS      |
| gamma-BHC (Lindane)                                       | ND< 0.0026         | 0.075   |
| Heptachlor  | ND< 0.0026         | 0.11    |
| Heptachlor epoxide  | ND< 0.0026         | 0.053   |
| Methoxychlor  | ND< 0.0026         | 16      |
| Toxaphene   | ND< 0.10           | 0.44    |
| <b>Dioxins and Furans / EPA Method 8290</b>               |                    |         |
| TEQ   | <b>0.00201939*</b> | 0.00024 |

Notes:

- Bold\*** The value exceeds the State of Hawaii Department of Health (HDOH) Tier 1 Environmental Action Levels (EAL) for unrestricted land use.
- HDOH Tier 1 EAL The State of Hawaii Department of Health (HDOH) Tier 1 Environmental Action Levels (EAL) for sites where groundwater is a current or potential source of drinking water, and the site is greater than 150 meters from a surface water body. (Fall 2011, updated January 2012).
- Italics* Reporting Limit is above the HDOH Tier 1 EAL.
- mg/kg milligrams per kilogram.
- ND< No detectable concentration. The number following the "less than" symbol is the laboratory reporting limit.
- NS No standard. The HDOH has not established a Tier 1 EAL for this analyte.
- TEQ Toxic equivalency. Calculating the total TEQ of a mixture involves multiplying the concentrations of individual congeners by their respective TEF, then adding the individual TEQ's to obtain a total TEQ concentration for the mixture.



**Table 2**  
**Dioxin Toxic Equivalency (TEQ) Calculations for Initial Soil Sample**  
**State of Hawaii Department of Education**  
**Radford High School Track and Field Investigation**  
**Honolulu, Oahu, Hawaii**

**Project No: 17012-012148.00 / Task 048**

| PCDD/F Analytes          | TEF    | RHS-01                    |                           |
|--------------------------|--------|---------------------------|---------------------------|
|                          |        | Analytical Results (pg/g) | TEF Adjusted Conc. (pg/g) |
| 2,3,7,8-TCDD             | 1      | 88                        | 88.0                      |
| 1,2,3,7,8-PeCDD          | 1      | 340                       | 340.00                    |
| 1,2,3,4,7,8-HxCDD        | 0.1    | 250                       | 25.00                     |
| 1,2,3,6,7,8-HxCDD        | 0.1    | 420                       | 42.00                     |
| 1,2,3,7,8,9-HxCDD        | 0.1    | 320                       | 32.00                     |
| 1,2,3,4,6,7,8-HpCDD      | 0.01   | 2,400                     | 24.0                      |
| OCDD                     | 0.0003 | 5,600                     | 1.68                      |
| 2,3,7,8-TCDF             | 0.1    | 1,400                     | 140.00                    |
| 1,2,3,7,8-PeCDF          | 0.03   | 1,100                     | 33.000                    |
| 2,3,4,7,8-PeCDF          | 0.3    | 2,300                     | 690.00                    |
| 1,2,3,4,7,8-HxCDF        | 0.1    | 1,900                     | 190.00                    |
| 1,2,3,6,7,8-HxCDF        | 0.1    | 1,400                     | 140.00                    |
| 1,2,3,7,8,9-HxCDF        | 0.1    | ND<100                    | 5.00                      |
| 2,3,4,6,7,8-HxCDF        | 0.1    | 1,900                     | 190.00                    |
| 1,2,3,4,6,7,8-HpCDF      | 0.01   | 7,600                     | 76.00                     |
| 1,2,3,4,7,8,9-HpCDF      | 0.01   | 220                       | 2.200                     |
| OCDF                     | 0.0003 | 1,700                     | 0.510                     |
| <b>Total TEQ (ng/kg)</b> |        |                           | <b>2,019.39</b>           |

**Notes:**

- ND< Analyte not detected. The value after the '<' is the laboratory Reporting Limit (RL)
- ng/kg nanograms per kilogram (equivalent to parts per trillion [ppt])
- TEF Toxicity equivalency factor. TEF is the ratio of the toxicity of one of the compounds to the toxicity of the two most toxic compounds.
- TEQ Toxic equivalency. Calculating the total TEQ of a mixture involves multiplying the concentrations of individual congeners by their respective TEF, then adding the individual TEQ's to obtain a total TEQ concentration for the mixture.
- HxCDD Hexachlorodibenzo-*p*-dioxin
- HxCDF Hexachlorodibenzofuran
- OCDD Octachlorodibenzo-*p*-dioxin
- OCDF Octachlorodibenzofuran
- PCDD Pentachlorodibenzo-*p*-dioxin
- PCDF Pentachlorodibenzofuran
- pg/g Picograms per liter (parts per quadrillion [ppq])
- TCDD Tetrachlorodibenzo-*p*-dioxin
- TCDF Tetrachlorodibenzofuran

**Table 3**  
**Analytical Results for Soil Samples from Decision Unit Areas**  
**State of Hawaii Department of Education**  
**Radford High School Track and Field Investigation**  
**Honolulu, Oahu, Hawaii**

Project No: 17012-012148.00 / Task 048

| Sample ID:<br>Date Sampled:<br>Units:                             | RHS-DU-1<br>2/4/2014<br>mg/kg | RHS-DU-2<br>2/4/2014<br>mg/kg | RHS-DU-3<br>2/4/2014<br>mg/kg | RHS-DU-4<br>2/4/2014<br>mg/kg | RHS-DU-5<br>2/4/2014<br>mg/kg | RHS-DU-6<br>2/4/2014<br>mg/kg | RHS-DU-7<br>2/4/2014<br>mg/kg | RHS-DU-8<br>2/4/2014<br>mg/kg | HDOH Tier 1<br>EAL<br>mg/kg |
|---|-------------------------------|-------------------------------|-------------------------------|-------------------------------|-------------------------------|-------------------------------|-------------------------------|-------------------------------|-----------------------------|
| <b>Dioxin / EPA Method 8290</b>                                   |                               |                               |                               |                               |                               |                               |                               |                               |                             |
| Total Dioxin Toxic Equivalent (TEQ) □                             | 710*                          | 290*                          | 25                            | 35                            | 33                            | 63                            | 110                           | 160                           | 240 □                       |
| <b>Total Petroleum Hydrocarbons (TPH) / EPA Method 8015B</b>      |                               |                               |                               |                               |                               |                               |                               |                               |                             |
| TPH-Diesel Range Organics   | ND< 50                        | ND< 50                        | ND< 50                        | ND< 50                        | ND< 50                        | ND< 50                        | ND< 50                        | ND< 50                        | 500                         |
| TPH-Residual Range Organics                                       | ND< 100                       | ND< 100                       | ND< 100                       | ND< 100                       | ND< 100                       | ND< 100                       | ND< 100                       | ND< 100                       | 500                         |
| <b>Total Metals / EPA Method 6010B/7471</b>                       |                               |                               |                               |                               |                               |                               |                               |                               |                             |
| Arsenic   | 35*                           | 11                            | 6.0                           | 7.5                           | 6.1                           | 7.6                           | 7.3                           | 6.8                           | 24                          |
| Barium  | 1,200*                        | 550                           | 410                           | 460                           | 280                           | 200                           | 280                           | 320                           | 1,000                       |
| Cadmium   | 14*                           | 5.9                           | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 1.0                           | 1.3                           | 14                          |
| Chromium  | 300                           | 150                           | 160                           | 170                           | 180                           | 240                           | 230                           | 170                           | 1,100                       |
| Lead  | 4,800*                        | 1,600*                        | 85                            | 140                           | 61                            | 66                            | 190                           | 260*                          | 200                         |
| Selenium  | ND< 20                        | ND< 20                        | ND< 20                        | ND< 20                        | ND< 20                        | ND< 20                        | ND< 20                        | ND< 20                        | 78                          |
| Silver  | ND< 20                        | ND< 20                        | ND< 20                        | ND< 20                        | ND< 20                        | ND< 20                        | ND< 20                        | ND< 20                        | 78                          |
| Mercury   | 5.2*                          | 15*                           | 5.7*                          | 3.5                           | 2.9                           | 3.3                           | 6.1*                          | 21*                           | 4.7                         |
| <b>Polychlorinated Biphenyls (PCBs) / EPA Method 8082</b>         |                               |                               |                               |                               |                               |                               |                               |                               |                             |
| PCB-1016  | ND< 0.10                      | ND< 0.10                      | ND< 0.10                      | ND< 0.10                      | ND< 0.10                      | ND< 0.10                      | ND< 0.10                      | ND< 0.10                      | 1.1                         |
| PCB-1221  | ND< 0.20                      | ND< 0.20                      | ND< 0.20                      | ND< 0.20                      | ND< 0.20                      | ND< 0.20                      | ND< 0.20                      | ND< 0.20                      | 1.1                         |
| PCB-1232  | ND< 0.20                      | ND< 0.20                      | ND< 0.20                      | ND< 0.20                      | ND< 0.20                      | ND< 0.20                      | ND< 0.20                      | ND< 0.20                      | 1.1                         |
| PCB-1242  | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | 1.1                         |
| PCB-1248  | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | 1.1                         |
| PCB-1254  | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | ND< 0.05                      | 1.1                         |
| PCB-1260  | ND< 0.05                      | 0.10                          | 0.21                          | ND< 0.05                      | 0.05                          | ND< 0.05                      | 0.09                          | 0.17                          | 1.1                         |
| <b>Semi-volatile Organic Compounds (SVOCs) / EPA Method 8270C</b> |                               |                               |                               |                               |                               |                               |                               |                               |                             |
| 1,2,4-Trichlorobenzene  | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 0.098                       |
| 1,2-Dichlorobenzene   | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 0.75                        |
| 1,2-Dinitrobenzene  | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| 1,3-Dichlorobenzene   | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 0.57                        |
| 1,3-Dinitrobenzene  | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | NS                          |
| 1,4-Dichlorobenzene   | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 0.047                       |
| 1,4-Dinitrobenzene  | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | NS                          |
| 1-Methylnapthalene  | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 1.8                         |
| 2,3,4,6-Tetrachlorophenol   | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 4.9                         |
| 2,3,5,6-Tetrachlorophenol   | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| 2,4,5-Trichlorophenol   | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | 29                          |
| 2,4,6-Trichlorophenol   | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | 1.8                         |

Notes on page 13





Table 3 (Continued)

| Sample ID:<br>Date Sampled:<br>Units: | RHS-DU-1<br>2/4/2014<br>mg/kg | RHS-DU-2<br>2/4/2014<br>mg/kg | RHS-DU-3<br>2/4/2014<br>mg/kg | RHS-DU-4<br>2/4/2014<br>mg/kg | RHS-DU-5<br>2/4/2014<br>mg/kg | RHS-DU-6<br>2/4/2014<br>mg/kg | RHS-DU-7<br>2/4/2014<br>mg/kg | RHS-DU-8<br>2/4/2014<br>mg/kg | HDOH Tier 1<br>EAL<br>mg/kg |
|---------------------------------------|-------------------------------|-------------------------------|-------------------------------|-------------------------------|-------------------------------|-------------------------------|-------------------------------|-------------------------------|-----------------------------|
| <b>SVOCs (Continued)</b>              |                               |                               |                               |                               |                               |                               |                               |                               |                             |
| 2,4-Dichlorophenol                    | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | 0.025                       |
| 2,4-Dimethylphenol                    | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 9.9                         |
| 2,4-Dinitrophenol                     | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | 5.6                         |
| 2,4-Dinitrotoluene                    | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 0.021                       |
| 2,6-Dinitrotoluene                    | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 3.6                         |
| 2-Chloronaphthalene                   | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| 2-Chlorophenol                        | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 0.0092                      |
| 2-Methylnaphthalene                   | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 4.1                         |
| 2-Methylphenol (o-cresol)             | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| 2-Nitroaniline                        | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | NS                          |
| 2-Nitrophenol                         | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | NS                          |
| 3,4-Methylphenol (m,p-cresol)         | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| 4-Bromophenylphenylether              | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| 4-Chloro-3-methylphenol               | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | NS                          |
| 4-Chloroaniline                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | 0.0063                      |
| 4-Chlorophenylphenylether             | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| 4-Nitroaniline                        | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | NS                          |
| 4-Nitrophenol                         | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | NS                          |
| Acenaphthene                          | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | 120                         |
| Acenaphthylene                        | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | 100                         |
| Aniline                               | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| Anthracene                            | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | 4.3                         |
| Azobenzene                            | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| Benzo(a)anthracene                    | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | 0.11                          | 1.5                         |
| Benzo(a)pyrene                        | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | 0.15                        |
| Benzo(b)fluoranthene                  | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | 0.12                          | 1.5                         |
| Benzo(ghi)perylene                    | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | 35                          |
| Benzo(k)fluoranthene                  | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | 0.20                          | 15                          |
| Benzyl alcohol                        | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| Bis (2-chloroethoxy) methane          | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| Bis (2-chloroethyl) ether             | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 0.000064                    |
| Bis (2-chloroisopropyl) ether         | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | 0.0035                      |
| Bis (2-ethylhexyl) phthalate          | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 35                          |
| Bis(2-ethylhexyl) adipate             | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| Butylbenzylphthalate                  | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| Carbazole                             | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| Chrysene                              | ND< 0.1                       | ND< 0.1                       | 0.13                          | 0.12                          | ND< 0.1                       | 0.20                          | 0.19                          | 0.34                          | 30                          |
| Dibenzo(a,h)anthracene                | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | 0.15                        |
| Dibenzofuran                          | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| Diethylphthalate                      | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 16                          |
| Dimethylphthalate                     | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 22                          |

Notes on page 13





Table 3 (Continued)

| Sample ID:<br>Date Sampled:<br>Units:               | RHS-DU-1<br>2/4/2014<br>mg/kg | RHS-DU-2<br>2/4/2014<br>mg/kg | RHS-DU-3<br>2/4/2014<br>mg/kg | RHS-DU-4<br>2/4/2014<br>mg/kg | RHS-DU-5<br>2/4/2014<br>mg/kg | RHS-DU-6<br>2/4/2014<br>mg/kg | RHS-DU-7<br>2/4/2014<br>mg/kg | RHS-DU-8<br>2/4/2014<br>mg/kg | HDOH Tier 1<br>EAL<br>mg/kg |
|---|-------------------------------|-------------------------------|-------------------------------|-------------------------------|-------------------------------|-------------------------------|-------------------------------|-------------------------------|-----------------------------|
| <b>SVOCs (Continued)</b>                            |                               |                               |                               |                               |                               |                               |                               |                               |                             |
| Di-n-butylphthalate                                 | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| Di-n-octyl phthalate                                | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| Fluoranthene  | ND< 0.1                       | 0.12                          | 0.12                          | ND< 0.1                       | ND< 0.1                       | 0.17                          | 0.21                          | 0.43                          | 460                         |
| Fluorene  | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | 100                         |
| Hexachlorobenzene                                   | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 0.3                         |
| Hexachlorobutadiene                                 | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 0.18                        |
| Hexachlorocyclopentadiene                           | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| Hexacholorethane                                    | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 0.27                        |
| Indeno(1,2,3-cd)pyrene                              | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | 1.5                         |
| Isophorone  | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 0.77                        |
| Naphthalene   | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 4.4                         |
| Nitrobenzene  | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 0.0046                      |
| N-Nitroso-di-n-propylamine                          | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| N-nitrosodiphenylamine                              | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| Pentachlorophenol                                   | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | ND< 5.0                       | 0.82                        |
| Phenanthrene  | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | ND< 0.1                       | 0.12                          | 440                         |
| Phenol  | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | 0.16                        |
| Pyrene  | ND< 0.1                       | 0.15                          | 0.14                          | 0.10                          | ND< 0.1                       | 0.17                          | 0.23                          | 0.42                          | 44                          |
| Pyridine  | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | ND< 1.0                       | NS                          |
| <b>Organochlorine Pesticides / EPA Method 8081A</b> |                               |                               |                               |                               |                               |                               |                               |                               |                             |
| Aldrin  | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | 0.92                        |
| Alpha-BHC   | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | NS                          |
| Alpha-Chlordane                                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | NS                          |
| Beta-BHC  | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | NS                          |
| Chlordane (technical)                               | ND< 0.050                     | ND< 0.050                     | ND< 0.050                     | ND< 0.050                     | ND< 0.050                     | ND< 0.050                     | ND< 0.050                     | ND< 0.050                     | 16                          |
| Delta-BHC   | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | NS                          |
| Dieldrin  | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | 0.017                         | 0.045                         | 1.5                         |
| Endosulfan I  | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | 18                          |
| Endosulfan II                                       | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | NS                          |
| Endosulfan sulfate                                  | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | NS                          |
| Endrin  | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | 3.7                         |
| Endrin aldehyde                                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | NS                          |
| Endrin ketone                                       | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | NS                          |
| Gamma-BHC (Lindane)                                 | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | 0.075                       |
| Gamma-Chlordane                                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | NS                          |
| Heptachlor  | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | 0.11                        |
| Heptachlor epoxide                                  | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | ND< 0.005                     | 0.053                       |
| Methoxychlor  | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | 16                          |
| p,p'-DDD  | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | 2                           |
| p,p'-DDE  | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | 0.010                         | ND< 0.010                     | 1.4                         |
| p,p'-DDT  | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | ND< 0.010                     | 0.011                         | 1.7                         |
| Toxaphene   | ND< 0.050                     | ND< 0.050                     | ND< 0.050                     | ND< 0.050                     | ND< 0.050                     | ND< 0.050                     | ND< 0.050                     | ND< 0.050                     | 0.44                        |

Notes on page 13



Table 3 (Continued)

| Sample ID:<br>Date Sampled:<br>Units:                             | RHS-DU-9<br>2/4/2014<br>mg/kg | RHS-DU-10<br>2/5/2014<br>mg/kg | RHS-DU-11<br>2/5/2014<br>mg/kg | RHS-DU-12<br>2/5/2014<br>mg/kg | RHS-DU-13<br>2/5/2014<br>mg/kg | RHS-DU-14<br>2/4/2014<br>mg/kg | RHS-DU-15<br>2/4/2014<br>mg/kg | HDOH Tier 1<br>EAL<br>mg/kg |
|---|-------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|-----------------------------|
| <b>Dioxin / EPA Method 8290</b>                                   |                               |                                |                                |                                |                                |                                |                                |                             |
| Total Dioxin Toxic Equivalent (TEQ) □                             | <b>290*</b>                   | 110                            | 220                            | 230                            | <b>380*</b>                    | 61                             | 110                            | 240 □                       |
| <b>Total Petroleum Hydrocarbons (TPH) / EPA Method 8015B</b>      |                               |                                |                                |                                |                                |                                |                                |                             |
| TPH-Diesel Range Organics   | ND< 50                        | ND< 50                         | ND< 50                         | ND< 50                         | ND< 50                         | ND< 50                         | ND< 50                         | 500                         |
| TPH-Residual Range Organics                                       | ND< 100                       | ND< 100                        | ND< 100                        | ND< 100                        | ND< 100                        | ND< 100                        | ND< 100                        | 500                         |
| <b>Total Metals / EPA Method 6010B/7471</b>                       |                               |                                |                                |                                |                                |                                |                                |                             |
| Arsenic   | 8.7                           | 12                             | 19                             | 23                             | 19                             | 12                             | 15                             | 24                          |
| Barium  | 320                           | 370                            | 522                            | 670                            | 740                            | 260                            | 270                            | 1,000                       |
| Cadmium   | 1.6                           | 1.7                            | 3.7                            | 7.1                            | 6.8                            | ND< 1.0                        | 1.4                            | 14                          |
| Chromium  | 110                           | 180                            | 200                            | 210                            | 240                            | 220                            | 190                            | 1,100                       |
| Lead  | <b>660*</b>                   | <b>490*</b>                    | <b>1,000*</b>                  | <b>1,300*</b>                  | <b>2,800*</b>                  | <b>240*</b>                    | <b>360*</b>                    | 200                         |
| Selenium  | ND< 20                        | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | 78                          |
| Silver  | ND< 20                        | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | 78                          |
| Mercury   | <b>18*</b>                    | <b>23*</b>                     | <b>68*</b>                     | <b>57*</b>                     | <b>23*</b>                     | <b>8.2*</b>                    | <b>14*</b>                     | 4.7                         |
| <b>Polychlorinated Biphenyls (PCBs) / EPA Method 8082</b>         |                               |                                |                                |                                |                                |                                |                                |                             |
| PCB-1016  | ND< 0.10                      | ND< 0.10                       | ND< 0.10                       | ND< 0.10                       | ND< 0.10                       | ND< 0.10                       | ND< 0.10                       | 1.1                         |
| PCB-1221  | ND< 0.20                      | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | 1.1                         |
| PCB-1232  | ND< 0.20                      | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | 1.1                         |
| PCB-1242  | ND< 0.05                      | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | 1.1                         |
| PCB-1248  | ND< 0.05                      | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | 1.1                         |
| PCB-1254  | ND< 0.05                      | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | 1.1                         |
| PCB-1260  | 0.32                          | 0.66                           | <b>3.27*</b>                   | <b>2.17*</b>                   | 1.00                           | 0.13                           | 0.23                           | 1.1                         |
| <b>Semi-volatile Organic Compounds (SVOCs) / EPA Method 8270C</b> |                               |                                |                                |                                |                                |                                |                                |                             |
| 1,2,4-Trichlorobenzene  | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.098                       |
| 1,2-Dichlorobenzene   | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.75                        |
| 1,2-Dinitrobenzene  | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 1,3-Dichlorobenzene   | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.57                        |
| 1,3-Dinitrobenzene  | ND< 5.0                       | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 1,4-Dichlorobenzene   | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.047                       |
| 1,4-Dinitrobenzene  | ND< 5.0                       | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 1-Methylnaphthalene   | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 1.8                         |
| 2,3,4,6-Tetrachlorophenol   | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 4.9                         |
| 2,3,5,6-Tetrachlorophenol   | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 2,4,5-Trichlorophenol   | ND< 5.0                       | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 29                          |
| 2,4,6-Trichlorophenol   | ND< 5.0                       | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 1.8                         |
| 2,4-Dichlorophenol  | ND< 5.0                       | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 0.025                       |
| 2,4-Dimethylphenol  | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 9.9                         |
| 2,4-Dinitrophenol   | ND< 5.0                       | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 5.6                         |
| 2,4-Dinitrotoluene  | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.021                       |
| 2,6-Dinitrotoluene  | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 3.6                         |

Notes on page 13



Table 3 (Continued)

| Sample ID:<br>Date Sampled:<br>Units: | RHS-DU-9<br>2/4/2014<br>mg/kg | RHS-DU-10<br>2/5/2014<br>mg/kg | RHS-DU-11<br>2/5/2014<br>mg/kg | RHS-DU-12<br>2/5/2014<br>mg/kg | RHS-DU-13<br>2/5/2014<br>mg/kg | RHS-DU-14<br>2/4/2014<br>mg/kg | RHS-DU-15<br>2/4/2014<br>mg/kg | HDOH Tier 1<br>EAL<br>mg/kg |
|---------------------------------------|-------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|-----------------------------|
| <b>SVOCs (Continued)</b>              |                               |                                |                                |                                |                                |                                |                                |                             |
| 2-Chloronaphthalene                   | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 2-Chlorophenol                        | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.0092                      |
| 2-Methylnaphthalene                   | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 4.1                         |
| 2-Methylphenol (o-cresol)             | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 2-Nitroaniline                        | ND< 5.0                       | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 2-Nitrophenol                         | ND< 5.0                       | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 3,4-Methylphenol (m,p-cresol)         | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 4-Bromophenylphenylether              | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 4-Chloro-3-methylphenol               | ND< 5.0                       | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 4-Chloroaniline                       | ND< 5.0                       | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 0.0063                      |
| 4-Chlorophenylphenylether             | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 4-Nitroaniline                        | ND< 5.0                       | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 4-Nitrophenol                         | ND< 5.0                       | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| Acenaphthene                          | ND< 0.1                       | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 120                         |
| Acenaphthylene                        | ND< 0.1                       | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 100                         |
| Aniline                               | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Anthracene                            | ND< 0.1                       | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 4.3                         |
| Azobenzene                            | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Benzo(a)anthracene                    | 0.26                          | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 1.5                         |
| Benzo(a)pyrene                        | 0.17*                         | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 0.15                        |
| Benzo(b)fluoranthene                  | 0.33                          | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 1.5                         |
| Benzo(ghi)perylene                    | 0.35                          | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 35                          |
| Benzo(k)fluoranthene                  | 0.35                          | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 15                          |
| Benzyl alcohol                        | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Bis (2-chloroethoxy) methane          | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Bis (2-chloroethyl) ether             | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.000064                    |
| Bis (2-chloroisopropyl) ether         | ND< 5.0                       | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 0.0035                      |
| Bis (2-ethylhexyl) phthalate          | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 35                          |
| Bis(2-ethylhexyl) adipate             | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Butylbenzylphthalate                  | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Carbazole                             | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Chrysene                              | 0.58                          | 0.14                           | 0.20                           | 0.13                           | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 30                          |
| Dibenzo(a,h)anthracene                | ND< 0.1                       | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 0.15                        |
| Dibenzofuran                          | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Diethylphthalate                      | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 16                          |
| Dimethylphthalate                     | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 22                          |
| Di-n-butylphthalate                   | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Di-n-octyl phthalate                  | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Fluoranthene                          | 0.50                          | 0.16                           | 0.18                           | 0.11                           | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 460                         |
| Fluorene                              | ND< 0.1                       | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 100                         |

Notes on page 13



Table 3 (Continued)

| Sample ID:<br>Date Sampled:<br>Units:               | RHS-DU-9<br>2/4/2014<br>mg/kg | RHS-DU-10<br>2/5/2014<br>mg/kg | RHS-DU-11<br>2/5/2014<br>mg/kg | RHS-DU-12<br>2/5/2014<br>mg/kg | RHS-DU-13<br>2/5/2014<br>mg/kg | RHS-DU-14<br>2/4/2014<br>mg/kg | RHS-DU-15<br>2/4/2014<br>mg/kg | HDOH Tier 1<br>EAL<br>mg/kg |
|---|-------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|-----------------------------|
| <b>SVOCs (Continued)</b>                            |                               |                                |                                |                                |                                |                                |                                |                             |
| Hexachlorobenzene                                   | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.3                         |
| Hexachlorobutadiene                                 | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.18                        |
| Hexachlorocyclopentadiene                           | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Hexachlorethane                                     | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.27                        |
| Indeno(1,2,3-cd)pyrene                              | 0.35                          | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 1.5                         |
| Isophorone  | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.77                        |
| Naphthalene   | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 4.4                         |
| Nitrobenzene  | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.0046                      |
| N-Nitroso-di-n-propylamine                          | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| N-nitrosodiphenylamine                              | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Pentachlorophenol                                   | ND< 5.0                       | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 0.82                        |
| Phenanthrene  | 0.13                          | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 440                         |
| Phenol  | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.16                        |
| Pyrene  | 0.50                          | 0.18                           | 0.23                           | 0.14                           | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 44                          |
| Pyridine  | ND< 1.0                       | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| <b>Organochlorine Pesticides / EPA Method 8081A</b> |                               |                                |                                |                                |                                |                                |                                |                             |
| Aldrin  | ND< 0.005                     | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | 0.92                        |
| Alpha-BHC   | ND< 0.005                     | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | NS                          |
| Alpha-Chlordane                                     | ND< 0.005                     | 0.006                          | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | 0.006                          | ND< 0.005                      | NS                          |
| Beta-BHC  | ND< 0.005                     | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | NS                          |
| Chlordane (technical)                               | ND< 0.050                     | ND< 0.050                      | 0.113                          | 0.127                          | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | 16                          |
| Delta-BHC   | ND< 0.005                     | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | NS                          |
| Dieldrin  | ND< 0.010                     | 0.027                          | 0.034                          | ND< 0.010                      | ND< 0.010                      | 0.014                          | ND< 0.010                      | 1.5                         |
| Endosulfan I  | ND< 0.005                     | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | 18                          |
| Endosulfan II                                       | ND< 0.010                     | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | NS                          |
| Endosulfan sulfate                                  | ND< 0.010                     | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | NS                          |
| Endrin  | ND< 0.010                     | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | 3.7                         |
| Endrin aldehyde                                     | ND< 0.010                     | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | NS                          |
| Endrin ketone                                       | ND< 0.010                     | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | NS                          |
| Gamma-BHC (Lindane)                                 | ND< 0.005                     | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | 0.075                       |
| Gamma-Chlordane                                     | ND< 0.005                     | 0.006                          | 0.024                          | 0.014                          | 0.022                          | ND< 0.005                      | ND< 0.005                      | NS                          |
| Heptachlor  | ND< 0.005                     | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | 0.11                        |
| Heptachlor epoxide                                  | ND< 0.005                     | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | 0.053                       |
| Methoxychlor  | ND< 0.010                     | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | 16                          |
| p,p'-DDD  | ND< 0.010                     | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | 2                           |
| p,p'-DDE  | ND< 0.010                     | ND< 0.010                      | ND< 0.010                      | 0.036                          | 0.016                          | 0.012                          | ND< 0.010                      | 1.4                         |
| p,p'-DDT  | 0.021                         | 0.040                          | 0.216                          | 0.174                          | 0.051                          | 0.012                          | 0.021                          | 1.7                         |
| Toxaphene   | ND< 0.050                     | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | 0.44                        |

Notes on page 13

Table 3 (Continued)

| Sample ID:<br>Date Sampled:<br>Units:                             | RHS-DU-16<br>2/4/2014<br>mg/kg | RHS-DU-17<br>2/4/2014<br>mg/kg | RHS-DU-18<br>2/4/2014<br>mg/kg | RHS-DU-19<br>2/4/2014<br>mg/kg | RHS-DU-20<br>2/4/2014<br>mg/kg | RHS-DU-21<br>2/4/2014<br>mg/kg | RHS-DU-22<br>2/4/2014<br>mg/kg | HDOH Tier 1<br>EAL<br>mg/kg |
|---|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|-----------------------------|
| <b>Dioxin / EPA Method 8290</b>                                   |                                |                                |                                |                                |                                |                                |                                |                             |
| Total Dioxin Toxic Equivalent (TEQ) □                             | 44                             | 230                            | 690*                           | 540*                           | 19                             | 18                             | 39                             | 240 □                       |
| <b>Total Petroleum Hydrocarbons (TPH) / EPA Method 8015B</b>      |                                |                                |                                |                                |                                |                                |                                |                             |
| TPH-Diesel Range Organics   | ND< 50                         | ND< 50                         | ND< 50                         | ND< 50                         | ND< 50                         | ND< 50                         | ND< 50                         | 500                         |
| TPH-Residual Range Organics                                       | ND< 100                        | ND< 100                        | ND< 100                        | ND< 100                        | ND< 100                        | ND< 100                        | ND< 100                        | 500                         |
| <b>Total Metals / EPA Method 6010B/7471</b>                       |                                |                                |                                |                                |                                |                                |                                |                             |
| Arsenic   | 5.4                            | 23                             | 26*                            | 26*                            | 6.4                            | 18                             | 6.9                            | 24                          |
| Barium  | 220                            | 590                            | 710                            | 610                            | 290                            | 810                            | 320                            | 1,000                       |
| Cadmium   | ND< 1.0                        | 4.5                            | 11                             | 7.8                            | ND< 1.0                        | 12                             | ND< 1.0                        | 14                          |
| Chromium  | 150                            | 210                            | 170                            | 160                            | 190                            | 240                            | 130                            | 1,100                       |
| Lead  | 95                             | 1,100*                         | 6,200*                         | 2,300*                         | 120                            | 2,300*                         | 110                            | 200                         |
| Selenium  | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | 78                          |
| Silver  | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | 78                          |
| Mercury   | 5.0*                           | 73*                            | 45*                            | 45*                            | 18*                            | 20*                            | 18*                            | 4.7                         |
| <b>Polychlorinated Biphenyls (PCBs) / EPA Method 8082</b>         |                                |                                |                                |                                |                                |                                |                                |                             |
| PCB-1016  | ND< 0.10                       | ND< 0.10                       | ND< 0.10                       | ND< 0.10                       | ND< 0.10                       | ND< 0.10                       | ND< 0.10                       | 1.1                         |
| PCB-1221  | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | 1.1                         |
| PCB-1232  | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | 1.1                         |
| PCB-1242  | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | 1.1                         |
| PCB-1248  | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | 1.1                         |
| PCB-1254  | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | 1.1                         |
| PCB-1260  | ND< 0.05                       | 2.23*                          | 1.19*                          | 0.47                           | 0.29                           | 0.20                           | 0.25                           | 1.1                         |
| <b>Semi-volatile Organic Compounds (SVOCs) / EPA Method 8270C</b> |                                |                                |                                |                                |                                |                                |                                |                             |
| 1,2,4-Trichlorobenzene  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.098                       |
| 1,2-Dichlorobenzene   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.75                        |
| 1,2-Dinitrobenzene  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 1,3-Dichlorobenzene   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.57                        |
| 1,3-Dinitrobenzene  | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 1,4-Dichlorobenzene   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.047                       |
| 1,4-Dinitrobenzene  | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 1-Methylnaphthalene   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 1.8                         |
| 2,3,4,6-Tetrachlorophenol   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 4.9                         |
| 2,3,5,6-Tetrachlorophenol   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 2,4,5-Trichlorophenol   | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 29                          |
| 2,4,6-Trichlorophenol   | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 1.8                         |
| 2,4-Dichlorophenol  | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 0.025                       |
| 2,4-Dimethylphenol  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 9.9                         |
| 2,4-Dinitrophenol   | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 5.6                         |
| 2,4-Dinitrotoluene  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.021                       |
| 2,6-Dinitrotoluene  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 3.6                         |

Notes on page 13





Table 3 (Continued)

| Sample ID:<br>Date Sampled:<br>Units: | RHS-DU-16<br>2/4/2014<br>mg/kg | RHS-DU-17<br>2/4/2014<br>mg/kg | RHS-DU-18<br>2/4/2014<br>mg/kg | RHS-DU-19<br>2/4/2014<br>mg/kg | RHS-DU-20<br>2/4/2014<br>mg/kg | RHS-DU-21<br>2/4/2014<br>mg/kg | RHS-DU-22<br>2/4/2014<br>mg/kg | HDOH Tier 1<br>EAL<br>mg/kg |
|---------------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|-----------------------------|
| <b>SVOCs (Continued)</b>              |                                |                                |                                |                                |                                |                                |                                |                             |
| 2-Chloronaphthalene                   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 2-Chlorophenol                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.0092                      |
| 2-Methylnaphthalene                   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 4.1                         |
| 2-Methylphenol (o-cresol)             | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 2-Nitroaniline                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 2-Nitrophenol                         | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 3,4-Methylphenol (m,p-cresol)         | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 4-Bromophenylphenylether              | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 4-Chloro-3-methylphenol               | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 4-Chloroaniline                       | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 0.0063                      |
| 4-Chlorophenylphenylether             | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 4-Nitroaniline                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 4-Nitrophenol                         | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| Acenaphthene                          | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 120                         |
| Acenaphthylene                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 100                         |
| Aniline                               | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Anthracene                            | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 4.3                         |
| Azobenzene                            | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Benzo(a)anthracene                    | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 1.5                         |
| Benzo(a)pyrene                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 0.15                        |
| Benzo(b)fluoranthene                  | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 1.5                         |
| Benzo(ghi)perylene                    | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 35                          |
| Benzo(k)fluoranthene                  | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 15                          |
| Benzyl alcohol                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Bis (2-chloroethoxy) methane          | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Bis (2-chloroethyl) ether             | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.000064                    |
| Bis (2-chloroisopropyl) ether         | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 0.0035                      |
| Bis (2-ethylhexyl) phthalate          | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 35                          |
| Bis(2-ethylhexyl) adipate             | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Butylbenzylphthalate                  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Carbazole                             | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Chrysene                              | 0.10                           | 0.16                           | 0.12                           | 0.11                           | 0.15                           | 0.12                           | 0.40                           | 30                          |
| Dibenzo(a,h)anthracene                | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 0.15                        |
| Dibenzofuran                          | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Diethylphthalate                      | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 16                          |
| Dimethylphthalate                     | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 22                          |
| Di-n-butylphthalate                   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Di-n-octyl phthalate                  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Fluoranthene                          | 0.10                           | 0.13                           | 0.11                           | ND< 0.1                        | 0.22                           | 0.15                           | 0.49                           | 460                         |
| Fluorene                              | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 100                         |

Notes on page 13



Table 3 (Continued)

| Sample ID:<br>Date Sampled:<br>Units:               | RHS-DU-16<br>2/4/2014<br>mg/kg | RHS-DU-17<br>2/4/2014<br>mg/kg | RHS-DU-18<br>2/4/2014<br>mg/kg | RHS-DU-19<br>2/4/2014<br>mg/kg | RHS-DU-20<br>2/4/2014<br>mg/kg | RHS-DU-21<br>2/4/2014<br>mg/kg | RHS-DU-22<br>2/4/2014<br>mg/kg | HDOH Tier 1<br>EAL<br>mg/kg |
|---|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|-----------------------------|
| <b>SVOCs (Continued)</b>                            |                                |                                |                                |                                |                                |                                |                                |                             |
| Hexachlorobenzene                                   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.3                         |
| Hexachlorobutadiene                                 | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.18                        |
| Hexachlorocyclopentadiene                           | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Hexachlorethane                                     | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.27                        |
| Indeno(1,2,3-cd)pyrene                              | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 1.5                         |
| Isophorone  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.77                        |
| Naphthalene   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 4.4                         |
| Nitrobenzene  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.0046                      |
| N-Nitroso-di-n-propylamine                          | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| N-nitrosodiphenylamine                              | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Pentachlorophenol                                   | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 0.82                        |
| Phenanthrene  | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 0.20                           | 440                         |
| Phenol  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.16                        |
| Pyrene  | 0.10                           | 0.17                           | 0.12                           | 0.12                           | 0.20                           | 0.16                           | 0.41                           | 44                          |
| Pyridine  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| <b>Organochlorine Pesticides / EPA Method 8081A</b> |                                |                                |                                |                                |                                |                                |                                |                             |
| Aldrin  | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | 0.92                        |
| Alpha-BHC   | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | NS                          |
| Alpha-Chlordane                                     | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | NS                          |
| Beta-BHC  | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | NS                          |
| Chlordane (technical)                               | 0.050                          | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | 16                          |
| Delta-BHC   | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | NS                          |
| Dieldrin  | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | 1.5                         |
| Endosulfan I  | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | 18                          |
| Endosulfan II                                       | ND< 0.010                      | ND< 0.010                      | 0.054                          | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | NS                          |
| Endosulfan sulfate                                  | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | NS                          |
| Endrin  | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | 3.7                         |
| Endrin aldehyde                                     | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | NS                          |
| Endrin ketone                                       | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | NS                          |
| Gamma-BHC (Lindane)                                 | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | 0.075                       |
| Gamma-Chlordane                                     | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | NS                          |
| Heptachlor  | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | 0.11                        |
| Heptachlor epoxide                                  | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | 0.053                       |
| Methoxychlor  | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | 16                          |
| p,p'-DDD  | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | 0.013                          | ND< 0.010                      | ND< 0.010                      | 2                           |
| p,p'-DDE  | ND< 0.010                      | ND< 0.010                      | 0.030                          | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | 1.4                         |
| p,p'-DDT  | ND< 0.010                      | 0.164                          | 0.132                          | 0.025                          | 0.019                          | 0.015                          | 0.022                          | 1.7                         |
| Toxaphene   | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | 0.44                        |

Notes on page 13

Table 3 (Continued)

| Sample ID:<br>Date Sampled:<br>Units:                             | RHS-DU-23<br>2/4/2014<br>mg/kg | RHS-DU-24<br>2/4/2014<br>mg/kg | RHS-DU-25<br>2/4/2014<br>mg/kg | RHS-DU-26<br>2/4/2014<br>mg/kg | RHS-DU-27<br>2/4/2014<br>mg/kg | RHS-DU-28<br>2/4/2014<br>mg/kg | RHS-DU-29<br>2/4/2014<br>mg/kg | HDOH Tier 1<br>EAL<br>mg/kg |
|---|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|-----------------------------|
| <b>Dioxin / EPA Method 8290</b>                                   |                                |                                |                                |                                |                                |                                |                                |                             |
| Total Dioxin Toxic Equivalent (TEQ) □                             | 18                             | 17                             | 20                             | 33                             | 22                             | 14                             | 21                             | 240 □                       |
| <b>Total Petroleum Hydrocarbons (TPH) / EPA Method 8015B</b>      |                                |                                |                                |                                |                                |                                |                                |                             |
| TPH-Diesel Range Organics   | ND< 50                         | ND< 50                         | ND< 50                         | ND< 50                         | ND< 50                         | ND< 50                         | ND< 50                         | 500                         |
| TPH-Residual Range Organics                                       | ND< 100                        | ND< 100                        | ND< 100                        | ND< 100                        | ND< 100                        | ND< 100                        | ND< 100                        | 500                         |
| <b>Total Metals / EPA Method 6010B/7471</b>                       |                                |                                |                                |                                |                                |                                |                                |                             |
| Arsenic   | 18                             | 5.1                            | 5.8                            | 5.3                            | 6.3                            | 4.6                            | 5.7                            | 24                          |
| Barium  | 340                            | 270                            | 280                            | 100                            | 85                             | 91                             | 99                             | 1,000                       |
| Cadmium   | 1.9                            | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 14                          |
| Chromium  | 150                            | 140                            | 190                            | 250                            | 230                            | 250                            | 270                            | 1,100                       |
| Lead  | 500*                           | 110                            | 120                            | 89                             | 59                             | 44                             | 68                             | 200                         |
| Selenium  | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | 78                          |
| Silver  | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | ND< 20                         | 78                          |
| Mercury   | 200*                           | 4.8*                           | 4.4                            | 3.5                            | 2.9                            | 3.1                            | 5.1*                           | 4.7                         |
| <b>Polychlorinated Biphenyls (PCBs) / EPA Method 8082</b>         |                                |                                |                                |                                |                                |                                |                                |                             |
| PCB-1016  | ND< 0.10                       | ND< 0.10                       | ND< 0.10                       | ND< 0.10                       | ND< 0.10                       | ND< 0.10                       | ND< 0.10                       | 1.1                         |
| PCB-1221  | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | 1.1                         |
| PCB-1232  | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | ND< 0.20                       | 1.1                         |
| PCB-1242  | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | 1.1                         |
| PCB-1248  | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | 1.1                         |
| PCB-1254  | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | 1.1                         |
| PCB-1260  | 1.46*                          | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | ND< 0.05                       | 1.1                         |
| <b>Semi-volatile Organic Compounds (SVOCs) / EPA Method 8270C</b> |                                |                                |                                |                                |                                |                                |                                |                             |
| 1,2,4-Trichlorobenzene  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.098                       |
| 1,2-Dichlorobenzene   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.75                        |
| 1,2-Dinitrobenzene  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 1,3-Dichlorobenzene   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.57                        |
| 1,3-Dinitrobenzene  | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 1,4-Dichlorobenzene   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.047                       |
| 1,4-Dinitrobenzene  | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 1-Methylnapthalene  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 1.8                         |
| 2,3,4,6-Tetrachlorophenol   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 4.9                         |
| 2,3,5,6-Tetrachlorophenol   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 2,4,5-Trichlorophenol   | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 29                          |
| 2,4,6-Trichlorophenol   | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 1.8                         |
| 2,4-Dichlorophenol  | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 0.025                       |
| 2,4-Dimethylphenol  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 9.9                         |
| 2,4-Dinitrophenol   | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 5.6                         |
| 2,4-Dinitrotoluene  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.021                       |
| 2,6-Dinitrotoluene  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 3.6                         |

Notes on page 13



Table 3 (Continued)

| Sample ID:<br>Date Sampled:<br>Units: | RHS-DU-23<br>2/4/2014<br>mg/kg | RHS-DU-24<br>2/4/2014<br>mg/kg | RHS-DU-25<br>2/4/2014<br>mg/kg | RHS-DU-26<br>2/4/2014<br>mg/kg | RHS-DU-27<br>2/4/2014<br>mg/kg | RHS-DU-28<br>2/4/2014<br>mg/kg | RHS-DU-29<br>2/4/2014<br>mg/kg | HDOH Tier 1<br>EAL<br>mg/kg |
|---------------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|-----------------------------|
| <b>SVOCs (Continued)</b>              |                                |                                |                                |                                |                                |                                |                                |                             |
| 2-Chloronaphthalene                   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 2-Chlorophenol                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.0092                      |
| 2-Methylnaphthalene                   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 4.1                         |
| 2-Methylphenol (o-cresol)             | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 2-Nitroaniline                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 2-Nitrophenol                         | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 3,4-Methylphenol (m,p-cresol)         | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 4-Bromophenylphenylether              | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 4-Chloro-3-methylphenol               | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 4-Chloroaniline                       | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 0.0063                      |
| 4-Chlorophenylphenylether             | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| 4-Nitroaniline                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| 4-Nitrophenol                         | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | NS                          |
| Acenaphthene                          | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 120                         |
| Acenaphthylene                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 100                         |
| Aniline                               | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Anthracene                            | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 4.3                         |
| Azobenzene                            | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Benzo(a)anthracene                    | 0.16                           | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 1.5                         |
| Benzo(a)pyrene                        | 0.15*                          | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 0.15                        |
| Benzo(b)fluoranthene                  | 0.31                           | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 1.5                         |
| Benzo(ghi)perylene                    | 0.44                           | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 35                          |
| Benzo(k)fluoranthene                  | 0.33                           | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 15                          |
| Benzyl alcohol                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Bis (2-chloroethoxy) methane          | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Bis (2-chloroethyl) ether             | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.000064                    |
| Bis (2-chloroisopropyl) ether         | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 0.0035                      |
| Bis (2-ethylhexyl) phthalate          | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 35                          |
| Bis(2-ethylhexyl) adipate             | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Butylbenzylphthalate                  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Carbazole                             | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Chrysene                              | 0.48                           | 0.21                           | 0.11                           | ND< 0.1                        | ND< 0.1                        | 0.10                           | 0.14                           | 30                          |
| Dibenzo(a,h)anthracene                | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 0.15                        |
| Dibenzofuran                          | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Diethylphthalate                      | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 16                          |
| Dimethylphthalate                     | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 22                          |
| Di-n-butylphthalate                   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Di-n-octyl phthalate                  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Fluoranthene                          | 0.49                           | 0.13                           | 0.11                           | ND< 0.1                        | ND< 0.1                        | 0.11                           | 0.10                           | 460                         |
| Fluorene                              | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 100                         |

Notes on page 13



Table 3 (Continued)

| Sample ID:<br>Date Sampled:<br>Units:               | RHS-DU-23<br>2/4/2014<br>mg/kg | RHS-DU-24<br>2/4/2014<br>mg/kg | RHS-DU-25<br>2/4/2014<br>mg/kg | RHS-DU-26<br>2/4/2014<br>mg/kg | RHS-DU-27<br>2/4/2014<br>mg/kg | RHS-DU-28<br>2/4/2014<br>mg/kg | RHS-DU-29<br>2/4/2014<br>mg/kg | HDOH Tier 1<br>EAL<br>mg/kg |
|---|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|--------------------------------|-----------------------------|
| <b>SVOCs (Continued)</b>                            |                                |                                |                                |                                |                                |                                |                                |                             |
| Hexachlorobenzene                                   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.3                         |
| Hexachlorobutadiene                                 | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.18                        |
| Hexachlorocyclopentadiene                           | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Hexachloroethane                                    | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.27                        |
| Indeno(1,2,3-cd)pyrene                              | 0.38                           | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 1.5                         |
| Isophorone  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.77                        |
| Naphthalene   | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 4.4                         |
| Nitrobenzene  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.0046                      |
| N-Nitroso-di-n-propylamine                          | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| N-nitrosodiphenylamine                              | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| Pentachlorophenol                                   | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | ND< 5.0                        | 0.82                        |
| Phenanthrene  | 0.16                           | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | ND< 0.1                        | 440                         |
| Phenol  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | 0.16                        |
| Pyrene  | 0.50                           | 0.14                           | 0.11                           | ND< 0.1                        | ND< 0.1                        | 0.10                           | 0.10                           | 44                          |
| Pyridine  | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | ND< 1.0                        | NS                          |
| <b>Organochlorine Pesticides / EPA Method 8081A</b> |                                |                                |                                |                                |                                |                                |                                |                             |
| Aldrin  | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | 0.92                        |
| Alpha-BHC   | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | NS                          |
| Alpha-Chlordane                                     | 0.011                          | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | NS                          |
| Beta-BHC  | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | NS                          |
| Chlordane (technical)                               | 0.057                          | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | 16                          |
| Delta-BHC   | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | NS                          |
| Dieldrin  | 0.017                          | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | 1.5                         |
| Endosulfan I  | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | 18                          |
| Endosulfan II                                       | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | NS                          |
| Endosulfan sulfate                                  | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | NS                          |
| Endrin  | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | 3.7                         |
| Endrin aldehyde                                     | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | NS                          |
| Endrin ketone                                       | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | NS                          |
| Gamma-BHC (Lindane)                                 | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | 0.075                       |
| Gamma-Chlordane                                     | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | NS                          |
| Heptachlor  | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | 0.11                        |
| Heptachlor epoxide                                  | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | ND< 0.005                      | 0.053                       |
| Methoxychlor  | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | 16                          |
| p,p'-DDD  | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | 2                           |
| p,p'-DDE  | 0.011                          | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | 1.4                         |
| p,p'-DDT  | 0.074                          | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | ND< 0.010                      | 1.7                         |
| Toxaphene   | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | ND< 0.050                      | 0.44                        |

Notes on page 13

Table 3 (Continued)

Notes:

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HDOH Tier 1 EAL

*Italics*

mg/kg

ND<

NS

Dioxin Toxic Equivalent (TEQ) reported in nanograms per kilogram (not milligrams per kilogram).

The value exceeds the State of Hawaii Department of Health (HDOH) Tier 1 Environmental Action Levels (EAL) .

The State of Hawaii Department of Health (HDOH) Tier 1 Environmental Action Level (EAL) for sites where groundwater is a current or potential source of drinking water, and the site is greater than 150 meters from a surface water body. (Fall 2011, updated January 2012).

Reporting Limit is above the HDOH Tier 1 EAL.

milligrams per kilogram.

No detectable concentration. The number following the "less than" symbol is the laboratory reporting limit .

No standard. The HDOH has not established a Tier 1 EAL for this analyte.

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**Table 4**  
**Analytical Results for Replicate Samples**  
**State of Hawaii Department of Education**  
**Radford High School Track and Field Investigation**  
**Honolulu, Oahu, Hawaii**

Project No: 17012-012148.00 / Task 048

| Sample ID:  | RHS-DU-1 | RHS-DU-1.2 Δ | RHS-DU-1.3 Δ | RHS-DU-17 | RHS-DU-17.2 Δ | RHS-DU-17.3 Δ | RHS-DU-26 | RHS-DU-26.2 Δ | RHS-DU-26.3 Δ | HDOH Tier 1 EAL |
|---|----------|--------------|--------------|-----------|---------------|---------------|-----------|---------------|---------------|-----------------|
| Date Sampled:   | 2/4/2014 | 2/5/2014     | 2/5/2014     | 2/4/2014  | 2/4/2014      | 2/4/2014      | 2/4/2014  | 2/4/2014      | 2/4/2014      | 2/4/2014        |
| Units:  | mg/kg    | mg/kg        | mg/kg        | mg/kg     | mg/kg         | mg/kg         | mg/kg     | mg/kg         | mg/kg         | mg/kg           |
| <b>Dioxin / EPA Method 8290</b>                                   |          |              |              |           |               |               |           |               |               |                 |
| Total Dioxin Toxic Equivalent (TEQ) □                             | 710*     | 890*         | 2,900*       | 230       | 290*          | 280*          | 33        | 19            | 20            | 240 □           |
| <b>Total Petroleum Hydrocarbons (TPH) / EPA Method 8015B</b>      |          |              |              |           |               |               |           |               |               |                 |
| TPH-Diesel Range Organics   | ND< 50   | ND< 50       | ND< 50       | ND< 50    | ND< 50        | ND< 50        | ND< 50    | ND< 50        | ND< 50        | 500             |
| TPH-Residual Range Organics                                       | ND< 100  | ND< 100      | ND< 100      | ND< 100   | ND< 100       | ND< 100       | ND< 100   | ND< 100       | ND< 100       | 500             |
| <b>Total Metals / EPA Method 6010B/7471</b>                       |          |              |              |           |               |               |           |               |               |                 |
| Arsenic   | 35*      | 38*          | 28*          | 23        | 24*           | 26*           | 5.3       | ND< 5.0       | 6.2           | 24              |
| Barium  | 1,200*   | 1,100*       | 1,100*       | 590       | 720           | 670           | 100       | 69            | 85            | 1,000           |
| Cadmium   | 14*      | 14*          | 26*          | 4.5       | 4.6           | 6.4           | ND< 1.0   | ND< 1.0       | ND< 1.0       | 14              |
| Chromium  | 300      | 230          | 310          | 210       | 220           | 150           | 250       | 220           | 190           | 1,100           |
| Lead  | 4,800*   | 5,800*       | 15,000*      | 1,100*    | 1,600*        | 1,600*        | 89        | 50            | 78            | 200             |
| Selenium  | ND< 20   | ND< 20       | ND< 20       | ND< 20    | ND< 20        | ND< 20        | ND< 20    | ND< 20        | ND< 20        | 78              |
| Silver  | ND< 20   | ND< 20       | ND< 20       | ND< 20    | ND< 20        | ND< 20        | ND< 20    | ND< 20        | ND< 20        | 78              |
| Mercury   | 5.2*     | 11*          | 18*          | 73*       | 66*           | 83*           | 3.5       | 2.7           | 2.9           | 4.7             |
| <b>Polychlorinated Biphenyls (PCBs) / EPA Method 8082</b>         |          |              |              |           |               |               |           |               |               |                 |
| PCB-1016  | ND< 0.10 | ND< 0.10     | ND< 0.10     | ND< 0.10  | ND< 0.10      | ND< 0.10      | ND< 0.10  | ND< 0.10      | ND< 0.10      | 1.1             |
| PCB-1221  | ND< 0.20 | ND< 0.20     | ND< 0.20     | ND< 0.20  | ND< 0.20      | ND< 0.20      | ND< 0.20  | ND< 0.20      | ND< 0.20      | 1.1             |
| PCB-1232  | ND< 0.20 | ND< 0.20     | ND< 0.20     | ND< 0.20  | ND< 0.20      | ND< 0.20      | ND< 0.20  | ND< 0.20      | ND< 0.20      | 1.1             |
| PCB-1242  | ND< 0.05 | ND< 0.05     | ND< 0.05     | ND< 0.05  | ND< 0.05      | ND< 0.05      | ND< 0.05  | ND< 0.05      | ND< 0.05      | 1.1             |
| PCB-1248  | ND< 0.05 | ND< 0.05     | ND< 0.05     | ND< 0.05  | ND< 0.05      | ND< 0.05      | ND< 0.05  | ND< 0.05      | ND< 0.05      | 1.1             |
| PCB-1254  | ND< 0.05 | ND< 0.05     | ND< 0.05     | ND< 0.05  | ND< 0.05      | ND< 0.05      | ND< 0.05  | ND< 0.05      | ND< 0.05      | 1.1             |
| PCB-1260  | ND< 0.05 | ND< 0.05     | ND< 0.05     | 2.23*     | 2.35*         | 2.24*         | ND< 0.05  | ND< 0.05      | ND< 0.05      | 1.1             |
| <b>Semi-volatile Organic Compounds (SVOCs) / EPA Method 8270C</b> |          |              |              |           |               |               |           |               |               |                 |
| 1,2,4-Trichlorobenzene  | ND< 1.0  | ND< 1.0      | ND< 1.0      | ND< 1.0   | ND< 1.0       | ND< 1.0       | ND< 1.0   | ND< 1.0       | ND< 1.0       | 0.098           |
| 1,2-Dichlorobenzene   | ND< 1.0  | ND< 1.0      | ND< 1.0      | ND< 1.0   | ND< 1.0       | ND< 1.0       | ND< 1.0   | ND< 1.0       | ND< 1.0       | 0.75            |
| 1,2-Dinitrobenzene  | ND< 1.0  | ND< 1.0      | ND< 1.0      | ND< 1.0   | ND< 1.0       | ND< 1.0       | ND< 1.0   | ND< 1.0       | ND< 1.0       | NS              |
| 1,3-Dichlorobenzene   | ND< 1.0  | ND< 1.0      | ND< 1.0      | ND< 1.0   | ND< 1.0       | ND< 1.0       | ND< 1.0   | ND< 1.0       | ND< 1.0       | 0.57            |
| 1,3-Dinitrobenzene  | ND< 5.0  | ND< 5.0      | ND< 5.0      | ND< 5.0   | ND< 5.0       | ND< 5.0       | ND< 5.0   | ND< 5.0       | ND< 5.0       | NS              |
| 1,4-Dichlorobenzene   | ND< 1.0  | ND< 1.0      | ND< 1.0      | ND< 1.0   | ND< 1.0       | ND< 1.0       | ND< 1.0   | ND< 1.0       | ND< 1.0       | 0.047           |
| 1,4-Dinitrobenzene  | ND< 5.0  | ND< 5.0      | ND< 5.0      | ND< 5.0   | ND< 5.0       | ND< 5.0       | ND< 5.0   | ND< 5.0       | ND< 5.0       | NS              |
| 1-Methylnapthalene  | ND< 1.0  | ND< 1.0      | ND< 1.0      | ND< 1.0   | ND< 1.0       | ND< 1.0       | ND< 1.0   | ND< 1.0       | ND< 1.0       | 1.8             |
| 2,3,4,6-Tetrachlorophenol   | ND< 1.0  | ND< 1.0      | ND< 1.0      | ND< 1.0   | ND< 1.0       | ND< 1.0       | ND< 1.0   | ND< 1.0       | ND< 1.0       | 4.9             |
| 2,3,5,6-Tetrachlorophenol   | ND< 1.0  | ND< 1.0      | ND< 1.0      | ND< 1.0   | ND< 1.0       | ND< 1.0       | ND< 1.0   | ND< 1.0       | ND< 1.0       | NS              |
| 2,4,5-Trichlorophenol   | ND< 5.0  | ND< 5.0      | ND< 5.0      | ND< 5.0   | ND< 5.0       | ND< 5.0       | ND< 5.0   | ND< 5.0       | ND< 5.0       | 29              |
| 2,4,6-Trichlorophenol   | ND< 5.0  | ND< 5.0      | ND< 5.0      | ND< 5.0   | ND< 5.0       | ND< 5.0       | ND< 5.0   | ND< 5.0       | ND< 5.0       | 1.8             |

Notes on page 4





Table 4 (Continued)

| Sample ID:<br>Date Sampled:<br>Units: | RHS-DU-1<br>2/4/2014<br>mg/kg | RHS-DU-1.2 Δ<br>2/5/2014<br>mg/kg | RHS-DU-1.3 Δ<br>2/5/2014<br>mg/kg | RHS-DU-17<br>2/4/2014<br>mg/kg | RHS-DU-17.2 Δ<br>2/4/2014<br>mg/kg | RHS-DU-17.3 Δ<br>2/4/2014<br>mg/kg | RHS-DU-26<br>2/4/2014<br>mg/kg | RHS-DU-26.2 Δ<br>2/4/2014<br>mg/kg | RHS-DU-26.3 Δ<br>2/4/2014<br>mg/kg | HDOH Tier 1<br>EAL<br>mg/kg |
|---------------------------------------|-------------------------------|-----------------------------------|-----------------------------------|--------------------------------|------------------------------------|------------------------------------|--------------------------------|------------------------------------|------------------------------------|-----------------------------|
| <b>SVOCs (Continued)</b>              |                               |                                   |                                   |                                |                                    |                                    |                                |                                    |                                    |                             |
| 2,4-Dichlorophenol                    | ND< 5.0                       | ND< 5.0                           | ND< 5.0                           | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | 0.025                       |
| 2,4-Dimethylphenol                    | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | 9.9                         |
| 2,4-Dinitrophenol                     | ND< 5.0                       | ND< 5.0                           | ND< 5.0                           | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | 5.6                         |
| 2,4-Dinitrotoluene                    | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | 0.021                       |
| 2,6-Dinitrotoluene                    | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | 3.6                         |
| 2-Chloronaphthalene                   | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| 2-Chlorophenol                        | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | 0.0092                      |
| 2-Methylnaphthalene                   | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | 4.1                         |
| 2-Methylphenol (o-cresol)             | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| 2-Nitroaniline                        | ND< 5.0                       | ND< 5.0                           | ND< 5.0                           | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | NS                          |
| 2-Nitrophenol                         | ND< 5.0                       | ND< 5.0                           | ND< 5.0                           | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | NS                          |
| 3,4-Methylphenol (m,p-cresol)         | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| 4-Bromophenylphenylether              | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| 4-Chloro-3-methylphenol               | ND< 5.0                       | ND< 5.0                           | ND< 5.0                           | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | NS                          |
| 4-Chloroaniline                       | ND< 5.0                       | ND< 5.0                           | ND< 5.0                           | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | 0.0063                      |
| 4-Chlorophenylphenylether             | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| 4-Nitroaniline                        | ND< 5.0                       | ND< 5.0                           | ND< 5.0                           | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | NS                          |
| 4-Nitrophenol                         | ND< 5.0                       | ND< 5.0                           | ND< 5.0                           | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | NS                          |
| Acenaphthene                          | ND< 0.1                       | ND< 0.1                           | ND< 0.1                           | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | 120                         |
| Acenaphthylene                        | ND< 0.1                       | ND< 0.1                           | ND< 0.1                           | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | 100                         |
| Aniline                               | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| Anthracene                            | ND< 0.1                       | ND< 0.1                           | ND< 0.1                           | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | 4.3                         |
| Azobenzene                            | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| Benzo(a)anthracene                    | ND< 0.1                       | ND< 0.1                           | ND< 0.1                           | ND< 0.1                        | ND< 0.1                            | 0.18                               | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | 1.5                         |
| Benzo(a)pyrene                        | ND< 0.1                       | ND< 0.1                           | ND< 0.1                           | ND< 0.1                        | ND< 0.1                            | 0.43*                              | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | 0.15                        |
| Benzo(b)fluoranthene                  | ND< 0.1                       | ND< 0.1                           | ND< 0.1                           | ND< 0.1                        | ND< 0.1                            | 0.40                               | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | 1.5                         |
| Benzo(ghi)perylene                    | ND< 0.1                       | ND< 0.1                           | ND< 0.1                           | ND< 0.1                        | ND< 0.1                            | 0.38                               | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | 35                          |
| Benzo(k)fluoranthene                  | ND< 0.1                       | ND< 0.1                           | ND< 0.1                           | ND< 0.1                        | ND< 0.1                            | 0.35                               | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | 15                          |
| Benzyl alcohol                        | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| Bis (2-chloroethoxy) methane          | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| Bis (2-chloroethyl) ether             | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | 0.000064                    |
| Bis (2-chloroisopropyl) ether         | ND< 5.0                       | ND< 5.0                           | ND< 5.0                           | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | 0.0035                      |
| Bis (2-ethylhexyl) phthalate          | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | 35                          |
| Bis(2-ethylhexyl) adipate             | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| Butylbenzylphthalate                  | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| Carbazole                             | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| Chrysene                              | ND< 0.1                       | ND< 0.1                           | ND< 0.1                           | 0.16                           | 0.15                               | 0.46                               | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | 30                          |
| Dibenzo(a,h)anthracene                | ND< 0.1                       | ND< 0.1                           | ND< 0.1                           | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | 0.15                        |
| Dibenzofuran                          | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| Diethylphthalate                      | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | 16                          |
| Dimethylphthalate                     | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | 22                          |

Notes on page 4

Table 4 (Continued)

| Sample ID:<br>Date Sampled:<br>Units:               | RHS-DU-1<br>2/4/2014<br>mg/kg | RHS-DU-1.2 Δ<br>2/5/2014<br>mg/kg | RHS-DU-1.3 Δ<br>2/5/2014<br>mg/kg | RHS-DU-17<br>2/4/2014<br>mg/kg | RHS-DU-17.2 Δ<br>2/4/2014<br>mg/kg | RHS-DU-17.3 Δ<br>2/4/2014<br>mg/kg | RHS-DU-26<br>2/4/2014<br>mg/kg | RHS-DU-26.2 Δ<br>2/4/2014<br>mg/kg | RHS-DU-26.3 Δ<br>2/4/2014<br>mg/kg | HDOH Tier 1<br>EAL<br>mg/kg |
|---|-------------------------------|-----------------------------------|-----------------------------------|--------------------------------|------------------------------------|------------------------------------|--------------------------------|------------------------------------|------------------------------------|-----------------------------|
| <b>SVOCs (Continued)</b>                            |                               |                                   |                                   |                                |                                    |                                    |                                |                                    |                                    |                             |
| Di-n-butylphthalate                                 | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| Di-n-octyl phthalate                                | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| Fluoranthene  | ND< 0.1                       | ND< 0.1                           | ND< 0.1                           | 0.13                           | 0.11                               | 0.31                               | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | 460                         |
| Fluorene  | ND< 0.1                       | ND< 0.1                           | ND< 0.1                           | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | 100                         |
| Hexachlorobenzene                                   | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | 0.3                         |
| Hexachlorobutadiene                                 | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | 0.18                        |
| Hexachlorocyclopentadiene                           | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| Hexachlorethane                                     | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | 0.27                        |
| Indeno(1,2,3-cd)pyrene                              | ND< 0.1                       | ND< 0.1                           | ND< 0.1                           | ND< 0.1                        | ND< 0.1                            | 0.40                               | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | 1.5                         |
| Isophorone  | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | 0.77                        |
| Naphthalene   | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | 4.4                         |
| Nitrobenzene  | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | 0.0046                      |
| N-Nitroso-di-n-propylamine                          | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| N-nitrosodiphenylamine                              | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| Pentachlorophenol                                   | ND< 5.0                       | ND< 5.0                           | ND< 5.0                           | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | ND< 5.0                        | ND< 5.0                            | ND< 5.0                            | 0.82                        |
| Phenanthrene  | ND< 0.1                       | ND< 0.1                           | ND< 0.1                           | ND< 0.1                        | ND< 0.1                            | 0.12                               | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | 440                         |
| Phenol  | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | 0.16                        |
| Pyrene  | ND< 0.1                       | ND< 0.1                           | ND< 0.1                           | 0.17                           | 0.14                               | 0.33                               | ND< 0.1                        | ND< 0.1                            | ND< 0.1                            | 44                          |
| Pyridine  | ND< 1.0                       | ND< 1.0                           | ND< 1.0                           | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | ND< 1.0                        | ND< 1.0                            | ND< 1.0                            | NS                          |
| <b>Organochlorine Pesticides / EPA Method 8081A</b> |                               |                                   |                                   |                                |                                    |                                    |                                |                                    |                                    |                             |
| Aldrin  | ND< 0.005                     | ND< 0.005                         | ND< 0.005                         | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | 0.92                        |
| Alpha-BHC   | ND< 0.005                     | ND< 0.005                         | ND< 0.005                         | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | NS                          |
| Alpha-Chlordane                                     | ND< 0.005                     | ND< 0.005                         | ND< 0.005                         | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | NS                          |
| Beta-BHC  | ND< 0.005                     | ND< 0.005                         | ND< 0.005                         | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | NS                          |
| Chlordane (technical)                               | ND< 0.050                     | ND< 0.050                         | ND< 0.050                         | ND< 0.050                      | ND< 0.050                          | 0.225                              | ND< 0.050                      | ND< 0.050                          | ND< 0.050                          | 16                          |
| Delta-BHC   | ND< 0.005                     | ND< 0.005                         | ND< 0.005                         | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | NS                          |
| Dieldrin  | ND< 0.010                     | ND< 0.010                         | ND< 0.010                         | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | 1.5                         |
| Endosulfan I  | ND< 0.005                     | ND< 0.005                         | ND< 0.005                         | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | 18                          |
| Endosulfan II                                       | ND< 0.010                     | ND< 0.010                         | ND< 0.010                         | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | NS                          |
| Endosulfan sulfate                                  | ND< 0.010                     | ND< 0.010                         | ND< 0.010                         | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | NS                          |
| Endrin  | ND< 0.010                     | ND< 0.010                         | ND< 0.010                         | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | 3.7                         |
| Endrin aldehyde                                     | ND< 0.010                     | ND< 0.010                         | ND< 0.010                         | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | NS                          |
| Endrin ketone                                       | ND< 0.010                     | ND< 0.010                         | ND< 0.010                         | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | NS                          |
| Gamma-BHC (Lindane)                                 | ND< 0.005                     | ND< 0.005                         | ND< 0.005                         | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | 0.075                       |
| Gamma-Chlordane                                     | ND< 0.005                     | ND< 0.005                         | ND< 0.005                         | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | NS                          |
| Heptachlor  | ND< 0.005                     | ND< 0.005                         | ND< 0.005                         | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | 0.11                        |
| Heptachlor epoxide                                  | ND< 0.005                     | ND< 0.005                         | ND< 0.005                         | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | ND< 0.005                      | ND< 0.005                          | ND< 0.005                          | 0.053                       |
| Methoxychlor  | ND< 0.010                     | ND< 0.010                         | ND< 0.010                         | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | 16                          |
| p,p'-DDD  | ND< 0.010                     | ND< 0.010                         | ND< 0.010                         | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | 2                           |
| p,p'-DDE  | ND< 0.010                     | ND< 0.010                         | ND< 0.010                         | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | 1.4                         |
| p,p'-DDT  | ND< 0.010                     | ND< 0.010                         | ND< 0.010                         | 0.164                          | 0.195                              | 0.175                              | ND< 0.010                      | ND< 0.010                          | ND< 0.010                          | 1.7                         |
| Toxaphene   | ND< 0.050                     | ND< 0.050                         | ND< 0.050                         | ND< 0.050                      | ND< 0.050                          | ND< 0.050                          | ND< 0.050                      | ND< 0.050                          | ND< 0.050                          | 0.44                        |

Notes on page 4

Table 4 (Continued)

Notes:

- Δ Indicates a duplicate sample.
- Dioxin Toxic Equivalent (TEQ) reported in nanograms per kilogram (not milligrams per kilogram).
- Bold\*** The value exceeds the State of Hawaii Department of Health (HDOH) Tier 1 Environmental Action Levels (EAL) .
- HDOH Tier 1 EAL The State of Hawaii Department of Health (HDOH) Tier 1 Environmental Action Level (EAL) for sites where groundwater is a current or potential source of drinking water, and the site is greater than 150 meters from a surface water body. (Fall 2011, updated January 2012).
- Italics* Reporting Limit is above the HDOH Tier 1 EAL.
- mg/kg milligrams per kilogram.
- ND< No detectable concentration. The number following the "less than" symbol is the laboratory reporting limit .
- NS No standard. The HDOH has not established a Tier 1 EAL for this analyte.

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**Table 5**  
**Statistical Calculations for Replicate Samples**  
**State of Hawaii Department of Education**  
**Radford High School Track and Field Investigation**  
**Honolulu, Oahu, Hawaii**

**Project No: 17012-012148.00 / Task 048**

| Analyte                                    | EPA Method | Sample Identification | Sample Type | Result | Units | RPD of Primary and Duplicate | RPD of Primary and Triplicate | Mean    | Standard Deviation | RSD   |
|--|------------|-----------------------|-------------|--------|-------|------------------------------|-------------------------------|---------|--------------------|-------|
| <b>Decision Unit 1 - Replicate Samples</b> |            |                       |             |        |       |                              |                               |         |                    |       |
| Dioxin                                     | 8290       | RHS-DU1               | Primary     | 710    | ng/kg | 22.5%                        | 121.3%                        | 1,500.0 | 1,215.8            | 81.1% |
|  |            | RHS-DU1.2             | Duplicate   | 890    | ng/kg |                              |                               |         |                    |       |
|  |            | RHS-DU1.3             | Triplicate  | 2,900  | ng/kg |                              |                               |         |                    |       |
| Arsenic                                    | 6010B      | RHS-DU1               | Primary     | 35     | mg/kg | 8.2%                         | 22.2%                         | 33.7    | 5.1                | 15.2% |
|  |            | RHS-DU1.2             | Duplicate   | 38     | mg/kg |                              |                               |         |                    |       |
|  |            | RHS-DU1.3             | Triplicate  | 28     | mg/kg |                              |                               |         |                    |       |
| Barium                                     | 6010B      | RHS-DU1               | Primary     | 1,200  | mg/kg | 8.7%                         | 8.7%                          | 1,133.3 | 57.7               | 5.1%  |
|  |            | RHS-DU1.2             | Duplicate   | 1,100  | mg/kg |                              |                               |         |                    |       |
|  |            | RHS-DU1.3             | Triplicate  | 1,100  | mg/kg |                              |                               |         |                    |       |
| Cadmium                                    | 6010B      | RHS-DU1               | Primary     | 14     | mg/kg | 0.0%                         | 60.0%                         | 18.0    | 6.9                | 38.5% |
|  |            | RHS-DU1.2             | Duplicate   | 14     | mg/kg |                              |                               |         |                    |       |
|  |            | RHS-DU1.3             | Triplicate  | 26     | mg/kg |                              |                               |         |                    |       |
| Chromium                                   | 6010B      | RHS-DU1               | Primary     | 300    | mg/kg | 26.4%                        | 3.3%                          | 280.0   | 43.6               | 15.6% |
|  |            | RHS-DU1.2             | Duplicate   | 230    | mg/kg |                              |                               |         |                    |       |
|  |            | RHS-DU1.3             | Triplicate  | 310    | mg/kg |                              |                               |         |                    |       |
| Lead                                       | 6010B      | RHS-DU1               | Primary     | 4,800  | mg/kg | 18.9%                        | 103.0%                        | 8,533.3 | 5,622.6            | 65.9% |
|  |            | RHS-DU1.2             | Duplicate   | 5,800  | mg/kg |                              |                               |         |                    |       |
|  |            | RHS-DU1.3             | Triplicate  | 15,000 | mg/kg |                              |                               |         |                    |       |
| Mercury                                    | 7471       | RHS-DU1               | Primary     | 5      | mg/kg | 71.6%                        | 110.3%                        | 11.4    | 6.4                | 56.2% |
|  |            | RHS-DU1.2             | Duplicate   | 11     | mg/kg |                              |                               |         |                    |       |
|  |            | RHS-DU1.3             | Triplicate  | 18     | mg/kg |                              |                               |         |                    |       |

Notes on page 4



Table 5 (Continued)

| Analyte                                     | EPA Method | Sample Identification | Sample Type | Result | Units | RPD of Primary and Duplicate | RPD of Primary and Triplicate | Mean    | Standard Deviation | RSD   |
|---|------------|-----------------------|-------------|--------|-------|------------------------------|-------------------------------|---------|--------------------|-------|
| <b>Decision Unit 17 - Replicate Samples</b> |            |                       |             |        |       |                              |                               |         |                    |       |
| Dioxin                                      | 8290       | RHS-DU17              | Primary     | 230    | ng/kg | 23.1%                        | 19.6%                         | 266.7   | 32.1               | 12.1% |
|   |            | RHS-DU17.2            | Duplicate   | 290    | ng/kg |                              |                               |         |                    |       |
|   |            | RHS-DU17.3            | Triplicate  | 280    | ng/kg |                              |                               |         |                    |       |
| Arsenic                                     | 6010B      | RHS-DU17              | Primary     | 23     | mg/kg | 4.3%                         | 12.2%                         | 24.3    | 1.5                | 6.3%  |
|   |            | RHS-DU17.2            | Duplicate   | 24     | mg/kg |                              |                               |         |                    |       |
|   |            | RHS-DU17.3            | Triplicate  | 26     | mg/kg |                              |                               |         |                    |       |
| Barium                                      | 6010B      | RHS-DU17              | Primary     | 590    | mg/kg | 19.8%                        | 12.7%                         | 660.0   | 65.6               | 9.9%  |
|   |            | RHS-DU17.2            | Duplicate   | 720    | mg/kg |                              |                               |         |                    |       |
|   |            | RHS-DU17.3            | Triplicate  | 670    | mg/kg |                              |                               |         |                    |       |
| Cadmium                                     | 6010B      | RHS-DU17              | Primary     | 4.5    | mg/kg | 2.2%                         | 34.9%                         | 5.2     | 1.1                | 20.7% |
|   |            | RHS-DU17.2            | Duplicate   | 4.6    | mg/kg |                              |                               |         |                    |       |
|   |            | RHS-DU17.3            | Triplicate  | 6.4    | mg/kg |                              |                               |         |                    |       |
| Chromium                                    | 6010B      | RHS-DU17              | Primary     | 210    | mg/kg | 4.7%                         | 33.3%                         | 193.3   | 37.9               | 19.6% |
|   |            | RHS-DU17.2            | Duplicate   | 220    | mg/kg |                              |                               |         |                    |       |
|   |            | RHS-DU17.3            | Triplicate  | 150    | mg/kg |                              |                               |         |                    |       |
| Lead  | 6010B      | RHS-DU17              | Primary     | 1,100  | mg/kg | 37.0%                        | 37.0%                         | 1,433.3 | 288.7              | 20.1% |
|   |            | RHS-DU17.2            | Duplicate   | 1,600  | mg/kg |                              |                               |         |                    |       |
|   |            | RHS-DU17.3            | Triplicate  | 1,600  | mg/kg |                              |                               |         |                    |       |
| Mercury                                     | 7471       | RHS-DU17              | Primary     | 73     | mg/kg | 10.1%                        | 12.8%                         | 74.0    | 8.5                | 11.5% |
|   |            | RHS-DU17.2            | Duplicate   | 66     | mg/kg |                              |                               |         |                    |       |
|   |            | RHS-DU17.3            | Triplicate  | 83     | mg/kg |                              |                               |         |                    |       |
| Chrysene                                    | 8270C      | RHS-DU17              | Primary     | 0.16   | mg/kg | 6.5%                         | 96.8%                         | 0.26    | 0.18               | 68.6% |
|   |            | RHS-DU17.2            | Duplicate   | 0.15   | mg/kg |                              |                               |         |                    |       |
|   |            | RHS-DU17.3            | Triplicate  | 0.46   | mg/kg |                              |                               |         |                    |       |
| Fluoroanthene                               | 8270C      | RHS-DU17              | Primary     | 0.13   | mg/kg | 16.7%                        | 81.8%                         | 0.18    | 0.11               | 60.1% |
|   |            | RHS-DU17.2            | Duplicate   | 0.11   | mg/kg |                              |                               |         |                    |       |
|   |            | RHS-DU17.3            | Triplicate  | 0.31   | mg/kg |                              |                               |         |                    |       |

Notes on page 4



Table 5 (Continued)

| Analyte   | EPA Method | Sample Identification | Sample Type | Result | Units | RPD of Primary and Duplicate | RPD of Primary and Triplicate | Mean   | Standard Deviation | RSD   |
|---|------------|-----------------------|-------------|--------|-------|------------------------------|-------------------------------|--------|--------------------|-------|
| <b>Decision Unit 17 - Replicate Samples (continued)</b> |            |                       |             |        |       |                              |                               |        |                    |       |
| Pyrene  | 8270C      | RHS-DU17              | Primary     | 0.17   | mg/kg | 19.4%                        | 64.0%                         | 0.21   | 0.10               | 47.9% |
|   |            | RHS-DU17.2            | Duplicate   | 0.14   | mg/kg |                              |                               |        |                    |       |
|   |            | RHS-DU17.3            | Triplicate  | 0.33   | mg/kg |                              |                               |        |                    |       |
| PCB-1260  | 8082       | RHS-DU17              | Primary     | 2.23   | mg/kg | 5.2%                         | 0.4%                          | 2.27   | 0.07               | 2.9%  |
|   |            | RHS-DU17.2            | Duplicate   | 2.35   | mg/kg |                              |                               |        |                    |       |
|   |            | RHS-DU17.3            | Triplicate  | 2.24   | mg/kg |                              |                               |        |                    |       |
| p,p'-DDT  | 8081A      | RHS-DU17              | Primary     | 0.164  | mg/kg | 17.3%                        | 6.5%                          | 0.18   | 0.02               | 8.8%  |
|   |            | RHS-DU17.2            | Duplicate   | 0.195  | mg/kg |                              |                               |        |                    |       |
|   |            | RHS-DU17.3            | Triplicate  | 0.175  | mg/kg |                              |                               |        |                    |       |
| <b>Decision Unit 26 - Replicate Samples</b>             |            |                       |             |        |       |                              |                               |        |                    |       |
| Dioxin  | 8290       | RHS-DU26              | Primary     | 33     | ng/kg | 53.8%                        | 49.1%                         | 24.00  | 7.81               | 32.5% |
|   |            | RHS-DU26.2            | Duplicate   | 19     | ng/kg |                              |                               |        |                    |       |
|   |            | RHS-DU26.3            | Triplicate  | 20     | ng/kg |                              |                               |        |                    |       |
| Barium  | 6010B      | RHS-DU26              | Primary     | 100    | mg/kg | 36.7%                        | 16.2%                         | 84.67  | 15.50              | 18.3% |
|   |            | RHS-DU26.2            | Duplicate   | 69     | mg/kg |                              |                               |        |                    |       |
|   |            | RHS-DU26.3            | Triplicate  | 85     | mg/kg |                              |                               |        |                    |       |
| Chromium  | 6010B      | RHS-DU26              | Primary     | 250    | mg/kg | 12.8%                        | 27.3%                         | 220.00 | 30.00              | 13.6% |
|   |            | RHS-DU26.2            | Duplicate   | 220    | mg/kg |                              |                               |        |                    |       |
|   |            | RHS-DU26.3            | Triplicate  | 190    | mg/kg |                              |                               |        |                    |       |
| Lead  | 6010B      | RHS-DU26              | Primary     | 89     | mg/kg | 56.1%                        | 13.2%                         | 72.33  | 20.11              | 27.8% |
|   |            | RHS-DU26.2            | Duplicate   | 50     | mg/kg |                              |                               |        |                    |       |
|   |            | RHS-DU26.3            | Triplicate  | 78     | mg/kg |                              |                               |        |                    |       |
| Mercury   | 7471       | RHS-DU26              | Primary     | 3.5    | mg/kg | 25.8%                        | 18.8%                         | 3.03   | 0.42               | 13.7% |
|   |            | RHS-DU26.2            | Duplicate   | 2.7    | mg/kg |                              |                               |        |                    |       |
|   |            | RHS-DU26.3            | Triplicate  | 2.9    | mg/kg |                              |                               |        |                    |       |

Notes on page 4





**Table 5 (Continued)**

Notes:

|       |                             |
|-------|-----------------------------|
| ng/kg | Nanograms per kilogram      |
| mg/kg | Milligrams per kilogram     |
| RPD   | Relative Percent Difference |
| RSD   | Relative Standard Deviation |

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## PHOTOGRAPHS

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|  |                    |  |                      |
|--|--------------------|--|----------------------|
| <b>Project No.</b><br>17012-012148.00<br>Task 48 | <b>Description</b> | View of the excavation pits located on the northeaster portion of the Radford High School track, looking north | <b>Photo 1</b>       |
|  | <b>Site Name</b>   | Radford High School, 4361 Salt Lake Blvd., Honolulu, Hawaii 96816  | <b>Photo Date</b>    |
|  | <b>Client</b>      | State of Hawaii Department of Education  | <b>Dec. 20, 2013</b> |



|  |                    |  |                      |
|--|--------------------|--|----------------------|
| <b>Project No.</b><br>17012-012148.00<br>Task 48 | <b>Description</b> | View of the excavation pit where asbestos containing material was found, looking north | <b>Photo 2</b>       |
|  | <b>Site Name</b>   | Radford High School, 4361 Salt Lake Blvd., Honolulu, Hawaii 96816                      | <b>Photo Date</b>    |
|  | <b>Client</b>      | State of Hawaii Department of Education  | <b>Dec. 20, 2013</b> |





|  |                    |   |                      |
|--|--------------------|---|----------------------|
| <b>Project No.</b><br>17012-012148.00<br>Task 48 | <b>Description</b> | Close up view of the asbestos containing material found under the Radford High School track | <b>Photo 3</b>       |
|  | <b>Site Name</b>   | Radford High School, 4361 Salt Lake Blvd., Honolulu, Hawaii 96816                           | <b>Photo Date</b>    |
|  | <b>Client</b>      | State of Hawaii Department of Education   | <b>Dec. 20, 2013</b> |



|  |                    |   |                      |
|--|--------------------|---|----------------------|
| <b>Project No.</b><br>17012-012148.00<br>Task 48 | <b>Description</b> | Close up view of the suspect asbestos containing material found under the Radford High School track | <b>Photo 4</b>       |
|  | <b>Site Name</b>   | Radford High School, 4361 Salt Lake Blvd., Honolulu, Hawaii 96816                                   | <b>Photo Date</b>    |
|  | <b>Client</b>      | State of Hawaii Department of Education   | <b>Dec. 20, 2013</b> |



|  |                    |   |                     |
|--|--------------------|---|---------------------|
| <b>Project No.</b><br>17012-012148.00<br>Task 48 | <b>Description</b> | View of the entrance to the site, looking north                   | <b>Photo 5</b>      |
|  | <b>Site Name</b>   | Radford High School, 4361 Salt Lake Blvd., Honolulu, Hawaii 96816 | <b>Photo Date</b>   |
|  | <b>Client</b>      | State of Hawaii Department of Education                           | <b>Feb. 5, 2014</b> |



|  |                    |  |                     |
|--|--------------------|--|---------------------|
| <b>Project No.</b><br>17012-012148.00<br>Task 48 | <b>Description</b> | View of the northern track portion and stockpiles covered by black fabric, looking north | <b>Photo 6</b>      |
|  | <b>Site Name</b>   | Radford High School, 4361 Salt Lake Blvd., Honolulu, Hawaii 96816                        | <b>Photo Date</b>   |
|  | <b>Client</b>      | State of Hawaii Department of Education  | <b>Feb. 5, 2014</b> |





|  |                    |   |                     |
|--|--------------------|---|---------------------|
| <b>Project No.</b><br>17012-012148.00<br>Task 48 | <b>Description</b> | View of Radford High School field, looking north                  | <b>Photo 7</b>      |
|  | <b>Site Name</b>   | Radford High School, 4361 Salt Lake Blvd., Honolulu, Hawaii 96816 | <b>Photo Date</b>   |
|  | <b>Client</b>      | State of Hawaii Department of Education                           | <b>Feb. 5, 2014</b> |



|  |                    |   |                     |
|--|--------------------|---|---------------------|
| <b>Project No.</b><br>17012-012148.00<br>Task 48 | <b>Description</b> | View of the southern portion of the track, looking west           | <b>Photo 8</b>      |
|  | <b>Site Name</b>   | Radford High School, 4361 Salt Lake Blvd., Honolulu, Hawaii 96816 | <b>Photo Date</b>   |
|  | <b>Client</b>      | State of Hawaii Department of Education                           | <b>Feb. 5, 2014</b> |





|  |                    |   |                     |
|--|--------------------|---|---------------------|
| <b>Project No.</b><br>17012-012148.00<br>Task 48 | <b>Description</b> | View of the western portion of the track, looking south           | <b>Photo 9</b>      |
|  | <b>Site Name</b>   | Radford High School, 4361 Salt Lake Blvd., Honolulu, Hawaii 96816 | <b>Photo Date</b>   |
|  | <b>Client</b>      | State of Hawaii Department of Education                           | <b>Feb. 5, 2014</b> |



|  |                    |   |                     |
|--|--------------------|---|---------------------|
| <b>Project No.</b><br>17012-012148.00<br>Task 48 | <b>Description</b> | View of the eastern portion of the track, looking south           | <b>Photo 10</b>     |
|  | <b>Site Name</b>   | Radford High School, 4361 Salt Lake Blvd., Honolulu, Hawaii 96816 | <b>Photo Date</b>   |
|  | <b>Client</b>      | State of Hawaii Department of Education                           | <b>Feb. 5, 2014</b> |



**APPENDIX A**

**LABORATORY ANALYTICAL REPORTS  
AND  
CHAIN OF CUSTODY FORMS (TESTAMERICA)**

# TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

## ANALYTICAL REPORT

TestAmerica Laboratories, Inc.  
TestAmerica Honolulu  
1946 Young St. Suite 400A  
Honolulu, HI 96826  
Tel: 808-486-5227

TestAmerica Job ID: HWL0089  
Client Project/Site: 17012-012148.00  
Client Project Description: Radford HS

For:  
Bureau Veritas  
841 Bishop Street, Suite 1100  
Honolulu, HI 96813

Attn: Kenney Gomes



Authorized for release by:  
1/24/2014 11:10:32 AM

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*This report has been electronically signed and authorized by the signatory. Electronic signature is intended to be the legally binding equivalent of a traditionally handwritten signature.*

*Results relate only to the items tested and the sample(s) as received by the laboratory.*

- 1
- 2
- 3
- 4
- 5
- 6
- 7
- 8
- 9
- 10
- 11
- 12
- 13
- 14
- 15
- 16



# Table of Contents

|                                    |    |
|------------------------------------|----|
| Cover Page . . . . .               | 1  |
| Table of Contents . . . . .        | 2  |
| Definitions . . . . .              | 3  |
| Case Narrative . . . . .           | 4  |
| Sample Summary . . . . .           | 7  |
| Detection Summary . . . . .        | 8  |
| Client Sample Results . . . . .    | 9  |
| Surrogate Summary . . . . .        | 13 |
| QC Sample Results . . . . .        | 15 |
| QC Association . . . . .           | 28 |
| Chronicle . . . . .                | 31 |
| Certification Summary . . . . .    | 32 |
| Method Summary . . . . .           | 34 |
| Chain of Custody . . . . .         | 35 |
| Toxicity Summary . . . . .         | 37 |
| Isotope Dilution Summary . . . . . | 38 |

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# Definitions/Glossary

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Qualifiers

### GC/MS Semi VOA

| Qualifier | Qualifier Description  |
|-----------|--|
| *         | RPD of the LCS and LCSD exceeds the control limits   |
| J         | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |

### GC Semi VOA

| Qualifier | Qualifier Description  |
|-----------|--|
| p         | The %RPD between the primary and confirmation column/detector is >40%. The lower value has been reported.      |
| J         | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |

### Dioxin

| Qualifier | Qualifier Description  |
|-----------|--|
| B         | Compound was found in the blank and sample.  |
| G         | The reported quantitation limit has been raised due to an exhibited elevated noise or matrix interference  |
| q         | The isomer is qualified as positively identified, but at an estimated quantity because the quantitation is based on the theoretical ratio for these samples. |
| J         | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value.   |

### Metals

| Qualifier | Qualifier Description   |
|-----------|---|
| 4         | MS, MSD: The analyte present in the original sample is greater than 4 times the matrix spike concentration; therefore, control limits are not applicable. |
| F1        | MS and/or MSD Recovery exceeds the control limits   |
| F2        | MS/MSD RPD exceeds control limits   |

## Glossary

| Abbreviation   | These commonly used abbreviations may or may not be present in this report.                                 |
|----------------|---|
| ▫              | Listed under the "D" column to designate that the result is reported on a dry weight basis                  |
| %R             | Percent Recovery  |
| CNF            | Contains no Free Liquid   |
| DER            | Duplicate error ratio (normalized absolute difference)  |
| Dil Fac        | Dilution Factor   |
| DL, RA, RE, IN | Indicates a Dilution, Re-analysis, Re-extraction, or additional Initial metals/anion analysis of the sample |
| DLC            | Decision level concentration  |
| MDA            | Minimum detectable activity   |
| EDL            | Estimated Detection Limit   |
| MDC            | Minimum detectable concentration  |
| MDL            | Method Detection Limit  |
| ML             | Minimum Level (Dioxin)  |
| NC             | Not Calculated  |
| ND             | Not detected at the reporting limit (or MDL or EDL if shown)  |
| PQL            | Practical Quantitation Limit  |
| QC             | Quality Control   |
| RER            | Relative error ratio  |
| RL             | Reporting Limit or Requested Limit (Radiochemistry)   |
| RPD            | Relative Percent Difference, a measure of the relative difference between two points                        |
| TEF            | Toxicity Equivalent Factor (Dioxin)   |
| TEQ            | Toxicity Equivalent Quotient (Dioxin)   |

TestAmerica Honolulu

# Case Narrative

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Job ID: HWL0089

### Laboratory: TestAmerica Honolulu

#### Narrative

The results listed within this Laboratory Report pertain only to the samples tested in the laboratory unless otherwise stated in the report. The analyses contained in this report were performed in accordance with the applicable certifications as noted. All soil samples are reported on a wet weight basis unless otherwise noted in the report. This Laboratory Report is confidential and is intended for the sole use of TestAmerica and its client. This report shall not be reproduced, except in full, without written permission from TestAmerica. TestAmerica Analytical Testing Corporation certifies that the analytical results contained herein apply only to the specific sample(s) analyzed.

The Chain(s) of Custody are included and are an integral part of this report. This entire report was reviewed and approved for release.

If you have any questions relating to this analytical report, please contact your Laboratory Project Manager at 1-(808)486-5227

#### LABORATORY REPORT

At sample receipt, the cooler/sample was 2 degrees C.

NELAC states that samples which require thermal preservation shall be considered acceptable if the arrival temperature is within 2 degrees C of the required temperature or the method specified range. For samples with a temperature requirement of 4 degrees C, an arrival temperature from 0 degrees C to 6 degrees C meets specifications. Samples that are delivered to the laboratory on the same day that they are collected may not meet these criteria. In these cases, the samples are considered acceptable if there is evidence that the chilling process has begun, such as arrival on ice.

The reported results were obtained in compliance with the 2003 NELAC standards unless otherwise noted.

Samples were prepared in accordance with the State of Hawai'i Department of Health Office of Hazard Evaluation and Emergency Response's Technical Guidance Manual for the Implementation of the Hawai'i State Contingency Plan 2009 edition Laboratory Preparation of Multi-Increment Samples.

### Laboratory: TestAmerica Irvine

#### Narrative

Job Narrative  
440-66219-1

#### Comments

No additional comments.

#### Receipt

The sample was received on 12/27/2013 9:15 AM; the sample arrived in good condition, properly preserved and, where required, on ice. The temperature of the cooler at receipt was 4.9° C.

#### GC/MS Semi VOA

Method(s) 8270C: Insufficient sample volume was available to perform batch matrix spike/matrix spike duplicate (MS/MSD) associated with batch 153305. The laboratory control sample (LCS) was performed in duplicate to provide precision data for this batch.

Method(s) 8270C: The %RPD of the laboratory control sample (LCS) and laboratory control standard duplicate (LCSD) for preparation batch 153305 recovered outside control limits for the following analytes: 4-nitrophenol, diethyl phthalate, 3-nitroaniline, 2,4-dinitrotoluene, and 4-nitroaniline. The individual LCS/LCSD percent recoveries met acceptance limit.

Method(s) 8270C: The following sample(s) was diluted due to the abundance of non-target analytes: HW0089-01 (440-66219-1). Elevated reporting limits (RLs) are provided.

No other analytical or quality issues were noted.



# Case Narrative

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Job ID: HWL0089 (Continued)

### Laboratory: TestAmerica Irvine (Continued)

#### GC Semi VOA

Method(s) 8081A: The continuing calibration verification (CCV) associated with batch 153523 recovered above the upper control limit for 4,4 DDD. The samples associated with this CCV were non-detects for the affected analytes; therefore, the data have been reported. The following samples are impacted: (CCVRT 440-153523/15), HW0089-01 (440-66219-1).

Method(s) 8081A: The capping continuing calibration verification (CCV) associated with batch 153523 did not meet control limits for 4,4 DDD, 4,4 DDT, and Methoxychlor. Samples were re-analyzed for confirmation of results. Sample matrix is suspected to have contributed to this failure. (CCV 440-153523/27), HW0089-01 (440-66219-1)

Method(s) 8082: The %RPD between the primary and confirmation column exceeded 40% for the DCBP for the following sample(s): (440-66257-2 MS), (440-66257-2 MSD). The lower value has been reported and qualified in accordance with the laboratory's SOP.

Method(s) 8082: The %RPD between the primary and confirmation column(s) exceeded 40% for 1260 for the following sample(s): HW0089-01 (440-66219-1). The lower value has been reported and qualified in accordance with the laboratory's SOP.

No other analytical or quality issues were noted.

#### Organic Prep

No analytical or quality issues were noted.

Laboratory: TestAmerica Sacramento

#### Narrative

Job Narrative  
320-5498-1

#### Receipt

The sample was received on 12/27/2013 8:55 AM; the sample arrived in good condition, properly preserved and, where required, on ice. The temperature of the cooler at receipt was 3.8° C.

#### Dioxin

Method(s) 8290:

The following samples were diluted to bring the concentration of one or more target analytes within the calibration range: HWL0089-01 (320-5498-1). Elevated reporting limits (RLs) are provided.

The following sample [320-5498-1] exhibited elevated noise or matrix interferences for one or more analytes requiring the detection limits to be raised appropriately. These analytes were flagged with the "G" qualifier.

The closing bracketing continuing calibration verification (CCV) associated with analytical batch 34312 has the analyte 2,3,7,8-TCDF with a percent difference value that is between the method criteria of 20% to 25% deviation from the initial calibration curve. Per method guidelines, an average relative response factor (RRF) is calculated from the opening and closing bracketing CCV's and is used to quantitate any positive results in the associated samples for the affected analyte.

Ion abundance ratios are outside criteria for one or more analytes in the following samples: (MB 320-33622/1-A), HWL0089-01 (320-5498-1). Quantitation is based on the theoretical ion abundance ratio; therefore, these analytes have been reported as an estimated maximum possible concentration (EMPC). The affected analytes have been flagged.

No other analytical or quality issues were noted.

#### General Chemistry

No analytical or quality issues were noted.

#### Dioxin Prep

No analytical or quality issues were noted.

Laboratory: TestAmerica Seattle

# Case Narrative

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

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## Job ID: HWL0089 (Continued)

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### Laboratory: TestAmerica Seattle (Continued)

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#### Narrative

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#### Receipt

The sample was received on 12/27/2013 9:55 AM; the sample arrived in good condition, properly preserved and, where required, on ice. The temperature of the cooler at receipt was 3.7° C.

#### Metals - Method(s) 6010B

The matrix spike / matrix spike duplicate (MS/MSD) recoveries for analysis batch 151504 were outside control limits. Sample matrix interference is suspected because the associated laboratory control sample (LCS) recovery was within acceptance limits. The data have been "F" qualified and reported.

No other analytical or quality issues were noted.

#### General Chemistry

No analytical or quality issues were noted.

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- 1
- 2
- 3
- 4
- 5
- 6
- 7
- 8
- 9
- 10
- 11
- 12
- 13
- 14
- 15
- 16

# Sample Summary

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

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| Lab Sample ID | Client Sample ID | Matrix     | Collected      | Received       |
|---------------|------------------|------------|----------------|----------------|
| HWL0089-01    | RHS-01           | Solid/Soil | 12/20/13 15:00 | 12/23/13 13:10 |

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- 1
- 2
- 3
- 4
- 5
- 6
- 7
- 8
- 9
- 10
- 11
- 12
- 13
- 14
- 15
- 16

# Detection Summary

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

**Client Sample ID: RHS-01**

**Lab Sample ID: HWL0089-01**

| Analyte                | Result  | Qualifier | RL     | EDL     | Unit  | Dil Fac | D | Method | Prep Type |
|------------------------|---------|-----------|--------|---------|-------|---------|---|--------|-----------|
| 2,3,7,8-TCDD           | 88      |           | 21     | 11      | pg/g  | 20      | * | 8290   | Total/NA  |
| 1,2,3,7,8-PeCDD        | 340     |           | 100    | 52      | pg/g  | 20      | * | 8290   | Total/NA  |
| 1,2,3,7,8-PeCDF        | 1100    |           | 100    | 49      | pg/g  | 20      | * | 8290   | Total/NA  |
| 2,3,4,7,8-PeCDF        | 2300    |           | 100    | 52      | pg/g  | 20      | * | 8290   | Total/NA  |
| 1,2,3,4,7,8-HxCDD      | 250     |           | 100    | 11      | pg/g  | 20      | * | 8290   | Total/NA  |
| 1,2,3,6,7,8-HxCDD      | 420     |           | 100    | 8.5     | pg/g  | 20      | * | 8290   | Total/NA  |
| 1,2,3,7,8,9-HxCDD      | 320     |           | 100    | 8.2     | pg/g  | 20      | * | 8290   | Total/NA  |
| 1,2,3,4,7,8-HxCDF      | 1900    | B         | 100    | 85      | pg/g  | 20      | * | 8290   | Total/NA  |
| 1,2,3,6,7,8-HxCDF      | 1400    | B         | 100    | 71      | pg/g  | 20      | * | 8290   | Total/NA  |
| 2,3,4,6,7,8-HxCDF      | 1900    | B         | 100    | 79      | pg/g  | 20      | * | 8290   | Total/NA  |
| 1,2,3,4,6,7,8-HpCDD    | 2400    | B         | 100    | 16      | pg/g  | 20      | * | 8290   | Total/NA  |
| 1,2,3,4,6,7,8-HpCDF    | 7600    | B         | 100    | 98      | pg/g  | 20      | * | 8290   | Total/NA  |
| 1,2,3,4,7,8,9-HpCDF    | 220     | G         | 120    | 120     | pg/g  | 20      | * | 8290   | Total/NA  |
| OCDD                   | 5600    | B         | 210    | 19      | pg/g  | 20      | * | 8290   | Total/NA  |
| OCDF                   | 1700    | B         | 210    | 5.7     | pg/g  | 20      | * | 8290   | Total/NA  |
| Total TCDD             | 4400    | q B       | 21     | 11      | pg/g  | 20      | * | 8290   | Total/NA  |
| Total TCDF             | 38000   | G         | 110    | 110     | pg/g  | 20      | * | 8290   | Total/NA  |
| Total PeCDD            | 5000    | q         | 100    | 52      | pg/g  | 20      | * | 8290   | Total/NA  |
| Total PeCDF            | 24000   | q         | 100    | 50      | pg/g  | 20      | * | 8290   | Total/NA  |
| Total HxCDD            | 5700    |           | 100    | 9.1     | pg/g  | 20      | * | 8290   | Total/NA  |
| Total HxCDF            | 16000   | q B       | 100    | 81      | pg/g  | 20      | * | 8290   | Total/NA  |
| Total HpCDD            | 4900    | B         | 100    | 16      | pg/g  | 20      | * | 8290   | Total/NA  |
| Total HpCDF            | 8900    | G B       | 110    | 110     | pg/g  | 20      | * | 8290   | Total/NA  |
| 2,3,7,8-TCDF - RA      | 1400    |           | 21     | 3.2     | pg/g  | 20      | * | 8290   | Total/NA  |
| Analyte                | Result  | Qualifier | RL     | MDL     | Unit  | Dil Fac | D | Method | Prep Type |
| 2-Methylnaphthalene    | 0.25    | J         | 0.69   | 0.15    | mg/Kg | 4       | * | 8270C  | Total/NA  |
| Benzo[b]fluoranthene   | 0.15    | J         | 0.69   | 0.14    | mg/Kg | 4       | * | 8270C  | Total/NA  |
| Indeno[1,2,3-cd]pyrene | 0.49    | J         | 0.69   | 0.27    | mg/Kg | 4       | * | 8270C  | Total/NA  |
| Naphthalene            | 2.7     |           | 0.69   | 0.14    | mg/Kg | 4       | * | 8270C  | Total/NA  |
| DRO (C10-C28)          | 24      |           | 5.2    | 2.6     | mg/Kg | 1       | * | 8015B  | Total/NA  |
| RRO(C29-C40)           | 23      |           | 5.2    | 2.6     | mg/Kg | 1       | * | 8015B  | Total/NA  |
| 4,4'-DDE               | 0.0046  |           | 0.0026 | 0.00078 | mg/Kg | 1       | * | 8081A  | Total/NA  |
| 4,4'-DDT               | 0.0045  | p         | 0.0026 | 0.00078 | mg/Kg | 1       | * | 8081A  | Total/NA  |
| Dieldrin               | 0.0035  |           | 0.0026 | 0.00078 | mg/Kg | 1       | * | 8081A  | Total/NA  |
| Endrin                 | 0.00086 | J p       | 0.0026 | 0.00078 | mg/Kg | 1       | * | 8081A  | Total/NA  |
| Endrin aldehyde        | 0.0035  | p         | 0.0026 | 0.00078 | mg/Kg | 1       | * | 8081A  | Total/NA  |
| Aroclor 1260           | 0.094   | p         | 0.026  | 0.0089  | mg/Kg | 1       | * | 8082   | Total/NA  |
| Arsenic                | 43      |           | 6.2    |         | mg/Kg | 10      | * | 6010B  | Total/NA  |
| Barium                 | 940     |           | 1.0    |         | mg/Kg | 10      | * | 6010B  | Total/NA  |
| Cadmium                | 26      |           | 2.1    |         | mg/Kg | 10      | * | 6010B  | Total/NA  |
| Chromium               | 210     |           | 2.7    |         | mg/Kg | 10      | * | 6010B  | Total/NA  |
| Lead                   | 5300    |           | 3.1    |         | mg/Kg | 10      | * | 6010B  | Total/NA  |
| Silver                 | 14      |           | 5.2    |         | mg/Kg | 10      | * | 6010B  | Total/NA  |
| Mercury                | 1.1     |           | 0.020  |         | mg/Kg | 10      | * | 7471A  | Total/NA  |

This Detection Summary does not include radiochemical test results.

TestAmerica Honolulu

# Client Sample Results

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

**Client Sample ID: RHS-01**

**Lab Sample ID: HWL0089-01**

**Date Collected: 12/20/13 15:00**

**Matrix: Solid/Soil**

**Date Received: 12/23/13 13:10**

**Percent Solids: 95.1**

**Method: 8270C - Semivolatile Organic Compounds (GC/MS)**

| Analyte                         | Result      | Qualifier | RL   | MDL  | Unit  | D | Prepared       | Analyzed       | Dil Fac |
|---------------------------------|-------------|-----------|------|------|-------|---|----------------|----------------|---------|
| 1,2,4-Trichlorobenzene          | ND          |           | 0.69 | 0.28 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 1,2-Dichlorobenzene             | ND          |           | 0.69 | 0.13 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 1,3-Dichlorobenzene             | ND          |           | 0.69 | 0.28 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 1,4-Dichlorobenzene             | ND          |           | 0.69 | 0.28 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 1-Methylnaphthalene             | ND          |           | 0.73 | 0.31 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 2,4,5-Trichlorophenol           | ND          |           | 0.69 | 0.27 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 2,4,6-Trichlorophenol           | ND          |           | 0.69 | 0.16 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 2,4-Dichlorophenol              | ND          |           | 0.69 | 0.14 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 2,4-Dimethylphenol              | ND          |           | 0.69 | 0.27 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 2,4-Dinitrophenol               | ND          |           | 1.4  | 0.69 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 2,4-Dinitrotoluene              | ND          | *         | 0.69 | 0.17 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 2,6-Dinitrotoluene              | ND          |           | 0.69 | 0.20 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 2-Chloronaphthalene             | ND          |           | 0.69 | 0.14 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 2-Chlorophenol                  | ND          |           | 0.69 | 0.15 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| <b>2-Methylnaphthalene</b>      | <b>0.25</b> | <b>J</b>  | 0.69 | 0.15 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 2-Methylphenol                  | ND          |           | 0.69 | 0.17 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 2-Nitroaniline                  | ND          |           | 0.69 | 0.14 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 2-Nitrophenol                   | ND          |           | 0.69 | 0.28 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 3,3'-Dichlorobenzidine          | ND          |           | 1.7  | 0.31 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 3-Methylphenol + 4-Methylphenol | ND          |           | 0.69 | 0.28 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 3-Nitroaniline                  | ND          | *         | 0.69 | 0.28 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 4,6-Dinitro-2-methylphenol      | ND          |           | 0.88 | 0.28 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 4-Bromophenyl phenyl ether      | ND          |           | 0.69 | 0.16 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 4-Chloro-3-methylphenol         | ND          |           | 0.69 | 0.15 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 4-Chloroaniline                 | ND          |           | 0.69 | 0.28 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 4-Chlorophenyl phenyl ether     | ND          |           | 0.69 | 0.18 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 4-Nitroaniline                  | ND          | *         | 1.7  | 0.28 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 4-Nitrophenol                   | ND          | *         | 1.7  | 0.29 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Acenaphthene                    | ND          |           | 0.69 | 0.14 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Acenaphthylene                  | ND          |           | 0.69 | 0.15 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Anthracene                      | ND          |           | 0.69 | 0.17 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Benzo[a]anthracene              | ND          |           | 0.69 | 0.15 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Benzo[a]pyrene                  | ND          |           | 0.69 | 0.14 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| <b>Benzo[b]fluoranthene</b>     | <b>0.15</b> | <b>J</b>  | 0.69 | 0.14 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Benzo[g,h,i]perylene            | ND          |           | 0.69 | 0.23 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Benzo[k]fluoranthene            | ND          |           | 0.69 | 0.15 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Benzoic acid                    | ND          |           | 1.7  | 0.31 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Benzyl alcohol                  | ND          |           | 0.69 | 0.42 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| bis (2-chloroisopropyl) ether   | ND          |           | 0.69 | 0.28 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Bis(2-chloroethoxy)methane      | ND          |           | 0.69 | 0.28 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Bis(2-chloroethyl)ether         | ND          |           | 0.69 | 0.13 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Bis(2-ethylhexyl) phthalate     | ND          |           | 0.69 | 0.19 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Butyl benzyl phthalate          | ND          |           | 0.69 | 0.17 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Carbazole                       | ND          |           | 0.69 | 0.28 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Chrysene                        | ND          |           | 0.69 | 0.16 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Dibenz(a,h)anthracene           | ND          |           | 0.88 | 0.21 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Dibenzofuran                    | ND          |           | 0.69 | 0.14 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Diethyl phthalate               | ND          | *         | 0.69 | 0.20 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Dimethyl phthalate              | ND          |           | 0.69 | 0.14 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |

TestAmerica Honolulu

# Client Sample Results

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

**Client Sample ID: RHS-01**

**Lab Sample ID: HWL0089-01**

**Date Collected: 12/20/13 15:00**

**Matrix: Solid/Soil**

**Date Received: 12/23/13 13:10**

**Percent Solids: 95.1**

**Method: 8270C - Semivolatile Organic Compounds (GC/MS) (Continued)**

| Analyte                       | Result      | Qualifier | RL   | MDL  | Unit  | D | Prepared       | Analyzed       | Dil Fac |
|-------------------------------|-------------|-----------|------|------|-------|---|----------------|----------------|---------|
| Di-n-butyl phthalate          | ND          |           | 0.69 | 0.19 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Di-n-octyl phthalate          | ND          |           | 0.69 | 0.19 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Fluoranthene                  | ND          |           | 0.69 | 0.15 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Fluorene                      | ND          |           | 0.69 | 0.15 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Hexachlorobenzene             | ND          |           | 0.69 | 0.15 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Hexachlorobutadiene           | ND          |           | 0.69 | 0.28 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Hexachlorocyclopentadiene     | ND          |           | 1.7  | 0.28 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Hexachloroethane              | ND          |           | 0.69 | 0.28 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| <b>Indeno[1,2,3-cd]pyrene</b> | <b>0.49</b> | <b>J</b>  | 0.69 | 0.27 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Isophorone                    | ND          |           | 0.69 | 0.14 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| <b>Naphthalene</b>            | <b>2.7</b>  |           | 0.69 | 0.14 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Nitrobenzene                  | ND          |           | 0.69 | 0.15 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| N-Nitrosodi-n-propylamine     | ND          |           | 0.52 | 0.15 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| N-Nitrosodiphenylamine        | ND          |           | 0.69 | 0.17 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Pentachlorophenol             | ND          |           | 1.7  | 0.31 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Phenanthrene                  | ND          |           | 0.69 | 0.14 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Phenol                        | ND          |           | 0.69 | 0.19 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Pyrene                        | ND          |           | 0.69 | 0.17 | mg/Kg | ☼ | 12/30/13 11:44 | 01/02/14 16:39 | 4       |

| Surrogate                   | %Recovery | Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|-----------------------------|-----------|-----------|----------|----------------|----------------|---------|
| 2,4,6-Tribromophenol (Surr) | 40        |           | 35 - 125 | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 2-Fluorophenol (Surr)       | 41        |           | 25 - 120 | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Nitrobenzene-d5 (Surr)      | 52        |           | 30 - 120 | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Phenol-d6 (Surr)            | 45        |           | 35 - 120 | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| Terphenyl-d14 (Surr)        | 79        |           | 40 - 135 | 12/30/13 11:44 | 01/02/14 16:39 | 4       |
| 2-Fluorobiphenyl            | 69        |           | 35 - 120 | 12/30/13 11:44 | 01/02/14 16:39 | 4       |

**Method: 8015B - Diesel Range Organics (DRO) (GC)**

| Analyte              | Result    | Qualifier | RL  | MDL | Unit  | D | Prepared       | Analyzed       | Dil Fac |
|----------------------|-----------|-----------|-----|-----|-------|---|----------------|----------------|---------|
| <b>DRO (C10-C28)</b> | <b>24</b> |           | 5.2 | 2.6 | mg/Kg | ☼ | 12/28/13 07:03 | 12/30/13 08:36 | 1       |
| <b>RRO(C29-C40)</b>  | <b>23</b> |           | 5.2 | 2.6 | mg/Kg | ☼ | 12/28/13 07:03 | 12/30/13 08:36 | 1       |

| Surrogate    | %Recovery | Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|--------------|-----------|-----------|----------|----------------|----------------|---------|
| n-Octacosane | 63        |           | 40 - 140 | 12/28/13 07:03 | 12/30/13 08:36 | 1       |

**Method: 8081A - Organochlorine Pesticides (GC)**

| Analyte               | Result         | Qualifier  | RL     | MDL     | Unit  | D | Prepared       | Analyzed       | Dil Fac |
|-----------------------|----------------|------------|--------|---------|-------|---|----------------|----------------|---------|
| 4,4'-DDD              | ND             |            | 0.0026 | 0.00078 | mg/Kg | ☼ | 12/27/13 17:32 | 12/31/13 14:53 | 1       |
| <b>4,4'-DDE</b>       | <b>0.0046</b>  |            | 0.0026 | 0.00078 | mg/Kg | ☼ | 12/27/13 17:32 | 12/31/13 14:53 | 1       |
| <b>4,4'-DDT</b>       | <b>0.0045</b>  | <b>p</b>   | 0.0026 | 0.00078 | mg/Kg | ☼ | 12/27/13 17:32 | 12/31/13 14:53 | 1       |
| Aldrin                | ND             |            | 0.0026 | 0.00078 | mg/Kg | ☼ | 12/27/13 17:32 | 12/31/13 14:53 | 1       |
| alpha-BHC             | ND             |            | 0.0026 | 0.00078 | mg/Kg | ☼ | 12/27/13 17:32 | 12/31/13 14:53 | 1       |
| beta-BHC              | ND             |            | 0.0026 | 0.00078 | mg/Kg | ☼ | 12/27/13 17:32 | 12/31/13 14:53 | 1       |
| Chlordane (technical) | ND             |            | 0.026  | 0.0052  | mg/Kg | ☼ | 12/27/13 17:32 | 12/31/13 14:53 | 1       |
| delta-BHC             | ND             |            | 0.0052 | 0.00078 | mg/Kg | ☼ | 12/27/13 17:32 | 12/31/13 14:53 | 1       |
| <b>Dieldrin</b>       | <b>0.0035</b>  |            | 0.0026 | 0.00078 | mg/Kg | ☼ | 12/27/13 17:32 | 12/31/13 14:53 | 1       |
| Endosulfan I          | ND             |            | 0.0026 | 0.00078 | mg/Kg | ☼ | 12/27/13 17:32 | 12/31/13 14:53 | 1       |
| Endosulfan II         | ND             |            | 0.0026 | 0.00078 | mg/Kg | ☼ | 12/27/13 17:32 | 12/31/13 14:53 | 1       |
| Endosulfan sulfate    | ND             |            | 0.0052 | 0.0010  | mg/Kg | ☼ | 12/27/13 17:32 | 12/31/13 14:53 | 1       |
| <b>Endrin</b>         | <b>0.00086</b> | <b>J p</b> | 0.0026 | 0.00078 | mg/Kg | ☼ | 12/27/13 17:32 | 12/31/13 14:53 | 1       |

TestAmerica Honolulu



# Client Sample Results

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

**Client Sample ID: RHS-01**

**Lab Sample ID: HWL0089-01**

**Date Collected: 12/20/13 15:00**

**Matrix: Solid/Soil**

**Date Received: 12/23/13 13:10**

**Percent Solids: 95.1**

**Method: 8081A - Organochlorine Pesticides (GC) (Continued)**

| Analyte                       | Result           | Qualifier        | RL            | MDL     | Unit  | D | Prepared        | Analyzed        | Dil Fac        |
|-------------------------------|------------------|------------------|---------------|---------|-------|---|-----------------|-----------------|----------------|
| <b>Endrin aldehyde</b>        | <b>0.0035</b>    | <b>p</b>         | 0.0026        | 0.00078 | mg/Kg | ☼ | 12/27/13 17:32  | 12/31/13 14:53  | 1              |
| Endrin ketone                 | ND               |                  | 0.0026        | 0.0010  | mg/Kg | ☼ | 12/27/13 17:32  | 12/31/13 14:53  | 1              |
| gamma-BHC (Lindane)           | ND               |                  | 0.0026        | 0.00078 | mg/Kg | ☼ | 12/27/13 17:32  | 12/31/13 14:53  | 1              |
| Heptachlor                    | ND               |                  | 0.0026        | 0.0010  | mg/Kg | ☼ | 12/27/13 17:32  | 12/31/13 14:53  | 1              |
| Heptachlor epoxide            | ND               |                  | 0.0026        | 0.0010  | mg/Kg | ☼ | 12/27/13 17:32  | 12/31/13 14:53  | 1              |
| Methoxychlor                  | ND               |                  | 0.0026        | 0.00078 | mg/Kg | ☼ | 12/27/13 17:32  | 12/31/13 14:53  | 1              |
| Toxaphene                     | ND               |                  | 0.10          | 0.026   | mg/Kg | ☼ | 12/27/13 17:32  | 12/31/13 14:53  | 1              |
| <b>Surrogate</b>              | <b>%Recovery</b> | <b>Qualifier</b> | <b>Limits</b> |         |       |   | <b>Prepared</b> | <b>Analyzed</b> | <b>Dil Fac</b> |
| Tetrachloro-m-xylene          | 67               |                  | 35 - 115      |         |       |   | 12/27/13 17:32  | 12/31/13 14:53  | 1              |
| DCB Decachlorobiphenyl (Surr) | 96               |                  | 45 - 120      |         |       |   | 12/27/13 17:32  | 12/31/13 14:53  | 1              |

**Method: 8082 - Polychlorinated Biphenyls (PCBs) by Gas Chromatography**

| Analyte                       | Result           | Qualifier        | RL            | MDL    | Unit  | D | Prepared        | Analyzed        | Dil Fac        |
|-------------------------------|------------------|------------------|---------------|--------|-------|---|-----------------|-----------------|----------------|
| Aroclor 1016                  | ND               |                  | 0.026         | 0.0089 | mg/Kg | ☼ | 12/30/13 11:45  | 12/30/13 20:23  | 1              |
| Aroclor 1221                  | ND               |                  | 0.026         | 0.0089 | mg/Kg | ☼ | 12/30/13 11:45  | 12/30/13 20:23  | 1              |
| Aroclor 1232                  | ND               |                  | 0.026         | 0.0089 | mg/Kg | ☼ | 12/30/13 11:45  | 12/30/13 20:23  | 1              |
| Aroclor 1242                  | ND               |                  | 0.026         | 0.0089 | mg/Kg | ☼ | 12/30/13 11:45  | 12/30/13 20:23  | 1              |
| Aroclor 1248                  | ND               |                  | 0.026         | 0.0089 | mg/Kg | ☼ | 12/30/13 11:45  | 12/30/13 20:23  | 1              |
| Aroclor 1254                  | ND               |                  | 0.026         | 0.0089 | mg/Kg | ☼ | 12/30/13 11:45  | 12/30/13 20:23  | 1              |
| <b>Aroclor 1260</b>           | <b>0.094</b>     | <b>p</b>         | 0.026         | 0.0089 | mg/Kg | ☼ | 12/30/13 11:45  | 12/30/13 20:23  | 1              |
| <b>Surrogate</b>              | <b>%Recovery</b> | <b>Qualifier</b> | <b>Limits</b> |        |       |   | <b>Prepared</b> | <b>Analyzed</b> | <b>Dil Fac</b> |
| DCB Decachlorobiphenyl (Surr) | 78               |                  | 45 - 120      |        |       |   | 12/30/13 11:45  | 12/30/13 20:23  | 1              |

**Method: 8290 - Dioxins and Furans (HRGC/HRMS)**

| Analyte                    | Result       | Qualifier  | RL  | EDL | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------------------------|--------------|------------|-----|-----|------|---|----------------|----------------|---------|
| <b>2,3,7,8-TCDD</b>        | <b>88</b>    |            | 21  | 11  | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>1,2,3,7,8-PeCDD</b>     | <b>340</b>   |            | 100 | 52  | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>1,2,3,7,8-PeCDF</b>     | <b>1100</b>  |            | 100 | 49  | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>2,3,4,7,8-PeCDF</b>     | <b>2300</b>  |            | 100 | 52  | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>1,2,3,4,7,8-HxCDD</b>   | <b>250</b>   |            | 100 | 11  | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>1,2,3,6,7,8-HxCDD</b>   | <b>420</b>   |            | 100 | 8.5 | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>1,2,3,7,8,9-HxCDD</b>   | <b>320</b>   |            | 100 | 8.2 | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>1,2,3,4,7,8-HxCDF</b>   | <b>1900</b>  | <b>B</b>   | 100 | 85  | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>1,2,3,6,7,8-HxCDF</b>   | <b>1400</b>  | <b>B</b>   | 100 | 71  | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>2,3,4,6,7,8-HxCDF</b>   | <b>1900</b>  | <b>B</b>   | 100 | 79  | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| 1,2,3,7,8,9-HxCDF          | ND           |            | 100 | 89  | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>1,2,3,4,6,7,8-HpCDD</b> | <b>2400</b>  | <b>B</b>   | 100 | 16  | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>1,2,3,4,6,7,8-HpCDF</b> | <b>7600</b>  | <b>B</b>   | 100 | 98  | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>1,2,3,4,7,8,9-HpCDF</b> | <b>220</b>   | <b>G</b>   | 120 | 120 | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>OCDD</b>                | <b>5600</b>  | <b>B</b>   | 210 | 19  | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>OCDF</b>                | <b>1700</b>  | <b>B</b>   | 210 | 5.7 | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>Total TCDD</b>          | <b>4400</b>  | <b>q B</b> | 21  | 11  | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>Total TCDF</b>          | <b>38000</b> | <b>G</b>   | 110 | 110 | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>Total PeCDD</b>         | <b>5000</b>  | <b>q</b>   | 100 | 52  | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>Total PeCDF</b>         | <b>24000</b> | <b>q</b>   | 100 | 50  | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>Total HxCDD</b>         | <b>5700</b>  |            | 100 | 9.1 | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>Total HxCDF</b>         | <b>16000</b> | <b>q B</b> | 100 | 81  | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |
| <b>Total HpCDD</b>         | <b>4900</b>  | <b>B</b>   | 100 | 16  | pg/g | ☼ | 01/10/14 11:33 | 01/20/14 16:28 | 20      |

TestAmerica Honolulu

# Client Sample Results

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

**Client Sample ID: RHS-01**

**Lab Sample ID: HWL0089-01**

**Date Collected: 12/20/13 15:00**

**Matrix: Solid/Soil**

**Date Received: 12/23/13 13:10**

**Percent Solids: 94.9**

**Method: 8290 - Dioxins and Furans (HRGC/HRMS) (Continued)**

| Analyte                 | Result           | Qualifier        | RL            | EDL | Unit | D | Prepared        | Analyzed        | Dil Fac        |
|-------------------------|------------------|------------------|---------------|-----|------|---|-----------------|-----------------|----------------|
| <b>Total HpCDF</b>      | <b>8900</b>      | <b>G B</b>       | 110           | 110 | pg/g | ☼ | 01/10/14 11:33  | 01/20/14 16:28  | 20             |
| <i>Isotope Dilution</i> | <i>%Recovery</i> | <i>Qualifier</i> | <i>Limits</i> |     |      |   | <i>Prepared</i> | <i>Analyzed</i> | <i>Dil Fac</i> |
| 13C-2,3,7,8-TCDD        | 73               |                  | 40 - 135      |     |      |   | 01/10/14 11:33  | 01/20/14 16:28  | 20             |
| 13C-2,3,7,8-TCDF        | 78               |                  | 40 - 135      |     |      |   | 01/10/14 11:33  | 01/20/14 16:28  | 20             |
| 13C-1,2,3,7,8-PeCDD     | 80               |                  | 40 - 135      |     |      |   | 01/10/14 11:33  | 01/20/14 16:28  | 20             |
| 13C-1,2,3,7,8-PeCDF     | 80               |                  | 40 - 135      |     |      |   | 01/10/14 11:33  | 01/20/14 16:28  | 20             |
| 13C-1,2,3,6,7,8-HxCDD   | 62               |                  | 40 - 135      |     |      |   | 01/10/14 11:33  | 01/20/14 16:28  | 20             |
| 13C-1,2,3,4,7,8-HxCDF   | 58               |                  | 40 - 135      |     |      |   | 01/10/14 11:33  | 01/20/14 16:28  | 20             |
| 13C-1,2,3,4,6,7,8-HpCDD | 49               |                  | 40 - 135      |     |      |   | 01/10/14 11:33  | 01/20/14 16:28  | 20             |
| 13C-1,2,3,4,6,7,8-HpCDF | 54               |                  | 40 - 135      |     |      |   | 01/10/14 11:33  | 01/20/14 16:28  | 20             |
| 13C-OCDD                | 50               |                  | 40 - 135      |     |      |   | 01/10/14 11:33  | 01/20/14 16:28  | 20             |

**Method: 8290 - Dioxins and Furans (HRGC/HRMS) - RA**

| Analyte                 | Result           | Qualifier        | RL            | EDL | Unit | D | Prepared        | Analyzed        | Dil Fac        |
|-------------------------|------------------|------------------|---------------|-----|------|---|-----------------|-----------------|----------------|
| <b>2,3,7,8-TCDF</b>     | <b>1400</b>      |                  | 21            | 3.2 | pg/g | ☼ | 01/10/14 11:33  | 01/17/14 20:40  | 20             |
| <i>Isotope Dilution</i> | <i>%Recovery</i> | <i>Qualifier</i> | <i>Limits</i> |     |      |   | <i>Prepared</i> | <i>Analyzed</i> | <i>Dil Fac</i> |
| 13C-2,3,7,8-TCDF        | 69               |                  | 40 - 135      |     |      |   | 01/10/14 11:33  | 01/17/14 20:40  | 20             |

**Method: 6010B - Metals (ICP)**

| Analyte         | Result      | Qualifier | RL  | MDL | Unit  | D | Prepared       | Analyzed       | Dil Fac |
|-----------------|-------------|-----------|-----|-----|-------|---|----------------|----------------|---------|
| <b>Arsenic</b>  | <b>43</b>   |           | 6.2 |     | mg/Kg | ☼ | 12/30/13 12:18 | 12/31/13 11:27 | 10      |
| <b>Barium</b>   | <b>940</b>  |           | 1.0 |     | mg/Kg | ☼ | 12/30/13 12:18 | 12/31/13 11:27 | 10      |
| <b>Cadmium</b>  | <b>26</b>   |           | 2.1 |     | mg/Kg | ☼ | 12/30/13 12:18 | 12/31/13 11:27 | 10      |
| <b>Chromium</b> | <b>210</b>  |           | 2.7 |     | mg/Kg | ☼ | 12/30/13 12:18 | 12/31/13 11:27 | 10      |
| <b>Lead</b>     | <b>5300</b> |           | 3.1 |     | mg/Kg | ☼ | 12/30/13 12:18 | 12/31/13 11:27 | 10      |
| Selenium        | ND          |           | 10  |     | mg/Kg | ☼ | 12/30/13 12:18 | 12/31/13 11:27 | 10      |
| <b>Silver</b>   | <b>14</b>   |           | 5.2 |     | mg/Kg | ☼ | 12/30/13 12:18 | 12/31/13 11:27 | 10      |

**Method: 7471A - Mercury (CVAA)**

| Analyte        | Result     | Qualifier | RL    | MDL | Unit  | D | Prepared       | Analyzed       | Dil Fac |
|----------------|------------|-----------|-------|-----|-------|---|----------------|----------------|---------|
| <b>Mercury</b> | <b>1.1</b> |           | 0.020 |     | mg/Kg | ☼ | 12/31/13 08:25 | 12/31/13 10:04 | 10      |

TestAmerica Honolulu

# Surrogate Summary

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Method: 8270C - Semivolatile Organic Compounds (GC/MS)

Matrix: Solid

Prep Type: Total/NA

| Lab Sample ID      | Client Sample ID       | Percent Surrogate Recovery (Acceptance Limits) |                 |                 |                 |                 |                 |
|--------------------|------------------------|--|-----------------|-----------------|-----------------|-----------------|-----------------|
|                    |                        | TBP<br>(35-125)                                | 2FP<br>(25-120) | NBZ<br>(30-120) | PHL<br>(35-120) | TPH<br>(40-135) | FBP<br>(35-120) |
| LCS 440-153305/2-A | Lab Control Sample     | 86   | 83              | 74              | 74              | 99              | 80              |
| LCS 440-153305/3-A | Lab Control Sample Dup | 88   | 86              | 77              | 76              | 94              | 92              |
| MB 440-153305/1-A  | Method Blank           | 89   | 87              | 76              | 75              | 99              | 90              |

### Surrogate Legend

TBP = 2,4,6-Tribromophenol (Surr)

2FP = 2-Fluorophenol (Surr)

NBZ = Nitrobenzene-d5 (Surr)

PHL = Phenol-d6 (Surr)

TPH = Terphenyl-d14 (Surr)

FBP = 2-Fluorobiphenyl

## Method: 8270C - Semivolatile Organic Compounds (GC/MS)

Matrix: Solid/Soil

Prep Type: Total/NA

| Lab Sample ID | Client Sample ID | Percent Surrogate Recovery (Acceptance Limits) |                 |                 |                 |                 |                 |
|---------------|------------------|--|-----------------|-----------------|-----------------|-----------------|-----------------|
|               |                  | TBP<br>(35-125)                                | 2FP<br>(25-120) | NBZ<br>(30-120) | PHL<br>(35-120) | TPH<br>(40-135) | FBP<br>(35-120) |
| HWL0089-01    | RHS-01           | 40   | 41              | 52              | 45              | 79              | 69              |

### Surrogate Legend

TBP = 2,4,6-Tribromophenol (Surr)

2FP = 2-Fluorophenol (Surr)

NBZ = Nitrobenzene-d5 (Surr)

PHL = Phenol-d6 (Surr)

TPH = Terphenyl-d14 (Surr)

FBP = 2-Fluorobiphenyl

## Method: 8015B - Diesel Range Organics (DRO) (GC)

Matrix: Solid

Prep Type: Total/NA

| Lab Sample ID       | Client Sample ID       | Percent Surrogate Recovery (Acceptance Limits) |
|---------------------|------------------------|--|
|                     |                        | OTC1<br>(40-140)                               |
| 440-65826-A-1-B MS  | Matrix Spike           | 78   |
| 440-65826-A-1-C MSD | Matrix Spike Duplicate | 76   |
| LCS 440-153196/2-A  | Lab Control Sample     | 72   |
| MB 440-153196/1-A   | Method Blank           | 73   |

### Surrogate Legend

OTC = n-Octacosane

## Method: 8015B - Diesel Range Organics (DRO) (GC)

Matrix: Solid/Soil

Prep Type: Total/NA

| Lab Sample ID | Client Sample ID | Percent Surrogate Recovery (Acceptance Limits) |
|---------------|------------------|--|
|               |                  | OTC1<br>(40-140)                               |
| HWL0089-01    | RHS-01           | 63   |

### Surrogate Legend

TestAmerica Honolulu

# Surrogate Summary

Client: Bureau Veritas  
Project/Site: 17012-012148.00  
OTC = n-Octacosane

TestAmerica Job ID: HWL0089

## Method: 8081A - Organochlorine Pesticides (GC)

Matrix: Solid

Prep Type: Total/NA

### Percent Surrogate Recovery (Acceptance Limits)

| Lab Sample ID      | Client Sample ID       | TCX2<br>(35-115) | DCB2<br>(45-120) |
|--------------------|------------------------|------------------|------------------|
| LCS 440-153161/2-A | Lab Control Sample     | 78               | 96               |
| LCS 440-153161/3-A | Lab Control Sample Dup | 80               | 98               |
| MB 440-153161/1-A  | Method Blank           | 72               | 96               |

#### Surrogate Legend

TCX = Tetrachloro-m-xylene  
DCB = DCB Decachlorobiphenyl (Surr)

## Method: 8081A - Organochlorine Pesticides (GC)

Matrix: Solid/Soil

Prep Type: Total/NA

### Percent Surrogate Recovery (Acceptance Limits)

| Lab Sample ID | Client Sample ID | TCX2<br>(35-115) | DCB2<br>(45-120) |
|---------------|------------------|------------------|------------------|
| HWL0089-01    | RHS-01           | 67               | 96               |

#### Surrogate Legend

TCX = Tetrachloro-m-xylene  
DCB = DCB Decachlorobiphenyl (Surr)

## Method: 8082 - Polychlorinated Biphenyls (PCBs) by Gas Chromatography

Matrix: Solid

Prep Type: Total/NA

### Percent Surrogate Recovery (Acceptance Limits)

| Lab Sample ID       | Client Sample ID       | DCB1<br>(45-120) |
|---------------------|------------------------|------------------|
| 440-66257-E-2-A MS  | Matrix Spike           | 75 p             |
| 440-66257-E-2-B MSD | Matrix Spike Duplicate | 74 p             |
| LCS 440-153161/4-A  | Lab Control Sample     | 94               |
| MB 440-152687/1-A   | Method Blank           | 97               |

#### Surrogate Legend

DCB = DCB Decachlorobiphenyl (Surr)

## Method: 8082 - Polychlorinated Biphenyls (PCBs) by Gas Chromatography

Matrix: Solid/Soil

Prep Type: Total/NA

### Percent Surrogate Recovery (Acceptance Limits)

| Lab Sample ID | Client Sample ID | DCB1<br>(45-120) |
|---------------|------------------|------------------|
| HWL0089-01    | RHS-01           | 78               |

#### Surrogate Legend

DCB = DCB Decachlorobiphenyl (Surr)

TestAmerica Honolulu

# QC Sample Results

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Method: 8270C - Semivolatile Organic Compounds (GC/MS)

**Lab Sample ID: MB 440-153305/1-A**

**Matrix: Solid**

**Analysis Batch: 153775**

**Client Sample ID: Method Blank**

**Prep Type: Total/NA**

**Prep Batch: 153305**

| Analyte                         | MB Result | MB Qualifier | RL   | MDL   | Unit  | D | Prepared       | Analyzed       | Dil Fac |
|---------------------------------|-----------|--------------|------|-------|-------|---|----------------|----------------|---------|
| 1,2,4-Trichlorobenzene          | ND        |              | 0.33 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 1,2-Dichlorobenzene             | ND        |              | 0.33 | 0.060 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 1,3-Dichlorobenzene             | ND        |              | 0.33 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 1,4-Dichlorobenzene             | ND        |              | 0.33 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 1-Methylnaphthalene             | ND        |              | 0.35 | 0.15  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 2,4,5-Trichlorophenol           | ND        |              | 0.33 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 2,4,6-Trichlorophenol           | ND        |              | 0.33 | 0.075 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 2,4-Dichlorophenol              | ND        |              | 0.33 | 0.067 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 2,4-Dimethylphenol              | ND        |              | 0.33 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 2,4-Dinitrophenol               | ND        |              | 0.66 | 0.33  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 2,4-Dinitrotoluene              | ND        |              | 0.33 | 0.080 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 2,6-Dinitrotoluene              | ND        |              | 0.33 | 0.095 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 2-Chloronaphthalene             | ND        |              | 0.33 | 0.067 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 2-Chlorophenol                  | ND        |              | 0.33 | 0.070 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 2-Methylnaphthalene             | ND        |              | 0.33 | 0.070 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 2-Methylphenol                  | ND        |              | 0.33 | 0.080 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 2-Nitroaniline                  | ND        |              | 0.33 | 0.067 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 2-Nitrophenol                   | ND        |              | 0.33 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 3,3'-Dichlorobenzidine          | ND        |              | 0.83 | 0.15  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 3-Methylphenol + 4-Methylphenol | ND        |              | 0.33 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 3-Nitroaniline                  | ND        |              | 0.33 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 4,6-Dinitro-2-methylphenol      | ND        |              | 0.42 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 4-Bromophenyl phenyl ether      | ND        |              | 0.33 | 0.075 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 4-Chloro-3-methylphenol         | ND        |              | 0.33 | 0.070 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 4-Chloroaniline                 | ND        |              | 0.33 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 4-Chlorophenyl phenyl ether     | ND        |              | 0.33 | 0.085 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 4-Nitroaniline                  | ND        |              | 0.83 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 4-Nitrophenol                   | ND        |              | 0.83 | 0.14  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Acenaphthene                    | ND        |              | 0.33 | 0.067 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Acenaphthylene                  | ND        |              | 0.33 | 0.070 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Anthracene                      | ND        |              | 0.33 | 0.080 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Benzo[a]anthracene              | ND        |              | 0.33 | 0.070 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Benzo[a]pyrene                  | ND        |              | 0.33 | 0.067 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Benzo[b]fluoranthene            | ND        |              | 0.33 | 0.067 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Benzo[g,h,i]perylene            | ND        |              | 0.33 | 0.11  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Benzo[k]fluoranthene            | ND        |              | 0.33 | 0.070 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Benzoic acid                    | ND        |              | 0.83 | 0.15  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Benzyl alcohol                  | ND        |              | 0.33 | 0.20  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| bis (2-chloroisopropyl) ether   | ND        |              | 0.33 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Bis(2-chloroethoxy)methane      | ND        |              | 0.33 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Bis(2-chloroethyl)ether         | ND        |              | 0.33 | 0.060 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Bis(2-ethylhexyl) phthalate     | ND        |              | 0.33 | 0.090 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Butyl benzyl phthalate          | ND        |              | 0.33 | 0.080 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Carbazole                       | ND        |              | 0.33 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Chrysene                        | ND        |              | 0.33 | 0.075 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Dibenz(a,h)anthracene           | ND        |              | 0.42 | 0.10  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Dibenzofuran                    | ND        |              | 0.33 | 0.067 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Diethyl phthalate               | ND        |              | 0.33 | 0.095 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |

TestAmerica Honolulu

# QC Sample Results

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Method: 8270C - Semivolatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: MB 440-153305/1-A**

**Matrix: Solid**

**Analysis Batch: 153775**

**Client Sample ID: Method Blank**

**Prep Type: Total/NA**

**Prep Batch: 153305**

| Analyte                   | MB Result | MB Qualifier | RL   | MDL   | Unit  | D | Prepared       | Analyzed       | Dil Fac |
|---------------------------|-----------|--------------|------|-------|-------|---|----------------|----------------|---------|
| Dimethyl phthalate        | ND        |              | 0.33 | 0.067 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Di-n-butyl phthalate      | ND        |              | 0.33 | 0.090 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Di-n-octyl phthalate      | ND        |              | 0.33 | 0.090 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Fluoranthene              | ND        |              | 0.33 | 0.070 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Fluorene                  | ND        |              | 0.33 | 0.070 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Hexachlorobenzene         | ND        |              | 0.33 | 0.070 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Hexachlorobutadiene       | ND        |              | 0.33 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Hexachlorocyclopentadiene | ND        |              | 0.83 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Hexachloroethane          | ND        |              | 0.33 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Indeno[1,2,3-cd]pyrene    | ND        |              | 0.33 | 0.13  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Isophorone                | ND        |              | 0.33 | 0.067 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Naphthalene               | ND        |              | 0.33 | 0.067 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Nitrobenzene              | ND        |              | 0.33 | 0.070 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| N-Nitrosodi-n-propylamine | ND        |              | 0.25 | 0.070 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| N-Nitrosodiphenylamine    | ND        |              | 0.33 | 0.080 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Pentachlorophenol         | 0.191     | J            | 0.83 | 0.15  | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Phenanthrene              | ND        |              | 0.33 | 0.067 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Phenol                    | ND        |              | 0.33 | 0.090 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Pyrene                    | ND        |              | 0.33 | 0.080 | mg/Kg |   | 12/30/13 08:15 | 01/02/14 13:01 | 1       |

| Surrogate                   | MB %Recovery | MB Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|-----------------------------|--------------|--------------|----------|----------------|----------------|---------|
| 2,4,6-Tribromophenol (Surr) | 89           |              | 35 - 125 | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 2-Fluorophenol (Surr)       | 87           |              | 25 - 120 | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Nitrobenzene-d5 (Surr)      | 76           |              | 30 - 120 | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Phenol-d6 (Surr)            | 75           |              | 35 - 120 | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| Terphenyl-d14 (Surr)        | 99           |              | 40 - 135 | 12/30/13 08:15 | 01/02/14 13:01 | 1       |
| 2-Fluorobiphenyl            | 90           |              | 35 - 120 | 12/30/13 08:15 | 01/02/14 13:01 | 1       |

**Lab Sample ID: LCS 440-153305/2-A**

**Matrix: Solid**

**Analysis Batch: 153775**

**Client Sample ID: Lab Control Sample**

**Prep Type: Total/NA**

**Prep Batch: 153305**

| Analyte                | Spike Added | LCS Result | LCS Qualifier | Unit  | D | %Rec | %Rec. Limits |
|------------------------|-------------|------------|---------------|-------|---|------|--------------|
| 1,2,4-Trichlorobenzene | 3.33        | 2.70       |               | mg/Kg |   | 81   | 40 - 120     |
| 1,2-Dichlorobenzene    | 3.33        | 2.56       |               | mg/Kg |   | 77   | 40 - 120     |
| 1,3-Dichlorobenzene    | 3.33        | 2.45       |               | mg/Kg |   | 74   | 35 - 120     |
| 1,4-Dichlorobenzene    | 3.33        | 2.53       |               | mg/Kg |   | 76   | 35 - 120     |
| 2,4,5-Trichlorophenol  | 3.33        | 3.37       |               | mg/Kg |   | 101  | 50 - 120     |
| 2,4,6-Trichlorophenol  | 3.33        | 3.26       |               | mg/Kg |   | 98   | 50 - 120     |
| 2,4-Dichlorophenol     | 3.33        | 3.20       |               | mg/Kg |   | 96   | 45 - 120     |
| 2,4-Dimethylphenol     | 3.33        | 2.82       |               | mg/Kg |   | 85   | 40 - 120     |
| 2,4-Dinitrophenol      | 3.33        | 3.23       |               | mg/Kg |   | 97   | 25 - 120     |
| 2,4-Dinitrotoluene     | 3.33        | 3.41       |               | mg/Kg |   | 102  | 55 - 125     |
| 2,6-Dinitrotoluene     | 3.33        | 3.41       |               | mg/Kg |   | 102  | 55 - 125     |
| 2-Chloronaphthalene    | 3.33        | 2.77       |               | mg/Kg |   | 83   | 45 - 120     |
| 2-Chlorophenol         | 3.33        | 2.86       |               | mg/Kg |   | 86   | 40 - 120     |
| 2-Methylnaphthalene    | 3.33        | 2.86       |               | mg/Kg |   | 86   | 45 - 120     |
| 2-Methylphenol         | 3.33        | 2.85       |               | mg/Kg |   | 85   | 40 - 120     |

TestAmerica Honolulu



# QC Sample Results

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Method: 8270C - Semivolatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: LCS 440-153305/2-A**

**Matrix: Solid**

**Analysis Batch: 153775**

**Client Sample ID: Lab Control Sample**

**Prep Type: Total/NA**

**Prep Batch: 153305**

| Analyte                         | Spike Added | LCS Result | LCS Qualifier | Unit  | D | %Rec | %Rec. Limits |
|---------------------------------|-------------|------------|---------------|-------|---|------|--------------|
| 2-Nitroaniline                  | 3.33        | 3.00       |               | mg/Kg |   | 90   | 50 - 125     |
| 2-Nitrophenol                   | 3.33        | 2.99       |               | mg/Kg |   | 90   | 45 - 120     |
| 3,3'-Dichlorobenzidine          | 3.33        | 2.52       |               | mg/Kg |   | 76   | 20 - 130     |
| 3-Methylphenol + 4-Methylphenol | 3.33        | 2.85       |               | mg/Kg |   | 85   | 50 - 120     |
| 3-Nitroaniline                  | 3.33        | 2.76       |               | mg/Kg |   | 83   | 35 - 120     |
| 4,6-Dinitro-2-methylphenol      | 3.33        | 3.12       |               | mg/Kg |   | 94   | 40 - 120     |
| 4-Bromophenyl phenyl ether      | 3.33        | 3.09       |               | mg/Kg |   | 93   | 45 - 120     |
| 4-Chloro-3-methylphenol         | 3.33        | 3.15       |               | mg/Kg |   | 95   | 50 - 125     |
| 4-Chloroaniline                 | 3.33        | 1.83       |               | mg/Kg |   | 55   | 20 - 120     |
| 4-Chlorophenyl phenyl ether     | 3.33        | 3.20       |               | mg/Kg |   | 96   | 55 - 120     |
| 4-Nitroaniline                  | 3.33        | 3.41       |               | mg/Kg |   | 102  | 45 - 125     |
| 4-Nitrophenol                   | 3.33        | 3.00       |               | mg/Kg |   | 90   | 40 - 125     |
| Acenaphthene                    | 3.33        | 2.97       |               | mg/Kg |   | 89   | 50 - 120     |
| Acenaphthylene                  | 3.33        | 3.22       |               | mg/Kg |   | 97   | 50 - 120     |
| Anthracene                      | 3.33        | 3.21       |               | mg/Kg |   | 96   | 55 - 120     |
| Benzo[a]anthracene              | 3.33        | 3.11       |               | mg/Kg |   | 93   | 55 - 120     |
| Benzo[a]pyrene                  | 3.33        | 3.05       |               | mg/Kg |   | 92   | 50 - 125     |
| Benzo[b]fluoranthene            | 3.33        | 2.81       |               | mg/Kg |   | 84   | 45 - 125     |
| Benzo[g,h,i]perylene            | 3.33        | 3.40       |               | mg/Kg |   | 102  | 35 - 130     |
| Benzo[k]fluoranthene            | 3.33        | 3.12       |               | mg/Kg |   | 93   | 45 - 125     |
| Benzoic acid                    | 3.33        | 3.07       |               | mg/Kg |   | 92   | 20 - 120     |
| Benzyl alcohol                  | 3.33        | 2.48       |               | mg/Kg |   | 75   | 35 - 120     |
| bis (2-chloroisopropyl) ether   | 3.33        | 2.08       |               | mg/Kg |   | 62   | 40 - 120     |
| Bis(2-chloroethoxy)methane      | 3.33        | 2.70       |               | mg/Kg |   | 81   | 45 - 120     |
| Bis(2-chloroethyl)ether         | 3.33        | 2.54       |               | mg/Kg |   | 76   | 35 - 120     |
| Bis(2-ethylhexyl) phthalate     | 3.33        | 3.58       |               | mg/Kg |   | 107  | 50 - 130     |
| Butyl benzyl phthalate          | 3.33        | 3.30       |               | mg/Kg |   | 99   | 50 - 125     |
| Chrysene                        | 3.33        | 3.13       |               | mg/Kg |   | 94   | 55 - 120     |
| Dibenz(a,h)anthracene           | 3.33        | 3.43       |               | mg/Kg |   | 103  | 40 - 135     |
| Dibenzofuran                    | 3.33        | 3.04       |               | mg/Kg |   | 91   | 55 - 120     |
| Diethyl phthalate               | 3.33        | 3.37       |               | mg/Kg |   | 101  | 50 - 125     |
| Dimethyl phthalate              | 3.33        | 3.24       |               | mg/Kg |   | 97   | 50 - 125     |
| Di-n-butyl phthalate            | 3.33        | 3.41       |               | mg/Kg |   | 102  | 50 - 125     |
| Di-n-octyl phthalate            | 3.33        | 3.55       |               | mg/Kg |   | 107  | 50 - 135     |
| Fluoranthene                    | 3.33        | 3.23       |               | mg/Kg |   | 97   | 55 - 120     |
| Fluorene                        | 3.33        | 3.21       |               | mg/Kg |   | 96   | 55 - 120     |
| Hexachlorobenzene               | 3.33        | 2.96       |               | mg/Kg |   | 89   | 50 - 120     |
| Hexachlorobutadiene             | 3.33        | 2.70       |               | mg/Kg |   | 81   | 40 - 120     |
| Hexachlorocyclopentadiene       | 3.33        | 2.78       |               | mg/Kg |   | 83   | 30 - 125     |
| Hexachloroethane                | 3.33        | 2.46       |               | mg/Kg |   | 74   | 40 - 120     |
| Indeno[1,2,3-cd]pyrene          | 3.33        | 3.74       |               | mg/Kg |   | 112  | 30 - 135     |
| Isophorone                      | 3.33        | 2.73       |               | mg/Kg |   | 82   | 40 - 120     |
| Naphthalene                     | 3.33        | 2.68       |               | mg/Kg |   | 81   | 45 - 120     |
| Nitrobenzene                    | 3.33        | 2.55       |               | mg/Kg |   | 77   | 45 - 120     |
| N-Nitrosodi-n-propylamine       | 3.33        | 2.28       |               | mg/Kg |   | 68   | 40 - 120     |
| N-Nitrosodiphenylamine          | 3.33        | 2.85       |               | mg/Kg |   | 86   | 50 - 120     |
| Pentachlorophenol               | 3.33        | 3.19       |               | mg/Kg |   | 96   | 40 - 120     |
| Phenanthrene                    | 3.33        | 3.08       |               | mg/Kg |   | 92   | 50 - 120     |

TestAmerica Honolulu



# QC Sample Results

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Method: 8270C - Semivolatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: LCS 440-153305/2-A**

**Matrix: Solid**

**Analysis Batch: 153775**

**Client Sample ID: Lab Control Sample**

**Prep Type: Total/NA**

**Prep Batch: 153305**

| Analyte | Spike Added | LCS Result | LCS Qualifier | Unit  | D | %Rec | %Rec. Limits |
|---------|-------------|------------|---------------|-------|---|------|--------------|
| Phenol  | 3.33        | 2.83       |               | mg/Kg |   | 85   | 40 - 120     |
| Pyrene  | 3.33        | 3.33       |               | mg/Kg |   | 100  | 45 - 125     |

| Surrogate                   | LCS %Recovery | LCS Qualifier | Limits   |
|-----------------------------|---------------|---------------|----------|
| 2,4,6-Tribromophenol (Surr) | 86            |               | 35 - 125 |
| 2-Fluorophenol (Surr)       | 83            |               | 25 - 120 |
| Nitrobenzene-d5 (Surr)      | 74            |               | 30 - 120 |
| Phenol-d6 (Surr)            | 74            |               | 35 - 120 |
| Terphenyl-d14 (Surr)        | 99            |               | 40 - 135 |
| 2-Fluorobiphenyl            | 80            |               | 35 - 120 |

**Lab Sample ID: LCSD 440-153305/3-A**

**Matrix: Solid**

**Analysis Batch: 153775**

**Client Sample ID: Lab Control Sample Dup**

**Prep Type: Total/NA**

**Prep Batch: 153305**

| Analyte                         | Spike Added | LCSD Result | LCSD Qualifier | Unit  | D | %Rec | %Rec. Limits | RPD | RPD Limit |
|---------------------------------|-------------|-------------|----------------|-------|---|------|--------------|-----|-----------|
| 1,2,4-Trichlorobenzene          | 3.33        | 2.73        |                | mg/Kg |   | 82   | 40 - 120     | 1   | 20        |
| 1,2-Dichlorobenzene             | 3.33        | 2.62        |                | mg/Kg |   | 79   | 40 - 120     | 2   | 20        |
| 1,3-Dichlorobenzene             | 3.33        | 2.49        |                | mg/Kg |   | 75   | 35 - 120     | 2   | 25        |
| 1,4-Dichlorobenzene             | 3.33        | 2.52        |                | mg/Kg |   | 75   | 35 - 120     | 1   | 25        |
| 2,4,5-Trichlorophenol           | 3.33        | 3.43        |                | mg/Kg |   | 103  | 50 - 120     | 2   | 20        |
| 2,4,6-Trichlorophenol           | 3.33        | 3.54        |                | mg/Kg |   | 106  | 50 - 120     | 8   | 20        |
| 2,4-Dichlorophenol              | 3.33        | 3.23        |                | mg/Kg |   | 97   | 45 - 120     | 1   | 20        |
| 2,4-Dimethylphenol              | 3.33        | 2.82        |                | mg/Kg |   | 84   | 40 - 120     | 0   | 20        |
| 2,4-Dinitrophenol               | 3.33        | 2.85        |                | mg/Kg |   | 85   | 25 - 120     | 13  | 25        |
| 2,4-Dinitrotoluene              | 3.33        | 2.62        | *              | mg/Kg |   | 79   | 55 - 125     | 26  | 20        |
| 2,6-Dinitrotoluene              | 3.33        | 2.98        |                | mg/Kg |   | 89   | 55 - 125     | 14  | 20        |
| 2-Chloronaphthalene             | 3.33        | 3.04        |                | mg/Kg |   | 91   | 45 - 120     | 9   | 20        |
| 2-Chlorophenol                  | 3.33        | 3.07        |                | mg/Kg |   | 92   | 40 - 120     | 7   | 20        |
| 2-Methylnaphthalene             | 3.33        | 2.70        |                | mg/Kg |   | 81   | 45 - 120     | 6   | 20        |
| 2-Methylphenol                  | 3.33        | 3.00        |                | mg/Kg |   | 90   | 40 - 120     | 5   | 20        |
| 2-Nitroaniline                  | 3.33        | 2.76        |                | mg/Kg |   | 83   | 50 - 125     | 8   | 20        |
| 2-Nitrophenol                   | 3.33        | 3.26        |                | mg/Kg |   | 98   | 45 - 120     | 9   | 20        |
| 3,3'-Dichlorobenzidine          | 3.33        | 2.29        |                | mg/Kg |   | 69   | 20 - 130     | 10  | 25        |
| 3-Methylphenol + 4-Methylphenol | 3.33        | 3.01        |                | mg/Kg |   | 90   | 50 - 120     | 5   | 20        |
| 3-Nitroaniline                  | 3.33        | 2.03        | *              | mg/Kg |   | 61   | 35 - 120     | 30  | 25        |
| 4,6-Dinitro-2-methylphenol      | 3.33        | 3.09        |                | mg/Kg |   | 93   | 40 - 120     | 1   | 25        |
| 4-Bromophenyl phenyl ether      | 3.33        | 3.33        |                | mg/Kg |   | 100  | 45 - 120     | 7   | 20        |
| 4-Chloro-3-methylphenol         | 3.33        | 2.77        |                | mg/Kg |   | 83   | 50 - 125     | 13  | 20        |
| 4-Chloroaniline                 | 3.33        | 1.52        |                | mg/Kg |   | 46   | 20 - 120     | 19  | 30        |
| 4-Chlorophenyl phenyl ether     | 3.33        | 2.84        |                | mg/Kg |   | 85   | 55 - 120     | 12  | 20        |
| 4-Nitroaniline                  | 3.33        | 2.41        | *              | mg/Kg |   | 72   | 45 - 125     | 34  | 20        |
| 4-Nitrophenol                   | 3.33        | 2.24        | *              | mg/Kg |   | 67   | 40 - 125     | 29  | 20        |
| Acenaphthene                    | 3.33        | 2.89        |                | mg/Kg |   | 87   | 50 - 120     | 2   | 20        |
| Acenaphthylene                  | 3.33        | 3.21        |                | mg/Kg |   | 96   | 50 - 120     | 0   | 20        |
| Anthracene                      | 3.33        | 3.11        |                | mg/Kg |   | 93   | 55 - 120     | 3   | 20        |
| Benzo[a]anthracene              | 3.33        | 3.04        |                | mg/Kg |   | 91   | 55 - 120     | 2   | 20        |
| Benzo[a]pyrene                  | 3.33        | 2.99        |                | mg/Kg |   | 90   | 50 - 125     | 2   | 20        |

TestAmerica Honolulu

# QC Sample Results

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Method: 8270C - Semivolatile Organic Compounds (GC/MS) (Continued)

Lab Sample ID: LCSD 440-153305/3-A

Matrix: Solid

Analysis Batch: 153775

Client Sample ID: Lab Control Sample Dup

Prep Type: Total/NA

Prep Batch: 153305

| Analyte                       | Spike Added | LCSD Result | LCSD Qualifier | Unit  | D | %Rec | %Rec.    |     | RPD | Limit |
|-------------------------------|-------------|-------------|----------------|-------|---|------|----------|-----|-----|-------|
|                               |             |             |                |       |   |      | Limits   | RPD |     |       |
| Benzo[b]fluoranthene          | 3.33        | 2.74        |                | mg/Kg |   | 82   | 45 - 125 | 3   | 25  |       |
| Benzo[g,h,i]perylene          | 3.33        | 3.39        |                | mg/Kg |   | 102  | 35 - 130 | 0   | 25  |       |
| Benzo[k]fluoranthene          | 3.33        | 3.09        |                | mg/Kg |   | 93   | 45 - 125 | 1   | 25  |       |
| Benzoic acid                  | 3.33        | 3.32        |                | mg/Kg |   | 100  | 20 - 120 | 8   | 30  |       |
| Benzyl alcohol                | 3.33        | 2.70        |                | mg/Kg |   | 81   | 35 - 120 | 8   | 25  |       |
| bis (2-chloroisopropyl) ether | 3.33        | 2.07        |                | mg/Kg |   | 62   | 40 - 120 | 0   | 20  |       |
| Bis(2-chloroethoxy)methane    | 3.33        | 2.78        |                | mg/Kg |   | 84   | 45 - 120 | 3   | 20  |       |
| Bis(2-chloroethyl)ether       | 3.33        | 2.59        |                | mg/Kg |   | 78   | 35 - 120 | 2   | 25  |       |
| Bis(2-ethylhexyl) phthalate   | 3.33        | 3.56        |                | mg/Kg |   | 107  | 50 - 130 | 1   | 20  |       |
| Butyl benzyl phthalate        | 3.33        | 3.20        |                | mg/Kg |   | 96   | 50 - 125 | 3   | 20  |       |
| Chrysene                      | 3.33        | 3.08        |                | mg/Kg |   | 92   | 55 - 120 | 2   | 20  |       |
| Dibenz(a,h)anthracene         | 3.33        | 3.42        |                | mg/Kg |   | 103  | 40 - 135 | 0   | 25  |       |
| Dibenzofuran                  | 3.33        | 2.85        |                | mg/Kg |   | 85   | 55 - 120 | 7   | 20  |       |
| Diethyl phthalate             | 3.33        | 2.72        | *              | mg/Kg |   | 81   | 50 - 125 | 21  | 20  |       |
| Dimethyl phthalate            | 3.33        | 2.92        |                | mg/Kg |   | 88   | 50 - 125 | 10  | 20  |       |
| Di-n-butyl phthalate          | 3.33        | 3.12        |                | mg/Kg |   | 94   | 50 - 125 | 9   | 20  |       |
| Di-n-octyl phthalate          | 3.33        | 3.52        |                | mg/Kg |   | 106  | 50 - 135 | 1   | 20  |       |
| Fluoranthene                  | 3.33        | 2.94        |                | mg/Kg |   | 88   | 55 - 120 | 9   | 20  |       |
| Fluorene                      | 3.33        | 2.74        |                | mg/Kg |   | 82   | 55 - 120 | 16  | 20  |       |
| Hexachlorobenzene             | 3.33        | 3.08        |                | mg/Kg |   | 92   | 50 - 120 | 4   | 20  |       |
| Hexachlorobutadiene           | 3.33        | 2.68        |                | mg/Kg |   | 80   | 40 - 120 | 1   | 20  |       |
| Hexachlorocyclopentadiene     | 3.33        | 3.55        |                | mg/Kg |   | 106  | 30 - 125 | 24  | 25  |       |
| Hexachloroethane              | 3.33        | 2.48        |                | mg/Kg |   | 74   | 40 - 120 | 1   | 20  |       |
| Indeno[1,2,3-cd]pyrene        | 3.33        | 3.72        |                | mg/Kg |   | 112  | 30 - 135 | 0   | 25  |       |
| Isophorone                    | 3.33        | 2.71        |                | mg/Kg |   | 81   | 40 - 120 | 0   | 20  |       |
| Naphthalene                   | 3.33        | 2.71        |                | mg/Kg |   | 81   | 45 - 120 | 1   | 20  |       |
| Nitrobenzene                  | 3.33        | 2.62        |                | mg/Kg |   | 79   | 45 - 120 | 3   | 20  |       |
| N-Nitrosodi-n-propylamine     | 3.33        | 2.43        |                | mg/Kg |   | 73   | 40 - 120 | 6   | 20  |       |
| N-Nitrosodiphenylamine        | 3.33        | 3.12        |                | mg/Kg |   | 94   | 50 - 120 | 9   | 20  |       |
| Pentachlorophenol             | 3.33        | 3.16        |                | mg/Kg |   | 95   | 40 - 120 | 1   | 20  |       |
| Phenanthrene                  | 3.33        | 3.00        |                | mg/Kg |   | 90   | 50 - 120 | 3   | 20  |       |
| Phenol                        | 3.33        | 3.02        |                | mg/Kg |   | 91   | 40 - 120 | 7   | 20  |       |
| Pyrene                        | 3.33        | 3.23        |                | mg/Kg |   | 97   | 45 - 125 | 3   | 25  |       |

| Surrogate                   | LCSD      |           | Limits   |
|-----------------------------|-----------|-----------|----------|
|                             | %Recovery | Qualifier |          |
| 2,4,6-Tribromophenol (Surr) | 88        |           | 35 - 125 |
| 2-Fluorophenol (Surr)       | 86        |           | 25 - 120 |
| Nitrobenzene-d5 (Surr)      | 77        |           | 30 - 120 |
| Phenol-d6 (Surr)            | 76        |           | 35 - 120 |
| Terphenyl-d14 (Surr)        | 94        |           | 40 - 135 |
| 2-Fluorobiphenyl            | 92        |           | 35 - 120 |

TestAmerica Honolulu

# QC Sample Results

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Method: 8015B - Diesel Range Organics (DRO) (GC)

**Lab Sample ID: MB 440-153196/1-A**

**Matrix: Solid**

**Analysis Batch: 153341**

**Client Sample ID: Method Blank**

**Prep Type: Total/NA**

**Prep Batch: 153196**

| Analyte       | MB Result | MB Qualifier | RL  | MDL | Unit  | D | Prepared       | Analyzed       | Dil Fac |
|---------------|-----------|--------------|-----|-----|-------|---|----------------|----------------|---------|
| DRO (C10-C28) | ND        |              | 5.0 | 2.5 | mg/Kg |   | 12/28/13 07:03 | 12/30/13 08:15 | 1       |
| RRO(C29-C40)  | ND        |              | 5.0 | 2.5 | mg/Kg |   | 12/28/13 07:03 | 12/30/13 08:15 | 1       |

| Surrogate    | MB %Recovery | MB Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|--------------|--------------|--------------|----------|----------------|----------------|---------|
| n-Octacosane | 73           |              | 40 - 140 | 12/28/13 07:03 | 12/30/13 08:15 | 1       |

**Lab Sample ID: LCS 440-153196/2-A**

**Matrix: Solid**

**Analysis Batch: 153341**

**Client Sample ID: Lab Control Sample**

**Prep Type: Total/NA**

**Prep Batch: 153196**

| Analyte       | Spike Added | LCS Result | LCS Qualifier | Unit  | D | %Rec | Limits   |
|---------------|-------------|------------|---------------|-------|---|------|----------|
| DRO (C10-C28) | 33.3        | 20.9       |               | mg/Kg |   | 63   | 45 - 115 |

| Surrogate    | LCS %Recovery | LCS Qualifier | Limits   |
|--------------|---------------|---------------|----------|
| n-Octacosane | 72            |               | 40 - 140 |

**Lab Sample ID: 440-65826-A-1-B MS**

**Matrix: Solid**

**Analysis Batch: 153341**

**Client Sample ID: Matrix Spike**

**Prep Type: Total/NA**

**Prep Batch: 153196**

| Analyte       | Sample Result | Sample Qualifier | Spike Added | MS Result | MS Qualifier | Unit  | D | %Rec | %Rec. Limits |
|---------------|---------------|------------------|-------------|-----------|--------------|-------|---|------|--------------|
| DRO (C10-C28) | ND            |                  | 33.4        | 22.1      |              | mg/Kg |   | 66   | 40 - 120     |

| Surrogate    | MS %Recovery | MS Qualifier | Limits   |
|--------------|--------------|--------------|----------|
| n-Octacosane | 78           |              | 40 - 140 |

**Lab Sample ID: 440-65826-A-1-C MSD**

**Matrix: Solid**

**Analysis Batch: 153341**

**Client Sample ID: Matrix Spike Duplicate**

**Prep Type: Total/NA**

**Prep Batch: 153196**

| Analyte       | Sample Result | Sample Qualifier | Spike Added | MSD Result | MSD Qualifier | Unit  | D | %Rec | %Rec. Limits | RPD | Limit |
|---------------|---------------|------------------|-------------|------------|---------------|-------|---|------|--------------|-----|-------|
| DRO (C10-C28) | ND            |                  | 33.4        | 21.8       |               | mg/Kg |   | 65   | 40 - 120     | 1   | 30    |

| Surrogate    | MSD %Recovery | MSD Qualifier | Limits   |
|--------------|---------------|---------------|----------|
| n-Octacosane | 76            |               | 40 - 140 |

## Method: 8081A - Organochlorine Pesticides (GC)

**Lab Sample ID: MB 440-153161/1-A**

**Matrix: Solid**

**Analysis Batch: 153523**

**Client Sample ID: Method Blank**

**Prep Type: Total/NA**

**Prep Batch: 153161**

| Analyte  | MB Result | MB Qualifier | RL     | MDL    | Unit  | D | Prepared       | Analyzed       | Dil Fac |
|----------|-----------|--------------|--------|--------|-------|---|----------------|----------------|---------|
| 4,4'-DDD | ND        |              | 0.0050 | 0.0015 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| 4,4'-DDE | ND        |              | 0.0050 | 0.0015 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| 4,4'-DDT | ND        |              | 0.0050 | 0.0015 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| Aldrin   | ND        |              | 0.0050 | 0.0015 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |

TestAmerica Honolulu

# QC Sample Results

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Method: 8081A - Organochlorine Pesticides (GC) (Continued)

**Lab Sample ID: MB 440-153161/1-A**  
**Matrix: Solid**  
**Analysis Batch: 153523**

**Client Sample ID: Method Blank**  
**Prep Type: Total/NA**  
**Prep Batch: 153161**

| Analyte               | MB Result | MB Qualifier | RL     | MDL    | Unit  | D | Prepared       | Analyzed       | Dil Fac |
|-----------------------|-----------|--------------|--------|--------|-------|---|----------------|----------------|---------|
| alpha-BHC             | ND        |              | 0.0050 | 0.0015 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| beta-BHC              | ND        |              | 0.0050 | 0.0015 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| Chlordane (technical) | ND        |              | 0.050  | 0.010  | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| delta-BHC             | ND        |              | 0.010  | 0.0015 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| Dieldrin              | ND        |              | 0.0050 | 0.0015 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| Endosulfan I          | ND        |              | 0.0050 | 0.0015 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| Endosulfan II         | ND        |              | 0.0050 | 0.0015 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| Endosulfan sulfate    | ND        |              | 0.010  | 0.0020 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| Endrin                | ND        |              | 0.0050 | 0.0015 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| Endrin aldehyde       | ND        |              | 0.0050 | 0.0015 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| Endrin ketone         | ND        |              | 0.0050 | 0.0020 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| gamma-BHC (Lindane)   | ND        |              | 0.0050 | 0.0015 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| Heptachlor            | ND        |              | 0.0050 | 0.0020 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| Heptachlor epoxide    | ND        |              | 0.0050 | 0.0020 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| Methoxychlor          | ND        |              | 0.0050 | 0.0015 | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| Toxaphene             | ND        |              | 0.20   | 0.050  | mg/Kg |   | 12/27/13 17:32 | 12/31/13 14:09 | 1       |

| Surrogate                     | MB %Recovery | MB Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|-------------------------------|--------------|--------------|----------|----------------|----------------|---------|
| Tetrachloro-m-xylene          | 72           |              | 35 - 115 | 12/27/13 17:32 | 12/31/13 14:09 | 1       |
| DCB Decachlorobiphenyl (Surr) | 96           |              | 45 - 120 | 12/27/13 17:32 | 12/31/13 14:09 | 1       |

**Lab Sample ID: LCS 440-153161/2-A**  
**Matrix: Solid**  
**Analysis Batch: 153523**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 153161**

| Analyte             | Spike Added | LCS Result | LCS Qualifier | Unit  | D | %Rec | %Rec. Limits |
|---------------------|-------------|------------|---------------|-------|---|------|--------------|
| 4,4'-DDD            | 0.0333      | 0.0344     |               | mg/Kg |   | 103  | 60 - 120     |
| 4,4'-DDE            | 0.0333      | 0.0318     |               | mg/Kg |   | 95   | 60 - 120     |
| 4,4'-DDT            | 0.0333      | 0.0298     |               | mg/Kg |   | 90   | 65 - 120     |
| Aldrin              | 0.0333      | 0.0283     |               | mg/Kg |   | 85   | 50 - 115     |
| alpha-BHC           | 0.0333      | 0.0288     |               | mg/Kg |   | 87   | 60 - 115     |
| beta-BHC            | 0.0333      | 0.0293     |               | mg/Kg |   | 88   | 60 - 115     |
| delta-BHC           | 0.0333      | 0.0317     |               | mg/Kg |   | 95   | 60 - 115     |
| Dieldrin            | 0.0333      | 0.0311     |               | mg/Kg |   | 93   | 65 - 115     |
| Endosulfan I        | 0.0333      | 0.0303     |               | mg/Kg |   | 91   | 40 - 120     |
| Endosulfan II       | 0.0333      | 0.0318     |               | mg/Kg |   | 95   | 55 - 120     |
| Endosulfan sulfate  | 0.0333      | 0.0332     |               | mg/Kg |   | 100  | 65 - 115     |
| Endrin              | 0.0333      | 0.0287     |               | mg/Kg |   | 86   | 55 - 120     |
| Endrin aldehyde     | 0.0333      | 0.0266     |               | mg/Kg |   | 80   | 55 - 115     |
| Endrin ketone       | 0.0333      | 0.0325     |               | mg/Kg |   | 97   | 65 - 115     |
| gamma-BHC (Lindane) | 0.0333      | 0.0291     |               | mg/Kg |   | 87   | 55 - 115     |
| Heptachlor          | 0.0333      | 0.0272     |               | mg/Kg |   | 82   | 55 - 115     |
| Heptachlor epoxide  | 0.0333      | 0.0297     |               | mg/Kg |   | 89   | 55 - 115     |
| Methoxychlor        | 0.0333      | 0.0292     |               | mg/Kg |   | 88   | 65 - 120     |

| Surrogate            | LCS %Recovery | LCS Qualifier | Limits   |
|----------------------|---------------|---------------|----------|
| Tetrachloro-m-xylene | 78            |               | 35 - 115 |

TestAmerica Honolulu

# QC Sample Results

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Method: 8081A - Organochlorine Pesticides (GC) (Continued)

**Lab Sample ID:** LCS 440-153161/2-A  
**Matrix:** Solid  
**Analysis Batch:** 153523

**Client Sample ID:** Lab Control Sample  
**Prep Type:** Total/NA  
**Prep Batch:** 153161

| Surrogate                     | LCS<br>%Recovery | LCS<br>Qualifier | Limits   |
|-------------------------------|------------------|------------------|----------|
| DCB Decachlorobiphenyl (Surr) | 96               |                  | 45 - 120 |

**Lab Sample ID:** LCSD 440-153161/3-A  
**Matrix:** Solid  
**Analysis Batch:** 153523

**Client Sample ID:** Lab Control Sample Dup  
**Prep Type:** Total/NA  
**Prep Batch:** 153161

| Analyte             | Spike<br>Added | LCSD<br>Result | LCSD<br>Qualifier | Unit  | D | %Rec | %Rec.<br>Limits | RPD | RPD<br>Limit |
|---------------------|----------------|----------------|-------------------|-------|---|------|-----------------|-----|--------------|
| 4,4'-DDD            | 0.0333         | 0.0352         |                   | mg/Kg |   | 106  | 60 - 120        | 2   | 30           |
| 4,4'-DDE            | 0.0333         | 0.0323         |                   | mg/Kg |   | 97   | 60 - 120        | 2   | 30           |
| 4,4'-DDT            | 0.0333         | 0.0298         |                   | mg/Kg |   | 89   | 65 - 120        | 0   | 30           |
| Aldrin              | 0.0333         | 0.0287         |                   | mg/Kg |   | 86   | 50 - 115        | 1   | 30           |
| alpha-BHC           | 0.0333         | 0.0292         |                   | mg/Kg |   | 88   | 60 - 115        | 1   | 30           |
| beta-BHC            | 0.0333         | 0.0299         |                   | mg/Kg |   | 90   | 60 - 115        | 2   | 30           |
| delta-BHC           | 0.0333         | 0.0324         |                   | mg/Kg |   | 97   | 60 - 115        | 2   | 30           |
| Dieldrin            | 0.0333         | 0.0316         |                   | mg/Kg |   | 95   | 65 - 115        | 2   | 30           |
| Endosulfan I        | 0.0333         | 0.0308         |                   | mg/Kg |   | 92   | 40 - 120        | 2   | 30           |
| Endosulfan II       | 0.0333         | 0.0325         |                   | mg/Kg |   | 98   | 55 - 120        | 2   | 30           |
| Endosulfan sulfate  | 0.0333         | 0.0343         |                   | mg/Kg |   | 103  | 65 - 115        | 3   | 30           |
| Endrin              | 0.0333         | 0.0287         |                   | mg/Kg |   | 86   | 55 - 120        | 0   | 30           |
| Endrin aldehyde     | 0.0333         | 0.0277         |                   | mg/Kg |   | 83   | 55 - 115        | 4   | 30           |
| Endrin ketone       | 0.0333         | 0.0338         |                   | mg/Kg |   | 102  | 65 - 115        | 4   | 30           |
| gamma-BHC (Lindane) | 0.0333         | 0.0295         |                   | mg/Kg |   | 89   | 55 - 115        | 1   | 30           |
| Heptachlor          | 0.0333         | 0.0274         |                   | mg/Kg |   | 82   | 55 - 115        | 1   | 30           |
| Heptachlor epoxide  | 0.0333         | 0.0300         |                   | mg/Kg |   | 90   | 55 - 115        | 1   | 30           |
| Methoxychlor        | 0.0333         | 0.0295         |                   | mg/Kg |   | 89   | 65 - 120        | 1   | 30           |

| Surrogate                     | LCSD<br>%Recovery | LCSD<br>Qualifier | Limits   |
|-------------------------------|-------------------|-------------------|----------|
| Tetrachloro-m-xylene          | 80                |                   | 35 - 115 |
| DCB Decachlorobiphenyl (Surr) | 98                |                   | 45 - 120 |

## Method: 8082 - Polychlorinated Biphenyls (PCBs) by Gas Chromatography

**Lab Sample ID:** MB 440-152687/1-A  
**Matrix:** Solid  
**Analysis Batch:** 153427

**Client Sample ID:** Method Blank  
**Prep Type:** Total/NA  
**Prep Batch:** 152687

| Analyte      | MB<br>Result | MB<br>Qualifier | RL    | MDL   | Unit  | D | Prepared       | Analyzed       | Dil Fac |
|--------------|--------------|-----------------|-------|-------|-------|---|----------------|----------------|---------|
| Aroclor 1016 | ND           |                 | 0.050 | 0.017 | mg/Kg |   | 12/24/13 18:18 | 12/30/13 19:07 | 1       |
| Aroclor 1221 | ND           |                 | 0.050 | 0.017 | mg/Kg |   | 12/24/13 18:18 | 12/30/13 19:07 | 1       |
| Aroclor 1232 | ND           |                 | 0.050 | 0.017 | mg/Kg |   | 12/24/13 18:18 | 12/30/13 19:07 | 1       |
| Aroclor 1242 | ND           |                 | 0.050 | 0.017 | mg/Kg |   | 12/24/13 18:18 | 12/30/13 19:07 | 1       |
| Aroclor 1248 | ND           |                 | 0.050 | 0.017 | mg/Kg |   | 12/24/13 18:18 | 12/30/13 19:07 | 1       |
| Aroclor 1254 | ND           |                 | 0.050 | 0.017 | mg/Kg |   | 12/24/13 18:18 | 12/30/13 19:07 | 1       |
| Aroclor 1260 | ND           |                 | 0.050 | 0.017 | mg/Kg |   | 12/24/13 18:18 | 12/30/13 19:07 | 1       |

| Surrogate                     | MB<br>%Recovery | MB<br>Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|-------------------------------|-----------------|-----------------|----------|----------------|----------------|---------|
| DCB Decachlorobiphenyl (Surr) | 97              |                 | 45 - 120 | 12/24/13 18:18 | 12/30/13 19:07 | 1       |

TestAmerica Honolulu



# QC Sample Results

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Method: 8082 - Polychlorinated Biphenyls (PCBs) by Gas Chromatography (Continued)

**Lab Sample ID: LCS 440-153161/4-A**  
**Matrix: Solid**  
**Analysis Batch: 153427**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 153161**

| Analyte                       | Spike Added | LCS    |           | Unit     | D | %Rec | %Rec. Limits |
|-------------------------------|-------------|--------|-----------|----------|---|------|--------------|
|                               |             | Result | Qualifier |          |   |      |              |
| Aroclor 1016                  | 0.267       | 0.195  |           | mg/Kg    |   | 73   | 65 - 115     |
| Aroclor 1260                  | 0.267       | 0.211  |           | mg/Kg    |   | 79   | 65 - 115     |
| Surrogate                     | %Recovery   |        | Qualifier | Limits   |   |      |              |
| DCB Decachlorobiphenyl (Surr) | 94          |        |           | 45 - 120 |   |      |              |

**Lab Sample ID: 440-66257-E-2-A MS**  
**Matrix: Solid**  
**Analysis Batch: 153427**

**Client Sample ID: Matrix Spike**  
**Prep Type: Total/NA**  
**Prep Batch: 153161**

| Analyte                       | Sample Result | Sample Qualifier | Spike Added | MS       |           | Unit  | D | %Rec | %Rec. Limits |
|-------------------------------|---------------|------------------|-------------|----------|-----------|-------|---|------|--------------|
|                               |               |                  |             | Result   | Qualifier |       |   |      |              |
| Aroclor 1016                  | ND            |                  | 0.264       | 0.190    |           | mg/Kg |   | 72   | 50 - 120     |
| Aroclor 1260                  | ND            |                  | 0.264       | 0.186    |           | mg/Kg |   | 70   | 50 - 125     |
| Surrogate                     | %Recovery     |                  | Qualifier   | Limits   |           |       |   |      |              |
| DCB Decachlorobiphenyl (Surr) | 75            |                  | p           | 45 - 120 |           |       |   |      |              |

**Lab Sample ID: 440-66257-E-2-B MSD**  
**Matrix: Solid**  
**Analysis Batch: 153427**

**Client Sample ID: Matrix Spike Duplicate**  
**Prep Type: Total/NA**  
**Prep Batch: 153161**

| Analyte                       | Sample Result | Sample Qualifier | Spike Added | MSD      |           | Unit  | D | %Rec | %Rec. Limits | RPD | Limit |
|-------------------------------|---------------|------------------|-------------|----------|-----------|-------|---|------|--------------|-----|-------|
|                               |               |                  |             | Result   | Qualifier |       |   |      |              |     |       |
| Aroclor 1016                  | ND            |                  | 0.263       | 0.189    |           | mg/Kg |   | 72   | 50 - 120     | 0   | 30    |
| Aroclor 1260                  | ND            |                  | 0.263       | 0.186    |           | mg/Kg |   | 71   | 50 - 125     | 0   | 30    |
| Surrogate                     | %Recovery     |                  | Qualifier   | Limits   |           |       |   |      |              |     |       |
| DCB Decachlorobiphenyl (Surr) | 74            |                  | p           | 45 - 120 |           |       |   |      |              |     |       |

## Method: 8290 - Dioxins and Furans (HRGC/HRMS)

**Lab Sample ID: MB 320-33622/1-A**  
**Matrix: Solid**  
**Analysis Batch: 33999**

**Client Sample ID: Method Blank**  
**Prep Type: Total/NA**  
**Prep Batch: 33622**

| Analyte           | MB     |           | RL  | EDL   | Unit | D | Prepared       | Analyzed       | Dil Fac |
|-------------------|--------|-----------|-----|-------|------|---|----------------|----------------|---------|
|                   | Result | Qualifier |     |       |      |   |                |                |         |
| 2,3,7,8-TCDD      | ND     |           | 1.0 | 0.017 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 2,3,7,8-TCDF      | ND     |           | 1.0 | 0.014 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 1,2,3,7,8-PeCDD   | ND     |           | 5.0 | 0.025 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 1,2,3,7,8-PeCDF   | ND     |           | 5.0 | 0.022 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 2,3,4,7,8-PeCDF   | ND     |           | 5.0 | 0.023 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 1,2,3,4,7,8-HxCDD | ND     |           | 5.0 | 0.017 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 1,2,3,6,7,8-HxCDD | ND     |           | 5.0 | 0.039 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 1,2,3,7,8,9-HxCDD | ND     |           | 5.0 | 0.012 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 1,2,3,4,7,8-HxCDF | 0.108  | J q       | 5.0 | 0.020 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 1,2,3,6,7,8-HxCDF | 0.0417 | J q       | 5.0 | 0.015 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 2,3,4,6,7,8-HxCDF | 0.0357 | J q       | 5.0 | 0.017 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 1,2,3,7,8,9-HxCDF | ND     |           | 5.0 | 0.019 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |

TestAmerica Honolulu

# QC Sample Results

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Method: 8290 - Dioxins and Furans (HRGC/HRMS) (Continued)

**Lab Sample ID: MB 320-33622/1-A**

**Matrix: Solid**

**Analysis Batch: 33999**

**Client Sample ID: Method Blank**

**Prep Type: Total/NA**

**Prep Batch: 33622**

| Analyte             | MB Result | MB Qualifier | RL  | EDL   | Unit | D | Prepared       | Analyzed       | Dil Fac |
|---------------------|-----------|--------------|-----|-------|------|---|----------------|----------------|---------|
| 1,2,3,4,6,7,8-HpCDD | 0.439     | J            | 5.0 | 0.030 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 1,2,3,4,6,7,8-HpCDF | 0.359     | J            | 5.0 | 0.027 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 1,2,3,4,7,8,9-HpCDF | ND        |              | 5.0 | 0.031 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| OCDD                | 6.38      | J            | 10  | 0.051 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| OCDF                | 0.486     | J q          | 10  | 0.043 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| Total TCDD          | 0.0915    | J q          | 1.0 | 0.017 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| Total TCDF          | ND        |              | 1.0 | 0.014 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| Total PeCDD         | ND        |              | 5.0 | 0.025 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| Total PeCDF         | ND        |              | 5.0 | 0.023 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| Total HxCDD         | ND        |              | 5.0 | 0.039 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| Total HxCDF         | 0.186     | J q          | 5.0 | 0.018 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| Total HpCDD         | 0.883     | J            | 5.0 | 0.030 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| Total HpCDF         | 0.492     | J q          | 5.0 | 0.029 | pg/g |   | 01/10/14 11:33 | 01/15/14 00:16 | 1       |

| Isotope Dilution        | MB %Recovery | MB Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|-------------------------|--------------|--------------|----------|----------------|----------------|---------|
| 13C-2,3,7,8-TCDD        | 65           |              | 40 - 135 | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 13C-2,3,7,8-TCDF        | 56           |              | 40 - 135 | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 13C-1,2,3,7,8-PeCDD     | 66           |              | 40 - 135 | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 13C-1,2,3,7,8-PeCDF     | 62           |              | 40 - 135 | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 13C-1,2,3,6,7,8-HxCDD   | 67           |              | 40 - 135 | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 13C-1,2,3,4,7,8-HxCDF   | 55           |              | 40 - 135 | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 13C-1,2,3,4,6,7,8-HpCDD | 78           |              | 40 - 135 | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 13C-1,2,3,4,6,7,8-HpCDF | 69           |              | 40 - 135 | 01/10/14 11:33 | 01/15/14 00:16 | 1       |
| 13C-OCDD                | 73           |              | 40 - 135 | 01/10/14 11:33 | 01/15/14 00:16 | 1       |

**Lab Sample ID: LCS 320-33622/2-A**

**Matrix: Solid**

**Analysis Batch: 33999**

**Client Sample ID: Lab Control Sample**

**Prep Type: Total/NA**

**Prep Batch: 33622**

| Analyte             | Spike Added | LCS Result | LCS Qualifier | Unit | D | %Rec | Limits   |
|---------------------|-------------|------------|---------------|------|---|------|----------|
| 2,3,7,8-TCDD        | 20.0        | 17.9       |               | pg/g |   | 89   | 60 - 138 |
| 2,3,7,8-TCDF        | 20.0        | 18.8       |               | pg/g |   | 94   | 56 - 158 |
| 1,2,3,7,8-PeCDD     | 100         | 101        |               | pg/g |   | 101  | 70 - 122 |
| 1,2,3,7,8-PeCDF     | 100         | 95.0       |               | pg/g |   | 95   | 69 - 134 |
| 2,3,4,7,8-PeCDF     | 100         | 94.5       |               | pg/g |   | 95   | 70 - 131 |
| 1,2,3,4,7,8-HxCDD   | 100         | 95.8       |               | pg/g |   | 96   | 60 - 138 |
| 1,2,3,6,7,8-HxCDD   | 100         | 102        |               | pg/g |   | 102  | 68 - 136 |
| 1,2,3,7,8,9-HxCDD   | 100         | 101        |               | pg/g |   | 101  | 68 - 138 |
| 1,2,3,4,7,8-HxCDF   | 100         | 99.6       |               | pg/g |   | 100  | 74 - 128 |
| 1,2,3,6,7,8-HxCDF   | 100         | 105        |               | pg/g |   | 105  | 67 - 140 |
| 2,3,4,6,7,8-HxCDF   | 100         | 109        |               | pg/g |   | 109  | 71 - 137 |
| 1,2,3,7,8,9-HxCDF   | 100         | 106        |               | pg/g |   | 106  | 72 - 134 |
| 1,2,3,4,6,7,8-HpCDD | 100         | 104        |               | pg/g |   | 104  | 71 - 128 |
| 1,2,3,4,6,7,8-HpCDF | 100         | 98.0       |               | pg/g |   | 98   | 71 - 134 |
| 1,2,3,4,7,8,9-HpCDF | 100         | 98.6       |               | pg/g |   | 99   | 68 - 129 |
| OCDD                | 200         | 204        |               | pg/g |   | 102  | 70 - 128 |
| OCDF                | 200         | 201        |               | pg/g |   | 100  | 63 - 141 |

TestAmerica Honolulu

# QC Sample Results

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Method: 8290 - Dioxins and Furans (HRGC/HRMS) (Continued)

Lab Sample ID: LCS 320-33622/2-A  
Matrix: Solid  
Analysis Batch: 33999

Client Sample ID: Lab Control Sample  
Prep Type: Total/NA  
Prep Batch: 33622

| Isotope Dilution        | LCS LCS   |           | Limits   |
|-------------------------|-----------|-----------|----------|
|                         | %Recovery | Qualifier |          |
| 13C-2,3,7,8-TCDD        | 51        |           | 40 - 135 |
| 13C-2,3,7,8-TCDF        | 41        |           | 40 - 135 |
| 13C-1,2,3,7,8-PeCDD     | 52        |           | 40 - 135 |
| 13C-1,2,3,7,8-PeCDF     | 48        |           | 40 - 135 |
| 13C-1,2,3,6,7,8-HxCDD   | 52        |           | 40 - 135 |
| 13C-1,2,3,4,7,8-HxCDF   | 45        |           | 40 - 135 |
| 13C-1,2,3,4,6,7,8-HpCDD | 62        |           | 40 - 135 |
| 13C-1,2,3,4,6,7,8-HpCDF | 56        |           | 40 - 135 |
| 13C-OCDD                | 58        |           | 40 - 135 |

## Method: 6010B - Metals (ICP)

Lab Sample ID: MB 580-151433/18-A  
Matrix: Solid  
Analysis Batch: 151504

Client Sample ID: Method Blank  
Prep Type: Total/NA  
Prep Batch: 151433

| Analyte  | MB MB  |           | RL   | MDL | Unit  | D | Prepared       | Analyzed       | Dil Fac |
|----------|--------|-----------|------|-----|-------|---|----------------|----------------|---------|
|          | Result | Qualifier |      |     |       |   |                |                |         |
| Arsenic  | ND     |           | 3.0  |     | mg/Kg |   | 12/30/13 12:18 | 12/31/13 11:13 | 1       |
| Barium   | ND     |           | 0.50 |     | mg/Kg |   | 12/30/13 12:18 | 12/31/13 11:13 | 1       |
| Cadmium  | ND     |           | 1.0  |     | mg/Kg |   | 12/30/13 12:18 | 12/31/13 11:13 | 1       |
| Chromium | ND     |           | 1.3  |     | mg/Kg |   | 12/30/13 12:18 | 12/31/13 11:13 | 1       |
| Lead     | ND     |           | 1.5  |     | mg/Kg |   | 12/30/13 12:18 | 12/31/13 11:13 | 1       |
| Selenium | ND     |           | 5.0  |     | mg/Kg |   | 12/30/13 12:18 | 12/31/13 11:13 | 1       |
| Silver   | ND     |           | 2.5  |     | mg/Kg |   | 12/30/13 12:18 | 12/31/13 11:13 | 1       |

Lab Sample ID: LCS 580-151433/19-A  
Matrix: Solid  
Analysis Batch: 151504

Client Sample ID: Lab Control Sample  
Prep Type: Total/NA  
Prep Batch: 151433

| Analyte  | Spike Added | LCS LCS |           | Unit  | D | %Rec | %Rec. Limits |
|----------|-------------|---------|-----------|-------|---|------|--------------|
|          |             | Result  | Qualifier |       |   |      |              |
| Arsenic  | 200         | 197     |           | mg/Kg |   | 99   | 80 - 120     |
| Barium   | 200         | 212     |           | mg/Kg |   | 106  | 80 - 120     |
| Cadmium  | 5.00        | 5.15    |           | mg/Kg |   | 103  | 80 - 120     |
| Chromium | 20.0        | 21.2    |           | mg/Kg |   | 106  | 80 - 120     |
| Lead     | 50.0        | 48.2    |           | mg/Kg |   | 96   | 80 - 120     |
| Selenium | 200         | 200     |           | mg/Kg |   | 100  | 80 - 120     |
| Silver   | 30.0        | 29.7    |           | mg/Kg |   | 99   | 75 - 120     |

Lab Sample ID: LCSD 580-151433/20-A  
Matrix: Solid  
Analysis Batch: 151504

Client Sample ID: Lab Control Sample Dup  
Prep Type: Total/NA  
Prep Batch: 151433

| Analyte  | Spike Added | LCSD LCSD |           | Unit  | D | %Rec | %Rec. Limits | RPD |       |
|----------|-------------|-----------|-----------|-------|---|------|--------------|-----|-------|
|          |             | Result    | Qualifier |       |   |      |              | RPD | Limit |
| Arsenic  | 200         | 196       |           | mg/Kg |   | 98   | 80 - 120     | 1   | 20    |
| Barium   | 200         | 208       |           | mg/Kg |   | 104  | 80 - 120     | 2   | 20    |
| Cadmium  | 5.00        | 5.00      |           | mg/Kg |   | 100  | 80 - 120     | 3   | 20    |
| Chromium | 20.0        | 20.8      |           | mg/Kg |   | 104  | 80 - 120     | 2   | 20    |
| Lead     | 50.0        | 47.4      |           | mg/Kg |   | 95   | 80 - 120     | 2   | 20    |
| Selenium | 200         | 198       |           | mg/Kg |   | 99   | 80 - 120     | 1   | 20    |

TestAmerica Honolulu

# QC Sample Results

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Method: 6010B - Metals (ICP) (Continued)

Lab Sample ID: LCSD 580-151433/20-A  
Matrix: Solid  
Analysis Batch: 151504

Client Sample ID: Lab Control Sample Dup  
Prep Type: Total/NA  
Prep Batch: 151433

| Analyte | Spike Added | LCSD Result | LCSD Qualifier | Unit  | D | %Rec | %Rec. Limits | RPD | RPD Limit |
|---------|-------------|-------------|----------------|-------|---|------|--------------|-----|-----------|
| Silver  | 30.0        | 29.6        |                | mg/Kg |   | 99   | 75 - 120     | 1   | 20        |

Lab Sample ID: 580-41776-1 MS  
Matrix: Solid  
Analysis Batch: 151504

Client Sample ID: HWL0089-01  
Prep Type: Total/NA  
Prep Batch: 151433

| Analyte  | Sample Result | Sample Qualifier | Spike Added | MS Result | MS Qualifier | Unit  | D | %Rec  | %Rec. Limits |
|----------|---------------|------------------|-------------|-----------|--------------|-------|---|-------|--------------|
| Arsenic  | 43            |                  | 41.9        | 82.2      |              | mg/Kg | * | 93    | 80 - 120     |
| Barium   | 940           |                  | 41.9        | 917       | 4            | mg/Kg | * | -55   | 80 - 120     |
| Cadmium  | 26            |                  | 1.05        | 31.1      | 4            | mg/Kg | * | 449   | 80 - 120     |
| Chromium | 210           |                  | 4.19        | 257       | 4            | mg/Kg | * | 1047  | 80 - 120     |
| Lead     | 5300          |                  | 10.5        | 4440      | 4            | mg/Kg | * | -7804 | 80 - 120     |
| Selenium | ND            |                  | 41.9        | 19.7      | F1           | mg/Kg | * | 47    | 80 - 120     |
| Silver   | 14            |                  | 6.29        | 20.1      |              | mg/Kg | * | 102   | 75 - 120     |

Lab Sample ID: 580-41776-1 MSD  
Matrix: Solid  
Analysis Batch: 151504

Client Sample ID: HWL0089-01  
Prep Type: Total/NA  
Prep Batch: 151433

| Analyte  | Sample Result | Sample Qualifier | Spike Added | MSD Result | MSD Qualifier | Unit  | D | %Rec  | %Rec. Limits | RPD | RPD Limit |
|----------|---------------|------------------|-------------|------------|---------------|-------|---|-------|--------------|-----|-----------|
| Arsenic  | 43            |                  | 41.0        | 84.9       |               | mg/Kg | * | 101   | 80 - 120     | 3   | 20        |
| Barium   | 940           |                  | 41.0        | 943        | 4             | mg/Kg | * | 7     | 80 - 120     | 3   | 20        |
| Cadmium  | 26            |                  | 1.02        | 29.4       | 4             | mg/Kg | * | 297   | 80 - 120     | 6   | 20        |
| Chromium | 210           |                  | 4.10        | 265        | 4             | mg/Kg | * | 1267  | 80 - 120     | 3   | 20        |
| Lead     | 5300          |                  | 10.2        | 6980       | 4 F2          | mg/Kg | * | 16766 | 80 - 120     | 44  | 20        |
| Selenium | ND            |                  | 41.0        | 20.1       | F1            | mg/Kg | * | 49    | 80 - 120     | 2   | 20        |
| Silver   | 14            |                  | 6.15        | 21.4       | F1            | mg/Kg | * | 127   | 75 - 120     | 7   | 20        |

## Method: 7471A - Mercury (CVAA)

Lab Sample ID: MB 580-151436/8-A ^10  
Matrix: Solid  
Analysis Batch: 151495

Client Sample ID: Method Blank  
Prep Type: Total/NA  
Prep Batch: 151436

| Analyte | MB Result | MB Qualifier | RL    | MDL | Unit  | D | Prepared       | Analyzed       | Dil Fac |
|---------|-----------|--------------|-------|-----|-------|---|----------------|----------------|---------|
| Mercury | ND        |              | 0.020 |     | mg/Kg |   | 12/31/13 08:25 | 12/31/13 09:57 | 10      |

Lab Sample ID: LCS 580-151436/9-A ^10  
Matrix: Solid  
Analysis Batch: 151495

Client Sample ID: Lab Control Sample  
Prep Type: Total/NA  
Prep Batch: 151436

| Analyte | Spike Added | LCS Result | LCS Qualifier | Unit  | D | %Rec | %Rec. Limits |
|---------|-------------|------------|---------------|-------|---|------|--------------|
| Mercury | 0.100       | 0.0994     |               | mg/Kg |   | 99   | 80 - 120     |

TestAmerica Honolulu

# QC Sample Results

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Method: 7471A - Mercury (CVAA) (Continued)

Lab Sample ID: LCSD 580-151436/10-A ^10

Matrix: Solid

Analysis Batch: 151495

Client Sample ID: Lab Control Sample Dup

Prep Type: Total/NA

Prep Batch: 151436

| Analyte | Spike Added | LCSD Result | LCSD Qualifier | Unit  | D | %Rec | %Rec. Limits | RPD | RPD Limit |
|---------|-------------|-------------|----------------|-------|---|------|--------------|-----|-----------|
| Mercury | 0.100       | 0.0993      |                | mg/Kg |   | 99   | 80 - 120     | 0   | 20        |

Lab Sample ID: 580-41776-1 MS

Matrix: Solid

Analysis Batch: 151495

Client Sample ID: HWL0089-01

Prep Type: Total/NA

Prep Batch: 151436

| Analyte | Sample Result | Sample Qualifier | Spike Added | MS Result | MS Qualifier | Unit  | D | %Rec | %Rec. Limits |
|---------|---------------|------------------|-------------|-----------|--------------|-------|---|------|--------------|
| Mercury | 1.1           |                  | 0.100       | 1.35      | 4            | mg/Kg | * | 277  | 80 - 120     |

Lab Sample ID: 580-41776-1 MSD

Matrix: Solid

Analysis Batch: 151495

Client Sample ID: HWL0089-01

Prep Type: Total/NA

Prep Batch: 151436

| Analyte | Sample Result | Sample Qualifier | Spike Added | MSD Result | MSD Qualifier | Unit  | D | %Rec | %Rec. Limits | RPD | RPD Limit |
|---------|---------------|------------------|-------------|------------|---------------|-------|---|------|--------------|-----|-----------|
| Mercury | 1.1           |                  | 0.103       | 1.44       | 4             | mg/Kg | * | 354  | 80 - 120     | 6   | 20        |

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TestAmerica Honolulu

# QC Association Summary

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## GC/MS Semi VOA

### Prep Batch: 153305

| Lab Sample ID      | Client Sample ID       | Prep Type | Matrix     | Method | Prep Batch |
|--------------------|------------------------|-----------|------------|--------|------------|
| HWL0089-01         | RHS-01                 | Total/NA  | Solid/Soil | 3546   |            |
| LCS 440-153305/2-A | Lab Control Sample     | Total/NA  | Solid      | 3546   |            |
| LCS 440-153305/3-A | Lab Control Sample Dup | Total/NA  | Solid      | 3546   |            |
| MB 440-153305/1-A  | Method Blank           | Total/NA  | Solid      | 3546   |            |

### Analysis Batch: 153775

| Lab Sample ID      | Client Sample ID       | Prep Type | Matrix     | Method | Prep Batch |
|--------------------|------------------------|-----------|------------|--------|------------|
| HWL0089-01         | RHS-01                 | Total/NA  | Solid/Soil | 8270C  | 153305     |
| LCS 440-153305/2-A | Lab Control Sample     | Total/NA  | Solid      | 8270C  | 153305     |
| LCS 440-153305/3-A | Lab Control Sample Dup | Total/NA  | Solid      | 8270C  | 153305     |
| MB 440-153305/1-A  | Method Blank           | Total/NA  | Solid      | 8270C  | 153305     |

## GC Semi VOA

### Prep Batch: 152687

| Lab Sample ID     | Client Sample ID | Prep Type | Matrix | Method | Prep Batch |
|-------------------|------------------|-----------|--------|--------|------------|
| MB 440-152687/1-A | Method Blank     | Total/NA  | Solid  | 3546   |            |

### Prep Batch: 153161

| Lab Sample ID       | Client Sample ID       | Prep Type | Matrix     | Method | Prep Batch |
|---------------------|------------------------|-----------|------------|--------|------------|
| 440-66257-E-2-A MS  | Matrix Spike           | Total/NA  | Solid      | 3546   |            |
| 440-66257-E-2-B MSD | Matrix Spike Duplicate | Total/NA  | Solid      | 3546   |            |
| HWL0089-01          | RHS-01                 | Total/NA  | Solid/Soil | 3546   |            |
| HWL0089-01          | RHS-01                 | Total/NA  | Solid/Soil | 3546   |            |
| LCS 440-153161/2-A  | Lab Control Sample     | Total/NA  | Solid      | 3546   |            |
| LCS 440-153161/4-A  | Lab Control Sample     | Total/NA  | Solid      | 3546   |            |
| LCS 440-153161/3-A  | Lab Control Sample Dup | Total/NA  | Solid      | 3546   |            |
| MB 440-153161/1-A   | Method Blank           | Total/NA  | Solid      | 3546   |            |

### Prep Batch: 153196

| Lab Sample ID       | Client Sample ID       | Prep Type | Matrix     | Method  | Prep Batch |
|---------------------|------------------------|-----------|------------|---------|------------|
| 440-65826-A-1-B MS  | Matrix Spike           | Total/NA  | Solid      | CA LUFT |            |
| 440-65826-A-1-C MSD | Matrix Spike Duplicate | Total/NA  | Solid      | CA LUFT |            |
| HWL0089-01          | RHS-01                 | Total/NA  | Solid/Soil | CA LUFT |            |
| LCS 440-153196/2-A  | Lab Control Sample     | Total/NA  | Solid      | CA LUFT |            |
| MB 440-153196/1-A   | Method Blank           | Total/NA  | Solid      | CA LUFT |            |

### Analysis Batch: 153341

| Lab Sample ID       | Client Sample ID       | Prep Type | Matrix | Method | Prep Batch |
|---------------------|------------------------|-----------|--------|--------|------------|
| 440-65826-A-1-B MS  | Matrix Spike           | Total/NA  | Solid  | 8015B  | 153196     |
| 440-65826-A-1-C MSD | Matrix Spike Duplicate | Total/NA  | Solid  | 8015B  | 153196     |
| LCS 440-153196/2-A  | Lab Control Sample     | Total/NA  | Solid  | 8015B  | 153196     |
| MB 440-153196/1-A   | Method Blank           | Total/NA  | Solid  | 8015B  | 153196     |

### Analysis Batch: 153342

| Lab Sample ID | Client Sample ID | Prep Type | Matrix     | Method | Prep Batch |
|---------------|------------------|-----------|------------|--------|------------|
| HWL0089-01    | RHS-01           | Total/NA  | Solid/Soil | 8015B  | 153196     |

TestAmerica Honolulu



# QC Association Summary

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## GC Semi VOA (Continued)

### Analysis Batch: 153427

| Lab Sample ID       | Client Sample ID       | Prep Type | Matrix     | Method | Prep Batch |
|---------------------|------------------------|-----------|------------|--------|------------|
| 440-66257-E-2-A MS  | Matrix Spike           | Total/NA  | Solid      | 8082   | 153161     |
| 440-66257-E-2-B MSD | Matrix Spike Duplicate | Total/NA  | Solid      | 8082   | 153161     |
| HWL0089-01          | RHS-01                 | Total/NA  | Solid/Soil | 8082   | 153161     |
| LCS 440-153161/4-A  | Lab Control Sample     | Total/NA  | Solid      | 8082   | 153161     |
| MB 440-152687/1-A   | Method Blank           | Total/NA  | Solid      | 8082   | 152687     |

### Analysis Batch: 153523

| Lab Sample ID      | Client Sample ID       | Prep Type | Matrix     | Method | Prep Batch |
|--------------------|------------------------|-----------|------------|--------|------------|
| HWL0089-01         | RHS-01                 | Total/NA  | Solid/Soil | 8081A  | 153161     |
| LCS 440-153161/2-A | Lab Control Sample     | Total/NA  | Solid      | 8081A  | 153161     |
| LCS 440-153161/3-A | Lab Control Sample Dup | Total/NA  | Solid      | 8081A  | 153161     |
| MB 440-153161/1-A  | Method Blank           | Total/NA  | Solid      | 8081A  | 153161     |

## Specialty Organics

### Prep Batch: 33622

| Lab Sample ID     | Client Sample ID   | Prep Type | Matrix     | Method | Prep Batch |
|-------------------|--------------------|-----------|------------|--------|------------|
| HWL0089-01 - RA   | RHS-01             | Total/NA  | Solid/Soil | 8290   |            |
| HWL0089-01        | RHS-01             | Total/NA  | Solid/Soil | 8290   |            |
| LCS 320-33622/2-A | Lab Control Sample | Total/NA  | Solid      | 8290   |            |
| MB 320-33622/1-A  | Method Blank       | Total/NA  | Solid      | 8290   |            |

### Analysis Batch: 33999

| Lab Sample ID     | Client Sample ID   | Prep Type | Matrix | Method | Prep Batch |
|-------------------|--------------------|-----------|--------|--------|------------|
| LCS 320-33622/2-A | Lab Control Sample | Total/NA  | Solid  | 8290   | 33622      |
| MB 320-33622/1-A  | Method Blank       | Total/NA  | Solid  | 8290   | 33622      |

### Analysis Batch: 34213

| Lab Sample ID   | Client Sample ID | Prep Type | Matrix     | Method | Prep Batch |
|-----------------|------------------|-----------|------------|--------|------------|
| HWL0089-01 - RA | RHS-01           | Total/NA  | Solid/Soil | 8290   | 33622      |

### Analysis Batch: 34312

| Lab Sample ID | Client Sample ID | Prep Type | Matrix     | Method | Prep Batch |
|---------------|------------------|-----------|------------|--------|------------|
| HWL0089-01    | RHS-01           | Total/NA  | Solid/Soil | 8290   | 33622      |

## Metals

### Prep Batch: 151433

| Lab Sample ID       | Client Sample ID       | Prep Type | Matrix     | Method | Prep Batch |
|---------------------|------------------------|-----------|------------|--------|------------|
| 580-41776-1 MS      | HWL0089-01             | Total/NA  | Solid      | 3050B  |            |
| 580-41776-1 MSD     | HWL0089-01             | Total/NA  | Solid      | 3050B  |            |
| HWL0089-01          | RHS-01                 | Total/NA  | Solid/Soil | 3050B  |            |
| LCS 580-151433/19-A | Lab Control Sample     | Total/NA  | Solid      | 3050B  |            |
| LCS 580-151433/20-A | Lab Control Sample Dup | Total/NA  | Solid      | 3050B  |            |
| MB 580-151433/18-A  | Method Blank           | Total/NA  | Solid      | 3050B  |            |

### Prep Batch: 151436

| Lab Sample ID   | Client Sample ID | Prep Type | Matrix | Method | Prep Batch |
|-----------------|------------------|-----------|--------|--------|------------|
| 580-41776-1 MS  | HWL0089-01       | Total/NA  | Solid  | 7471A  |            |
| 580-41776-1 MSD | HWL0089-01       | Total/NA  | Solid  | 7471A  |            |

TestAmerica Honolulu

# QC Association Summary

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Metals (Continued)

### Prep Batch: 151436 (Continued)

| Lab Sample ID            | Client Sample ID       | Prep Type | Matrix     | Method | Prep Batch |
|--------------------------|------------------------|-----------|------------|--------|------------|
| HWL0089-01               | RHS-01                 | Total/NA  | Solid/Soil | 7471A  |            |
| LCS 580-151436/9-A ^10   | Lab Control Sample     | Total/NA  | Solid      | 7471A  |            |
| LCSD 580-151436/10-A ^10 | Lab Control Sample Dup | Total/NA  | Solid      | 7471A  |            |
| MB 580-151436/8-A ^10    | Method Blank           | Total/NA  | Solid      | 7471A  |            |

### Analysis Batch: 151495

| Lab Sample ID            | Client Sample ID       | Prep Type | Matrix     | Method | Prep Batch |
|--------------------------|------------------------|-----------|------------|--------|------------|
| 580-41776-1 MS           | HWL0089-01             | Total/NA  | Solid      | 7471A  | 151436     |
| 580-41776-1 MSD          | HWL0089-01             | Total/NA  | Solid      | 7471A  | 151436     |
| HWL0089-01               | RHS-01                 | Total/NA  | Solid/Soil | 7471A  | 151436     |
| LCS 580-151436/9-A ^10   | Lab Control Sample     | Total/NA  | Solid      | 7471A  | 151436     |
| LCSD 580-151436/10-A ^10 | Lab Control Sample Dup | Total/NA  | Solid      | 7471A  | 151436     |
| MB 580-151436/8-A ^10    | Method Blank           | Total/NA  | Solid      | 7471A  | 151436     |

### Analysis Batch: 151504

| Lab Sample ID        | Client Sample ID       | Prep Type | Matrix     | Method | Prep Batch |
|----------------------|------------------------|-----------|------------|--------|------------|
| 580-41776-1 MS       | HWL0089-01             | Total/NA  | Solid      | 6010B  | 151433     |
| 580-41776-1 MSD      | HWL0089-01             | Total/NA  | Solid      | 6010B  | 151433     |
| HWL0089-01           | RHS-01                 | Total/NA  | Solid/Soil | 6010B  | 151433     |
| LCS 580-151433/19-A  | Lab Control Sample     | Total/NA  | Solid      | 6010B  | 151433     |
| LCSD 580-151433/20-A | Lab Control Sample Dup | Total/NA  | Solid      | 6010B  | 151433     |
| MB 580-151433/18-A   | Method Blank           | Total/NA  | Solid      | 6010B  | 151433     |

## General Chemistry

### Analysis Batch: 34148

| Lab Sample ID   | Client Sample ID | Prep Type | Matrix     | Method | Prep Batch |
|-----------------|------------------|-----------|------------|--------|------------|
| 320-5681-A-1 DU | Duplicate        | Total/NA  | Solid      | D 2216 |            |
| HWL0089-01      | RHS-01           | Total/NA  | Solid/Soil | D 2216 |            |

### Analysis Batch: 152388

| Lab Sample ID    | Client Sample ID | Prep Type | Matrix     | Method | Prep Batch |
|------------------|------------------|-----------|------------|--------|------------|
| 580-41946-A-1 DU | Duplicate        | Total/NA  | Solid      | D 2216 |            |
| HWL0089-01       | RHS-01           | Total/NA  | Solid/Soil | D 2216 |            |

### Analysis Batch: 156504

| Lab Sample ID     | Client Sample ID | Prep Type | Matrix     | Method   | Prep Batch |
|-------------------|------------------|-----------|------------|----------|------------|
| 440-67522-A-20 DU | Duplicate        | Total/NA  | Solid      | Moisture |            |
| HWL0089-01        | RHS-01           | Total/NA  | Solid/Soil | Moisture |            |

TestAmerica Honolulu

# Lab Chronicle

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

**Client Sample ID: RHS-01**

**Lab Sample ID: HWL0089-01**

**Date Collected: 12/20/13 15:00**

**Matrix: Solid/Soil**

**Date Received: 12/23/13 13:10**

**Percent Solids: 95.1**

| Prep Type | Batch Type | Batch Method | Run | Dilution Factor | Batch Number | Prepared or Analyzed | Analyst | Lab     |
|-----------|------------|--------------|-----|-----------------|--------------|----------------------|---------|---------|
| Total/NA  | Prep       | 3546         |     |                 | 153305       | 12/30/13 11:44       | HN      | TAL IRV |
| Total/NA  | Analysis   | 8270C        |     | 4               | 153775       | 01/02/14 16:39       | DF      | TAL IRV |
| Total/NA  | Prep       | CA LUFT      |     |                 | 153196       | 12/28/13 07:03       | HN      | TAL IRV |
| Total/NA  | Analysis   | 8015B        |     | 1               | 153342       | 12/30/13 08:36       | KW      | TAL IRV |
| Total/NA  | Prep       | 3546         |     |                 | 153161       | 12/30/13 11:45       | QCT     | TAL IRV |
| Total/NA  | Analysis   | 8082         |     | 1               | 153427       | 12/30/13 20:23       | JM      | TAL IRV |
| Total/NA  | Analysis   | 8081A        |     | 1               | 153523       | 12/31/13 14:53       | KS      | TAL IRV |
| Total/NA  | Prep       | 3546         |     |                 | 153161       | 12/27/13 17:32       | QCT     | TAL IRV |
| Total/NA  | Prep       | 8290         | RA  |                 | 33622        | 01/10/14 11:33       | TGL     | TAL SAC |
| Total/NA  | Analysis   | 8290         | RA  | 20              | 34213        | 01/17/14 20:40       | JRB     | TAL SAC |
| Total/NA  | Prep       | 8290         |     |                 | 33622        | 01/10/14 11:33       | TGL     | TAL SAC |
| Total/NA  | Analysis   | 8290         |     | 20              | 34312        | 01/20/14 16:28       | JRB     | TAL SAC |
| Total/NA  | Prep       | 7471A        |     |                 | 151436       | 12/31/13 08:25       | PAB     | TAL SEA |
| Total/NA  | Analysis   | 7471A        |     | 10              | 151495       | 12/31/13 10:04       | FCW     | TAL SEA |
| Total/NA  | Prep       | 3050B        |     |                 | 151433       | 12/30/13 12:18       | PAB     | TAL SEA |
| Total/NA  | Analysis   | 6010B        |     | 10              | 151504       | 12/31/13 11:27       | HJM     | TAL SEA |
| Total/NA  | Analysis   | D 2216       |     | 1               | 34148        | 01/17/14 13:30       | CV1     | TAL SAC |
| Total/NA  | Analysis   | Moisture     |     | 1               | 156504       | 01/16/14 13:00       | SP      | TAL IRV |
| Total/NA  | Analysis   | D 2216       |     | 1               | 152388       | 01/21/14 10:47       | RMB     | TAL SEA |

**Laboratory References:**

TAL IRV = TestAmerica Irvine, 17461 Derian Ave, Suite 100, Irvine, CA 92614-5817, TEL (949)261-1022

TAL SAC = TestAmerica Sacramento, 880 Riverside Parkway, West Sacramento, CA 95605, TEL (916)373-5600

TAL SEA = TestAmerica Seattle, 5755 8th Street East, Tacoma, WA 98424, TEL (253)922-2310

# Certification Summary

Client: Bureau Veritas  
 Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Laboratory: TestAmerica Honolulu

All certifications held by this laboratory are listed. Not all certifications are applicable to this report.

| Authority | Program       | EPA Region | Certification ID | Expiration Date |
|-----------|---------------|------------|------------------|-----------------|
| Florida   | NELAP         | 4          | E87907           | 06-30-14        |
| Hawaii    | State Program | 9          | N/A              | 06-28-14        |
| USDA      | Federal       |            | HON-S-206        | 01-31-15        |

## Laboratory: TestAmerica Irvine

All certifications held by this laboratory are listed. Not all certifications are applicable to this report.

| Authority                | Program                     | EPA Region | Certification ID  | Expiration Date |
|--------------------------|-----------------------------|------------|-------------------|-----------------|
| Alaska                   | State Program               | 10         | CA01531           | 06-30-14        |
| Arizona                  | State Program               | 9          | AZ0671            | 10-13-14        |
| California               | LA Cty Sanitation Districts | 9          | 10256             | 01-31-15        |
| California               | NELAP                       | 9          | 1108CA            | 01-31-14        |
| California               | State Program               | 9          | 2706              | 06-30-14        |
| Guam                     | State Program               | 9          | Cert. No. 12.002r | 01-23-14 *      |
| Hawaii                   | State Program               | 9          | N/A               | 01-31-14        |
| Nevada                   | State Program               | 9          | CA015312007A      | 07-31-14        |
| New Mexico               | State Program               | 6          | N/A               | 01-31-14        |
| Northern Mariana Islands | State Program               | 9          | MP0002            | 01-31-14        |
| Oregon                   | NELAP                       | 10         | 4005              | 09-12-14        |
| USDA                     | Federal                     |            | P330-09-00080     | 06-06-14        |
| USEPA UCMR               | Federal                     | 1          | CA01531           | 01-31-15        |

## Laboratory: TestAmerica Sacramento

All certifications held by this laboratory are listed. Not all certifications are applicable to this report.

| Authority                | Program       | EPA Region | Certification ID | Expiration Date |
|--------------------------|---------------|------------|------------------|-----------------|
| A2LA                     | A2LA          |            | NE-OS-22-13      | 01-31-14        |
| A2LA                     | DoD ELAP      |            | 2928-01          | 03-31-14        |
| Alaska (UST)             | State Program | 10         | UST-055          | 02-03-14 *      |
| Arizona                  | State Program | 9          | AZ0708           | 08-11-14        |
| Arkansas DEQ             | State Program | 6          | 88-0691          | 06-17-14        |
| California               | NELAP         | 9          | 1119CA           | 01-31-14        |
| Colorado                 | State Program | 8          | N/A              | 08-31-14        |
| Connecticut              | State Program | 1          | PH-0691          | 06-30-15        |
| Florida                  | NELAP         | 4          | E87570           | 06-30-14        |
| Guam                     | State Program | 9          | N/A              | 08-31-14        |
| Hawaii                   | State Program | 9          | N/A              | 01-31-14        |
| Illinois                 | NELAP         | 5          | 200060           | 03-17-15        |
| Kansas                   | NELAP         | 7          | E-10375          | 10-31-14        |
| Louisiana                | NELAP         | 6          | 30612            | 06-30-14        |
| Michigan                 | State Program | 5          | 9947             | 01-31-14        |
| Nebraska                 | State Program | 7          | NE-OS-22-13      | 01-31-14        |
| Nevada                   | State Program | 9          | CA44             | 07-31-14        |
| New Jersey               | NELAP         | 2          | CA005            | 06-30-14        |
| New York                 | NELAP         | 2          | 11666            | 03-31-14        |
| Northern Mariana Islands | State Program | 9          | MP0007           | 02-01-14        |
| Oregon                   | NELAP         | 10         | CA200005         | 03-28-14        |
| Pennsylvania             | NELAP         | 3          | 68-01272         | 03-31-14        |
| South Carolina           | State Program | 4          | 87014            | 06-30-14        |
| Texas                    | NELAP         | 6          | T104704399-08-TX | 05-31-14        |
| US Fish & Wildlife       | Federal       |            | LE148388-0       | 12-31-14        |

\* Expired certification is currently pending renewal and is considered valid.

# Certification Summary

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Laboratory: TestAmerica Sacramento (Continued)

All certifications held by this laboratory are listed. Not all certifications are applicable to this report.

| Authority          | Program       | EPA Region | Certification ID | Expiration Date |
|--------------------|---------------|------------|------------------|-----------------|
| USDA               | Federal       |            | P330-11-00436    | 12-30-14        |
| USEPA UCMR         | Federal       | 1          | CA00044          | 11-06-14        |
| Utah               | NELAP         | 8          | QUAN1            | 01-31-14        |
| Washington         | State Program | 10         | C581             | 05-05-14        |
| West Virginia DHHR | State Program | 3          | 9930C            | 02-03-14 *      |
| Wyoming            | State Program | 8          | 8TMS-Q           | 01-31-14        |

## Laboratory: TestAmerica Seattle

All certifications held by this laboratory are listed. Not all certifications are applicable to this report.

| Authority     | Program       | EPA Region | Certification ID | Expiration Date |
|---------------|---------------|------------|------------------|-----------------|
| Alaska (UST)  | State Program | 10         | UST-022          | 03-04-14        |
| California    | NELAP         | 9          | 01115CA          | 01-31-14        |
| L-A-B         | DoD ELAP      |            | L2236            | 01-19-16        |
| L-A-B         | ISO/IEC 17025 |            | L2236            | 01-19-16        |
| Montana (UST) | State Program | 8          | N/A              | 04-30-20        |
| Oregon        | NELAP         | 10         | WA100007         | 11-06-14        |
| USDA          | Federal       |            | P330-11-00222    | 05-20-14        |
| Washington    | State Program | 10         | C553             | 02-17-14        |

\* Expired certification is currently pending renewal and is considered valid.

# Method Summary

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

| Method   | Method Description                                     | Protocol | Laboratory |
|----------|--|----------|------------|
| 8270C    | Semivolatile Organic Compounds (GC/MS)                 | SW846    | TAL IRV    |
| 8015B    | Diesel Range Organics (DRO) (GC)                       | SW846    | TAL IRV    |
| 8081A    | Organochlorine Pesticides (GC)                         | SW846    | TAL IRV    |
| 8082     | Polychlorinated Biphenyls (PCBs) by Gas Chromatography | SW846    | TAL IRV    |
| 8290     | Dioxins and Furans (HRGC/HRMS)                         | SW846    | TAL SAC    |
| 6010B    | Metals (ICP)   | SW846    | TAL SEA    |
| 7471A    | Mercury (CVAA)   | SW846    | TAL SEA    |
| D 2216   | Percent Moisture                                       | ASTM     | TAL SAC    |
| D 2216   | Percent Moisture                                       | ASTM     | TAL SEA    |
| Moisture | Percent Moisture                                       | EPA      | TAL IRV    |

#### Protocol References:

ASTM = ASTM International

EPA = US Environmental Protection Agency

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

#### Laboratory References:

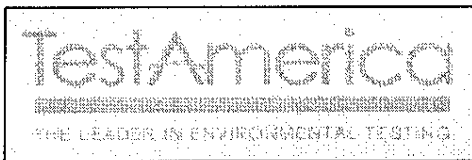
TAL IRV = TestAmerica Irvine, 17461 Derian Ave, Suite 100, Irvine, CA 92614-5817, TEL (949)261-1022

TAL SAC = TestAmerica Sacramento, 880 Riverside Parkway, West Sacramento, CA 95605, TEL (916)373-5600

TAL SEA = TestAmerica Seattle, 5755 8th Street East, Tacoma, WA 98424, TEL (253)922-2310







Rush TAT Confirmation (Initial/Date) \_\_\_\_\_

### Sample Receipt Checklist

Client Name: Bureau Veritas Date/ Time Received: 12/23/13 1310

Received By: [Signature]

Matrices: Soil

Carrier: Client

Airbill# :

- Shipping container/cooler in good condition? Yes  No  Not Present
- Chain of Custody present? Yes  No
- Chain of Custody Signed when relinquished and received? Yes  No
- Chain of Custody agrees with sample labels? Yes  No
- Samples in proper container/bottle? Yes  No
- Sample containers intact? Yes  No
- Sample containers on ice? Yes  No  Type: Wet
- Sufficient sample volume for indicated test? Yes  No
- All samples received within holding time? Yes  No
- Water - VOA Vials have Zero Headspace? Yes  No  No VOA vials present:
- Water - pH acceptable upon receipt? Yes  No  Not Checked:
- pH Adjusted? Yes  No  Final pH: \_\_\_\_\_
- Encores / MI-VOC / 5035 Vials Present? Yes  No  Location: \_\_\_\_\_
- Sample Filtration Needed? Yes  No  Filtered in Field:
- Dry Weight Corrected Results? Yes  No  Take Action:
- DODQSM / QAPP Project? Yes  No  Type: \_\_\_\_\_
- Temperature Blank Present? Yes  No
- Sample Container Temperature: 2 °C

### Comments/ Sampling Handling Notes:

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# Toxicity Summary

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

Client Sample ID: RHS-01

Lab Sample ID: HWL0089-01

| Analyte                | Result | Qualifier | NONE |     | Unit | WHO 2005<br>ND = 0 |                    | Method |        |
|------------------------|--------|-----------|------|-----|------|--------------------|--------------------|--------|--------|
|                        |        |           | TEF  | TEQ |      | TEF                | TEQ                |        |        |
| Total Dioxin/Furan TEQ |        |           |      |     | pg/g |                    | 2000               | TEQ    |        |
| Analyte                | Result | Qualifier | RL   |     | EDL  | Unit               | WHO 2005<br>ND = 0 |        | Method |
|                        |        |           | TEF  | TEQ |      |                    | TEF                | TEQ    |        |
| 2,3,7,8-TCDD           | 88     |           | 21   | 11  | pg/g | 1                  | 88                 | 8290   |        |
| 1,2,3,7,8-PeCDD        | 340    |           | 100  | 52  | pg/g | 1                  | 340                | 8290   |        |
| 1,2,3,7,8-PeCDF        | 1100   |           | 100  | 49  | pg/g | 0.03               | 33                 | 8290   |        |
| 2,3,4,7,8-PeCDF        | 2300   |           | 100  | 52  | pg/g | 0.3                | 690                | 8290   |        |
| 1,2,3,4,7,8-HxCDD      | 250    |           | 100  | 11  | pg/g | 0.1                | 25                 | 8290   |        |
| 1,2,3,6,7,8-HxCDD      | 420    |           | 100  | 8.5 | pg/g | 0.1                | 42                 | 8290   |        |
| 1,2,3,7,8,9-HxCDD      | 320    |           | 100  | 8.2 | pg/g | 0.1                | 32                 | 8290   |        |
| 1,2,3,4,7,8-HxCDF      | 1900   | B         | 100  | 85  | pg/g | 0.1                | 190                | 8290   |        |
| 1,2,3,6,7,8-HxCDF      | 1400   | B         | 100  | 71  | pg/g | 0.1                | 140                | 8290   |        |
| 2,3,4,6,7,8-HxCDF      | 1900   | B         | 100  | 79  | pg/g | 0.1                | 190                | 8290   |        |
| 1,2,3,7,8,9-HxCDF      | ND     |           | 100  | 89  | pg/g | 0.1                | 0.00               | 8290   |        |
| 1,2,3,4,6,7,8-HpCDD    | 2400   | B         | 100  | 16  | pg/g | 0.01               | 24                 | 8290   |        |
| 1,2,3,4,6,7,8-HpCDF    | 7600   | B         | 100  | 98  | pg/g | 0.01               | 76                 | 8290   |        |
| 1,2,3,4,7,8,9-HpCDF    | 220    | G         | 120  | 120 | pg/g | 0.01               | 2.2                | 8290   |        |
| OCDD                   | 5600   | B         | 210  | 19  | pg/g | 0.0003             | 1.7                | 8290   |        |
| OCDF                   | 1700   | B         | 210  | 5.7 | pg/g | 0.0003             | 0.51               | 8290   |        |
| 2,3,7,8-TCDF - RA      | 1400   |           | 21   | 3.2 | pg/g | 0.1                | 140                | 8290   |        |

**TEF Reference:**

WHO 2005 = World Health Organization (WHO) 2005 TEF, Dioxins, Furans and PCB Congeners

TestAmerica Honolulu

# Isotope Dilution Summary

Client: Bureau Veritas  
Project/Site: 17012-012148.00

TestAmerica Job ID: HWL0089

## Method: 8290 - Dioxins and Furans (HRGC/HRMS)

Matrix: Solid

Prep Type: Total/NA

|                   |                    | Percent Isotope Dilution Recovery (Acceptance Limits) |                  |                   |                    |                    |                    |                   |                    |
|-------------------|--------------------|---|------------------|-------------------|--------------------|--------------------|--------------------|-------------------|--------------------|
| Lab Sample ID     | Client Sample ID   | TCDD<br>(40-135)                                      | TCDF<br>(40-135) | PeCDD<br>(40-135) | PeCDF1<br>(40-135) | HxCDD2<br>(40-135) | HxCDF1<br>(40-135) | HpCDD<br>(40-135) | HpCDF1<br>(40-135) |
| LCS 320-33622/2-A | Lab Control Sample | 51  | 41               | 52                | 48                 | 52                 | 45                 | 62                | 56                 |
| MB 320-33622/1-A  | Method Blank       | 65  | 56               | 66                | 62                 | 67                 | 55                 | 78                | 69                 |

|                   |                    | Percent Isotope Dilution Recovery (Acceptance Limits) |  |  |  |  |  |  |  |
|-------------------|--------------------|---|--|--|--|--|--|--|--|
| Lab Sample ID     | Client Sample ID   | OCDD<br>(40-135)                                      |  |  |  |  |  |  |  |
| LCS 320-33622/2-A | Lab Control Sample | 58  |  |  |  |  |  |  |  |
| MB 320-33622/1-A  | Method Blank       | 73  |  |  |  |  |  |  |  |

**Surrogate Legend**

- TCDD = 13C-2,3,7,8-TCDD
- TCDF = 13C-2,3,7,8-TCDF
- PeCDD = 13C-1,2,3,7,8-PeCDD
- PeCDF1 = 13C-1,2,3,7,8-PeCDF
- HxCDD2 = 13C-1,2,3,6,7,8-HxCDD
- HxCDF1 = 13C-1,2,3,4,7,8-HxCDF
- HpCDD = 13C-1,2,3,4,6,7,8-HpCDD
- HpCDF1 = 13C-1,2,3,4,6,7,8-HpCDF
- OCDD = 13C-OCDD

## Method: 8290 - Dioxins and Furans (HRGC/HRMS)

Matrix: Solid/Soil

Prep Type: Total/NA

|                 |                  | Percent Isotope Dilution Recovery (Acceptance Limits) |                  |                   |                    |                    |                    |                   |                    |
|-----------------|------------------|---|------------------|-------------------|--------------------|--------------------|--------------------|-------------------|--------------------|
| Lab Sample ID   | Client Sample ID | TCDD<br>(40-135)                                      | TCDF<br>(40-135) | PeCDD<br>(40-135) | PeCDF1<br>(40-135) | HxCDD2<br>(40-135) | HxCDF1<br>(40-135) | HpCDD<br>(40-135) | HpCDF1<br>(40-135) |
| HWL0089-01 - RA | RHS-01           |   | 69               |                   |                    |                    |                    |                   |                    |
| HWL0089-01      | RHS-01           | 73  | 78               | 80                | 80                 | 62                 | 58                 | 49                | 54                 |

|                 |                  | Percent Isotope Dilution Recovery (Acceptance Limits) |  |  |  |  |  |  |  |
|-----------------|------------------|---|--|--|--|--|--|--|--|
| Lab Sample ID   | Client Sample ID | OCDD<br>(40-135)                                      |  |  |  |  |  |  |  |
| HWL0089-01 - RA | RHS-01           |   |  |  |  |  |  |  |  |
| HWL0089-01      | RHS-01           | 50  |  |  |  |  |  |  |  |

**Surrogate Legend**

- TCDD = 13C-2,3,7,8-TCDD
- TCDF = 13C-2,3,7,8-TCDF
- PeCDD = 13C-1,2,3,7,8-PeCDD
- PeCDF1 = 13C-1,2,3,7,8-PeCDF
- HxCDD2 = 13C-1,2,3,6,7,8-HxCDD
- HxCDF1 = 13C-1,2,3,4,7,8-HxCDF
- HpCDD = 13C-1,2,3,4,6,7,8-HpCDD
- HpCDF1 = 13C-1,2,3,4,6,7,8-HpCDF
- OCDD = 13C-OCDD

TestAmerica Honolulu



**APPENDIX B**

**LABORATORY ANALYTICAL REPORTS  
AND  
CHAIN OF CUSTODY FORMS (NVL)**

December 27, 2013

Kenney Gomes  
Bureau Veritas North America, Inc.-Hawaii  
841 Bishop Street, Suite 1100  
Honolulu, HI 96813



Laboratory | Management | Training

**RE: Bulk Asbestos Fiber Analysis, NVL Batch # 1321182.00**

Dear Mr. Gomes,

Enclosed please find test results for the bulk samples submitted to our laboratory for analysis. Examination of these samples was conducted for the presence of identifiable asbestos fibers using polarized light microscopy (PLM) with dispersion staining in accordance with both U.S. EPA 600/M4-82-020, Interim Method for Determination of Asbestos in Bulk Insulation Samples, as found in 40 CFR, Part 763, Subpart E, Appendix E (formerly Subpart F, Appendix A), and U.S. EPA 600/R-93/116 (July 1993) Test Methods.

For samples containing more than one separable layer of materials, the report will include findings for each layer (labeled Layer 1 and Layer 2, etc. for each individual layer). The asbestos concentration in the sample is determined by visual estimation.

For those samples with asbestos concentrations between 1 and 10 percent based on visual estimation, the EPA recommends a procedure known as point counting (NESHAPS, 40 CFR Part 61). Point counting is a statistically more accurate means of quantification for samples with low concentrations of asbestos. If you would like us to further refine the concentration estimates of asbestos in these samples using point counting, please let me know.

This report is considered highly confidential and will not be released without your approval. Samples are archived for two weeks following analysis. Samples that are not retrieved by the client are discarded after two weeks.

Thank you for using our laboratory services. Please do not hesitate to call if there is anything further we can assist you with.

Sincerely,

A handwritten signature in black ink, appearing to read "Nick Ly".

Nick Ly, Technical Director



Lab Code: 102083-0

1.888.NVL.LABS  
1.888.(685.5227)  
www.nvllabs.com

Enc.: Sample Results

NVL Laboratories, Inc.

4708 Aurora Ave N, Seattle, WA 98103

www.carrollcox.com 808-782-6627 | f 206.547.0100 | 206.634.1936



**Bulk Asbestos Fibers Analysis**

By Polarized Light Microscopy

Client: Bureau Veritas North America, Inc.-Hawaii  
 Address: 841 Bishop Street, Suite 1100  
 Honolulu, HI 96813

**Batch #: 1321182.00**  
 Client Project #: 1702-012148.00  
 Date Received: 12/26/2013  
 Samples Received: 6  
 Samples Analyzed: 6  
 Method: EPA/600/R-93/116  
 & EPA/600/M4-82-020

**Attention: Mr. Kenney Gomes**  
 Project Location: Radford HS

**Lab ID: 13147580 Client Sample #: RHS-1**

Location: Radford HS

Layer 1 of 1 Description: Off-white fibrous material

|                               |                           |                         |
|-------------------------------|---------------------------|-------------------------|
| Non-Fibrous Materials:        | Other Fibrous Materials:% | <b>Asbestos Type: %</b> |
| Binder/Filler, Fine particles | Cellulose 33%             | <b>Chrysotile 55%</b>   |

**Lab ID: 13147581 Client Sample #: RHS-2**

Location: Radford HS

Layer 1 of 1 Description: Off-white fibrous material

|                               |                           |                         |
|-------------------------------|---------------------------|-------------------------|
| Non-Fibrous Materials:        | Other Fibrous Materials:% | <b>Asbestos Type: %</b> |
| Binder/Filler, Fine particles | Cellulose 37%             | <b>Chrysotile 51%</b>   |

**Lab ID: 13147582 Client Sample #: RHS-3**

Location: Radford HS

Layer 1 of 1 Description: Off-white fibrous material

|                               |                           |                         |
|-------------------------------|---------------------------|-------------------------|
| Non-Fibrous Materials:        | Other Fibrous Materials:% | <b>Asbestos Type: %</b> |
| Binder/Filler, Fine particles | Cellulose 35%             | <b>Chrysotile 53%</b>   |

**Lab ID: 13147583 Client Sample #: RHS-4**

Location: Radford HS

Layer 1 of 3 Description: Off-white/brown fibrous material with debris

|                             |                           |                         |
|-----------------------------|---------------------------|-------------------------|
| Non-Fibrous Materials:      | Other Fibrous Materials:% | <b>Asbestos Type: %</b> |
| Binder/Filler, Glass debris | Glass fibers 68%          | <b>None Detected ND</b> |
|                             | Cellulose 13%             |                         |

Layer 2 of 3 Description: Off-white brittle material with woven fibrous material and silver paper

|                                  |                           |                         |
|----------------------------------|---------------------------|-------------------------|
| Non-Fibrous Materials:           | Other Fibrous Materials:% | <b>Asbestos Type: %</b> |
| Binder/Filler, Calcareous binder | Cellulose 28%             | <b>None Detected ND</b> |

**Sampled by:** Client

**Analyzed by:** Jacob Laugeson

**Reviewed by:** Nick Ly

**Date:** 12/27/2013

**Date:** 12/27/2013

Nick Ly, Technical Director

Note: If samples are not homogeneous, then subsamples of the components were analyzed separately. All bulk samples are analyzed using both EPA 600/R-93/116 and 600/M4-82-020 Methods with the following measurement uncertainties for the reported % Asbestos (1%=0-3%, 5%=1-9%, 10%=5-15%, 20%=10-30%, 50%=40-60%). This report relates only to the items tested. If sample was not collected by NVL personnel, then the accuracy of the results is limited by the methodology and acuity of the sample collector. This report shall not be reproduced except in full, without written approval of NVL Laboratories, Inc. It shall not be used to claim product endorsement by NVLAP or any other agency of the US Government.

### Bulk Asbestos Fibers Analysis

By Polarized Light Microscopy

Client: Bureau Veritas North America, Inc.-Hawaii  
Address: 841 Bishop Street, Suite 1100  
Honolulu, HI 96813

**Batch #: 1321182.00**  
Client Project #: 1702-012148.00  
Date Received: 12/26/2013  
Samples Received: 6  
Samples Analyzed: 6  
Method: EPA/600/R-93/116  
& EPA/600/M4-82-020

**Attention: Mr. Kenney Gomes**  
Project Location: Radford HS

**Layer 3 of 3** Description: White brittle material

Non-Fibrous Materials:  
Calcareous binder, Mineral grains

Other Fibrous Materials:%  
Mineral fibers 8%  
Cellulose 3%

**Asbestos Type: %  
None Detected ND**

**Lab ID: 13147584** Client Sample #: **RHS-5**

Location: Radford HS

**Layer 1 of 1** Description: Orange fibrous material with debris

Non-Fibrous Materials:  
Binder/Filler, Fine particles, Glass debris

Other Fibrous Materials:%  
Glass fibers 52%  
Cellulose 33%

**Asbestos Type: %  
None Detected ND**

**Lab ID: 13147585** Client Sample #: **RHS-6**

Location: Radford HS

**Layer 1 of 1** Description: Brown fibrous material with debris

Non-Fibrous Materials:  
Binder/Filler, Calcareous particles, Glass debris

Other Fibrous Materials:%  
Glass fibers 61%  
Cellulose 19%

**Asbestos Type: %  
None Detected ND**

**Sampled by: Client**

**Analyzed by: Jacob Laugeson**

**Reviewed by: Nick Ly**

**Date: 12/27/2013**

**Date: 12/27/2013**

Nick Ly, Technical Director

Note: If samples are not homogeneous, then subsamples of the components were analyzed separately. All bulk samples are analyzed using both EPA 600/R-93/116 and 600/M4-82-020 Methods with the following measurement uncertainties for the reported % Asbestos (1%=0-3%, 5%=1-9%, 10%=5-15%, 20%=10-30%, 50%=40-60%). This report relates only to the items tested. If sample was not collected by NVL personnel, then the accuracy of the results is limited by the methodology and acuity of the sample collector. This report shall not be reproduced except in full, without written approval of NVL Laboratories, Inc. It shall not be used to claim product endorsement by NVLAP or any other agency of the US Government.

**NVL Laboratories, Inc.**

4708 Aurora Ave N, Seattle, WA 98103  
 Tel: 206.547.0100 Emerg. Pager: 206.344.1878  
 Fax: 206.634.1936 1.888.NVL.LABS (685.5227)

**CHAIN of CUSTODY  
 SAMPLE LOG**

**NVL Batch ID  
 1321182**

**Client** Bureau Veritas North America, Inc.  
**Street** 841 Bishop Street, Suite 1100  
 Honolulu, HI 96813

**NVL Batch Number** \_\_\_\_\_  
**Client Job Number** 17012-02/49,00  
**Total Samples** 6

**Project Manager** Ken Gomes  
**Project Location** Radford HS

**Turn Around Time**  1-Hr  24-Hrs  4 Days  
 2-Hrs  2 Days  5 Days  
 4-Hrs  3 Days  6 to 10 Days  
 Please call for TAT less than 24 Hrs

**Email address** kenney.gomes@us.bureauveritas.com

**Phone:** (808) 531-6708 **Fax:** (808) 537-4084 **Cell** \_\_\_\_\_

|   |  |  |   |  |   |
|---|--|--|---|--|---|
| <input type="checkbox"/> Asbestos Air   | <input type="checkbox"/> PCM (NIOSH 7400)  | <input type="checkbox"/> TEM (NIOSH 7402)  | <input type="checkbox"/> TEM (AHERA)  | <input type="checkbox"/> TEM (EPA Level II)  | <input type="checkbox"/> Other _____  |
| <input checked="" type="checkbox"/> Asbestos Bulk                                       | <input checked="" type="checkbox"/> PLM (EPA/600/R-93/116)                                     | <input type="checkbox"/> PLM (EPA Point Count)   | <input type="checkbox"/> PLM (EPA Gravimetry)   | <input type="checkbox"/> TEM Bulk  |   |
| <b>METALS</b><br><input type="checkbox"/> Total Metals<br><input type="checkbox"/> TCLP | <b>Det. Limit</b><br><input type="checkbox"/> ppm (AAS)<br><input type="checkbox"/> ppb (GFAA) | <b>Matrix</b><br><input type="checkbox"/> Air Filter<br><input type="checkbox"/> Drinking water<br><input type="checkbox"/> Dust/wipe<br><input type="checkbox"/> Soil | <input type="checkbox"/> Paint Chips<br><input type="checkbox"/> Paint Chips (Area)<br><input type="checkbox"/> Waste Water | <b>RCRA Metals</b><br><input type="checkbox"/> Arsenic (As)<br><input type="checkbox"/> Barium (Ba)<br><input type="checkbox"/> Cadmium (Cd)<br><input type="checkbox"/> Chromium (Cr) | <input type="checkbox"/> All 8<br><input type="checkbox"/> Lead (Pb)<br><input type="checkbox"/> Mercury (Hg)<br><input type="checkbox"/> Selenium (Se)<br><input type="checkbox"/> Silver (Ag) |
| <input type="checkbox"/> Other Types of Analysis  | <input type="checkbox"/> Fiberglass<br><input type="checkbox"/> Silica                         | <input type="checkbox"/> Nuisance Dust<br><input type="checkbox"/> Respirable Dust   | <input type="checkbox"/> Other (Specify) _____  |  |   |

**Condition of Package:**  Good  Damaged (no spillage)  Severe damage (spillage)

| Seq. # | Lab ID | Client Sample Number | Comments                       | A/R |
|--------|--------|----------------------|--------------------------------|-----|
| 1      |        | RMS-1                | See the attached data sheet(s) |     |
| 2      |        | RMS-2                |                                |     |
| 3      |        | RMS-3                |                                |     |
| 4      |        | RMS-4                |                                |     |
| 5      |        | RMS-5                |                                |     |
| 6      |        | RMS-6                |                                |     |
| 7      |        |                      |                                |     |
| 8      |        |                      |                                |     |
| 9      |        |                      |                                |     |
| 10     |        |                      |                                |     |
| 11     |        |                      |                                |     |
| 12     |        |                      |                                |     |
| 13     |        |                      |                                |     |
| 14     |        |                      |                                |     |
| 15     |        |                      |                                |     |

|                   | Print Below   | Sign Below         | Company       | Date     | Time |
|-------------------|---------------|--------------------|---------------|----------|------|
| Sampled by        | Ken Gomes     | <i>[Signature]</i> | BVNA - Hawaii | 12/20/13 | 1500 |
| Relinquished by   | Ken Gomes     | <i>[Signature]</i> | BVNA - Hawaii | 12/23/13 | 1230 |
| Received by       | Madon Kiki    | <i>[Signature]</i> | NL            | 12/20/13 | 1530 |
| Analyzed by       | Jacob Luyssen | <i>[Signature]</i> | NL            | 12-27-13 | 1315 |
| Results Called by |               |                    |               |          |      |
| Results Faxed by  |               |                    |               |          |      |

**Special Instructions:** Unless requested in writing, all samples will be disposed of two (2) weeks after analysis.  
 \*Please provide separate reports for each project number





**APPENDIX C**

**LABORATORY ANALYTICAL REPORTS  
AND  
CHAIN OF CUSTODY FORMS (ESN)**



## **Environmental Services Network**

February 8, 2014

Ken Gomes  
Bureau Veritas North America, Inc  
970 N. Kalaheo, Ste C-316  
Kailua, HI 96734

**SUBJECT:** DATA REPORT – 17012.012148.48, Radford High School

**ESN Project # D1402050037**

Mr. Gomes:

Please find enclosed a data report for the samples analyzed from the above referenced project for Bureau Veritas North America, Inc. The samples were received intact. Applicable detection limits, QA/QC data, and any issues encountered during analysis are included in the report.

The following tests were conducted:

- Multi-increment sample processing by Hawaii DOH Method.
- Analyses for TPH as diesel by EPA 8015 mod.
- Analyses for TPH as oil by EPA 8015 mod.
- Analyses for organochlorine pesticides by EPA 8081 mod.
- Analyses for polychlorinated biphenyls by EPA 8082 mod.
- Analyses for semi-volatile organic compounds by EPA method 8270.
- Analyses for total RCRA 8 by EPA 6020.
- Analyses for dioxins & furans by EPA 8290.

ESN appreciates the opportunity to have provided analytical services to Bureau Veritas North America, Inc on this project. If you have any further questions relating to the data or report, please do not hesitate to contact us.

Sincerely,

A handwritten signature in blue ink that reads 'Karen Carvalho'.

Karen Carvalho  
Operations Manager

ESN Pacific  
2020-B Kahai Street  
Honolulu, HI 96819

Ph: (808) 847-0067  
esn@esnpacific.com



# ESN PACIFIC'S CHAIN-OF-CUSTODY RECORD

CLIENT: Bureau Venues TAT (circle one): 5-day 48-hr. or Other: \_\_\_\_\_ PAGE 1 OF 2  
 ADDRESS: 841 Bishop Street DATE: 4/5/14  
 PHONE: 531-6708 FAX: 537-4084 ESN PROJECT #: 01402050037  
 EMAIL: kenney.gomes@bureauvenues.com LOCATION/PROJECT NAME: Rutherford High School / A012-012140.408  
 CLIENT PROJECT #: \_\_\_\_\_ COLLECTOR: K6/RE/MC/NO DATE COLLECTED: 3/4/14 & 2/5/14  
 Project Manager: Ken Gomes

| Sample ID# | Sample Type | Time        | Container Type | 8260 HVOX ext 5035? Y N | 8260 VOC ext 5035? Y N | 8260 MBE ext 5035? Y N | 8015 Fuel Scan only Y N | 8015 TPH-Diesel Y N | 8015 TPH-Oil Y N | 8081 Chlor. Pesticides Y N | 8270 PCB Y N | 8270 FAH EPA-16 Y N | 8270 FAH DOH-4 Y N | 1010 SVOC Y N | 1010 FlashPoint (Ignitibility) Y N | Total Metals Y N | TCLP Y N | MI Sample Prep Y N | Comments | # of Containers |
|------------|-------------|-------------|----------------|-------------------------|------------------------|------------------------|-------------------------|---------------------|------------------|----------------------------|--------------|---------------------|--------------------|---------------|------------------------------------|------------------|----------|--------------------|----------|-----------------|
| 1          | RHS-DU-1    | 2/4/14 1525 |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 2          | RHS-DU-1.2  | 2/5/14 0933 |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 3          | RHS-DU-1.3  | ↓ 1030      |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 4          | RHS-DU-2    | 2/4/14 1500 |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 5          | RHS-DU-3    | ↓ 1403      |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 6          | RHS-DU-4    | ↓ 1145      |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 7          | RHS-DU-5    | ↓ 1220      |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 8          | RHS-DU-6    | ↓ 1402      |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 9          | RHS-DU-7    | ↓ 1131      |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 10         | RHS-DU-8    | ↓ 1424      |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 11         | RHS-DU-9    | ↓ 1455      |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 12         | RHS-DU-10   | 4/5/14 0915 |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 13         | RHS-DU-11   | ↓ 0900      |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 14         | RHS-DU-12   | ↓ 0945      |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 15         | RHS-DU-13   | ↓ 0745      |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 16         | RHS-DU-14   | 2/4/14 1252 |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 17         | RHS-DU-15   | ↓ 1245      |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 18         | RHS-DU-16   | ↓ 1411      |                | X                       | X                      | X                      | X                       | X                   | X                | X                          | X            | X                   | X                  | X             | X                                  | X                | X        | X                  |          |                 |
| 19         |             |             |                |                         |                        |                        |                         |                     |                  |                            |              |                     |                    |               |                                    |                  |          |                    |          |                 |
| 20         |             |             |                |                         |                        |                        |                         |                     |                  |                            |              |                     |                    |               |                                    |                  |          |                    |          |                 |

RELINQUISHED BY: (Signature) [Signature] DATE/TIME 2/5/14 1030 RECEIVED BY (Signature) [Signature] DATE/TIME 2/5  
 RELINQUISHED BY: (Signature) \_\_\_\_\_ DATE/TIME \_\_\_\_\_ RECEIVED BY (Signature) \_\_\_\_\_ DATE/TIME \_\_\_\_\_  
 LABORATORY NOTES: MF Samples  
RUSH DIOXINS  
Hold samples for TCLP / SPLP Testing  
 SAMPLE RECEIPT: TOTAL # OF CONTAINERS 18 COC SEALS Y N / N A SEALS INTACT Y N / N A RECEIVED TEMP: 30C  
 SAMPLE DISPOSAL INSTRUCTIONS: \_\_\_\_\_ ESN @ \$5.00/sample or \_\_\_\_\_ Return to Client

# ESN PACIFIC'S CHAIN-OF-CUSTODY RECORD

CLIENT: Bureau Veritas  
 ADDRESS: 841 Bishop Street Suite 1100  
 PHONE: 531-6708 FAX: 537-4084  
 EMAIL: Kennedy.gomez@bureauveritas.com  
 CLIENT PROJECT #: 1701-912148-93 Project Manager: Ken Gomez  
 TAT (circle one): 5-day 48-hr. or Other: \_\_\_\_\_  
 DATE: 2/17/14 PAGE 2 OF 2  
 ESN PROJECT #: D1402050037  
 LOCATION/PROJECT NAME: Portland High School  
 COLLECTOR: Ky Hume DATE COLLECTED: 2/14/14

| Sample ID# | Depth | Time | Sample Type | Container Type | 8290 HVOC ext 5035? Y N | 8290 MBE ext 5035? Y N | 8015 TRH-Gas only | 8015 TRH-Diesel | 8081 Chlor. Pesticides | 8270 PAH EPA-16 | 1010 Flashpoint (Ignitibility) | Total Metals | RCRA 8 Metals | MI Sample Prep | Comments | # of Containers |
|------------|-------|------|-------------|----------------|-------------------------|------------------------|-------------------|-----------------|------------------------|-----------------|--------------------------------|--------------|---------------|----------------|----------|-----------------|
| 1          | 24114 | 1426 | DV-17       |                | X                       | X                      | X                 | X               | X                      | X               | X                              | X            | X             | X              |          |                 |
| 2          | 24115 | 1427 | DV-17.2     |                | X                       | X                      | X                 | X               | X                      | X               | X                              | X            | X             | X              |          |                 |
| 3          | 24115 | 1302 | DV-17.3     |                | X                       | X                      | X                 | X               | X                      | X               | X                              | X            | X             | X              |          |                 |
| 4          | 24115 | 1303 | DV-18       |                | X                       | X                      | X                 | X               | X                      | X               | X                              | X            | X             | X              |          |                 |
| 5          | 24115 | 1422 | DV-19       |                | X                       | X                      | X                 | X               | X                      | X               | X                              | X            | X             | X              |          |                 |
| 6          | 24115 | 1416 | DV-20       |                | X                       | X                      | X                 | X               | X                      | X               | X                              | X            | X             | X              |          |                 |
| 7          | 24115 | 1226 | DV-21       |                | X                       | X                      | X                 | X               | X                      | X               | X                              | X            | X             | X              |          |                 |
| 8          | 24115 | 1015 | DV-22       |                | X                       | X                      | X                 | X               | X                      | X               | X                              | X            | X             | X              |          |                 |
| 9          | 24115 | 1023 | DV-23       |                | X                       | X                      | X                 | X               | X                      | X               | X                              | X            | X             | X              |          |                 |
| 10         | 24115 | 1050 | DV-24       |                | X                       | X                      | X                 | X               | X                      | X               | X                              | X            | X             | X              |          |                 |
| 11         | 24115 | 1115 | DV-25       |                | X                       | X                      | X                 | X               | X                      | X               | X                              | X            | X             | X              |          |                 |
| 12         | 24115 | 1020 | DV-26       |                | X                       | X                      | X                 | X               | X                      | X               | X                              | X            | X             | X              |          |                 |
| 13         | 24115 | 1019 | DV-26.2     |                | X                       | X                      | X                 | X               | X                      | X               | X                              | X            | X             | X              |          |                 |
| 14         | 24115 | 1145 | DV-26.3     |                | X                       | X                      | X                 | X               | X                      | X               | X                              | X            | X             | X              |          |                 |
| 15         | 24115 | 1000 | DV-27       |                | X                       | X                      | X                 | X               | X                      | X               | X                              | X            | X             | X              |          |                 |
| 16         | 24115 | 1007 | DV-28       |                | X                       | X                      | X                 | X               | X                      | X               | X                              | X            | X             | X              |          |                 |
| 17         | 24115 | 0945 | DV-29       |                | X                       | X                      | X                 | X               | X                      | X               | X                              | X            | X             | X              |          |                 |
| 18         |       |      |             |                |                         |                        |                   |                 |                        |                 |                                |              |               |                |          |                 |
| 19         |       |      |             |                |                         |                        |                   |                 |                        |                 |                                |              |               |                |          |                 |
| 20         |       |      |             |                |                         |                        |                   |                 |                        |                 |                                |              |               |                |          |                 |

RELINQUISHED BY: (Signature) [Signature] DATE/TIME 2/17/14 12:30  
 RECEIVED BY (Signature) [Signature] DATE/TIME 2/15  
 LABORATORY NOTES: MSI samples RUSH DIOXINS  
Hold sample for TUP / SPL Testing  
 SAMPLE RECEIPT: TOTAL # OF CONTAINERS 17  
 COC SEALS Y / N / NA 0 / 17 / 0  
 SEALS INTACT Y / N / NA 0 / 17 / 0  
 RECEIVED TEMP: 30C  
 SAMPLE DISPOSAL INSTRUCTIONS: ESN @ \$5.00/sample or Return to Client



## NON-VOLATILE MULTI-INCREMENT PREPARATION

DATE: 2/6 and 2/7/2014

PREPARED BY: A. Crowell, N. Kippen, Z. Tullis

ESN PROJECT: D1402050037

Bureau Veritas North America, Inc., personnel collected multi-increment samples in the field for analysis. The samples were delivered to ESN's lab for preparation and analyses, according to DOH's Multi-Increment Sampling Plan (HEER TGM, Section 4.2).

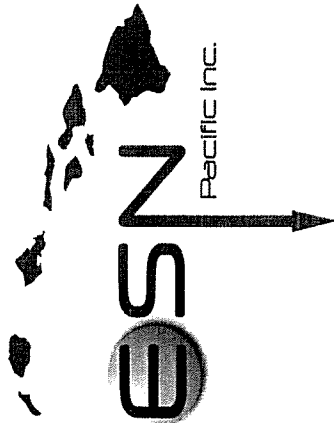
(Dry weights)

| Sample ID#  | Sample Weight (g) |
|-------------|-------------------|
| RHS-DU-1    | 1655.9            |
| RHS-DU-1.2  | 2490.1            |
| RHS-DU-1.3  | 2749.6            |
| RHS-DU-2    | 2403.1            |
| RHS-DU-3    | 1348.7            |
| RHS-DU-4    | 1361.4            |
| RHS-DU-5    | 2393.6            |
| RHS-DU-6    | 1982.6            |
| RHS-DU-7    | 2190.7            |
| RHS-DU-8    | 1738.5            |
| RHS-DU-9    | 1346.9            |
| RHS-DU-10   | 1895.3            |
| RHS-DU-11   | 1361.0            |
| RHS-DU-12   | 2363.9            |
| RHS-DU-13   | 2176.6            |
| RHS-DU-14   | 1584.5            |
| RHS-DU-15   | 1880.3            |
| RHS-DU-16   | 2119.8            |
| RHS-DU-17   | 1802.0            |
| RHS-DU-17.2 | 1897.2            |
| RHS-DU-17.3 | 2067.8            |
| RHS-DU-18   | 1770.5            |
| RHS-DU-19   | 1451.4            |
| RHS-DU-20   | 1934.5            |
| RHS-DU-21   | 1588.3            |
| RHS-DU-22   | 1857.2            |

|             |        |
|-------------|--------|
| RHS-DU-23   | 2096.1 |
| RHS-DU-24   | 1690.8 |
| RHS-DU-25   | 1412.7 |
| RHS-DU-26   | 1806.5 |
| RHS-DU-26.2 | 2268.0 |
| RHS-DU-26.3 | 2205.2 |
| RHS-DU-27   | 2204.1 |
| RHS-DU-28   | 974.8  |
| RHS-DU-29   | 1061.8 |

**SAMPLE PREP PROCESS SUMMARY:**

- Each sample was emptied from their containers on to its own clean, flat baking pan. The sample was spread out evenly across the pan no more than 1” high in thickness.
- Sample trays were placed on separate oven racks and air-dried overnight or until no significant moisture remained as determined by appearance.
- Sample trays were removed and weighed to determine sample’s total dry weight (see above table).
- Sieved entire dried sample through a #10 mesh sieve rendering particles less than two-millimeters in size. Clumps of dirt were crushed until able to pass through the sieve, rocks and other debris were set aside.
- Sieved samples were spread out evenly on flat baking pans.
- A small flat spatula was used to randomly scoop up 30 to 50 small increments to fill sample containers. Care was taken during this step to ensure that samples contained a good distribution of particle sizes. If duplicate or triplicate analysis is requested, additional samples are collected at this time. Remaining sample was returned to resealable (i.e. Zip-Loc) bags.



# Environmental Services Network

Bureau Veritas North America, Inc. PROJECT #17012.012148.48  
 Radford High School

ESN Project #D1402050037

| SAMPLE NUMBER | DATE SAMPLED | EXTRACTED DATE | ANALYZED DATE | PCB-1016 (mg/kg) | PCB-1221 (mg/kg) | PCB-1232 (mg/kg) | PCB-1242 (mg/kg) | PCB-1248 (mg/kg) | PCB-1254 (mg/kg) | PCB-1260 (mg/kg) | SURROGATE RECOVERY(%) | FLAGS |
|---------------|--------------|----------------|---------------|------------------|------------------|------------------|------------------|------------------|------------------|------------------|-----------------------|-------|
| Method Blank  |              |                |               |                  |                  |                  |                  |                  |                  |                  |                       |       |
| RHS-DU-1      | 2/4/2014     | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 91%                   |       |
| RHS-DU-1 Dup  | 2/4/2014     | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 67%                   |       |
| RHS-DU-1.2    | 2/5/2014     | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 78%                   |       |
| RHS-DU-1.3    | 2/5/2014     | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 68%                   |       |
| RHS-DU-2      | 2/4/2014     | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 0.10             | 84%                   |       |
| RHS-DU-3      | 2/4/2014     | 2/6/2014       | 2/12/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 0.21             | 88%                   |       |
| RHS-DU-4      | 2/4/2014     | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 0.05             | 87%                   |       |
| RHS-DU-5      | 2/4/2014     | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 0.09             | 74%                   |       |
| RHS-DU-6      | 2/4/2014     | 2/6/2014       | 2/7/2014      | nd               | nd               | nd               | nd               | nd               | nd               | 0.17             | 88%                   |       |
| RHS-DU-7      | 2/4/2014     | 2/6/2014       | 2/7/2014      | nd               | nd               | nd               | nd               | nd               | nd               | 0.32             | 82%                   |       |
| RHS-DU-8      | 2/4/2014     | 2/6/2014       | 2/7/2014      | nd               | nd               | nd               | nd               | nd               | nd               | 0.66             | 83%                   |       |
| RHS-DU-9      | 2/4/2014     | 2/6/2014       | 2/12/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 3.27             | 84%                   |       |
| RHS-DU-10     | 2/5/2014     | 2/6/2014       | 2/12/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 2.17             | DO                    | DF 2  |
| RHS-DU-11     | 2/5/2014     | 2/6/2014       | 2/14/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 1.00             | DO                    | DF 2  |
| RHS-DU-12     | 2/5/2014     | 2/6/2014       | 2/14/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 0.13             | 83%                   |       |
| RHS-DU-13     | 2/5/2014     | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 0.23             | 80%                   |       |
| RHS-DU-14     | 2/4/2014     | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 80%                   |       |
| RHS-DU-15     | 2/4/2014     | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 80%                   |       |
| RHS-DU-16     | 2/4/2014     | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 80%                   |       |

| MDL   | 0.10 | 0.20 | 0.20 | 0.05 | 0.05 | 0.05 | 0.05 | 0.05 | 0.05 | 0.05 |  |  |
|---|------|------|------|------|------|------|------|------|------|------|--|--|
| 2005 HI DOH EAL   | 0.05 | 0.15 | 0.10 | 0.04 | 0.02 | 0.04 | 0.04 | 0.04 | 0.04 | 0.04 |  |  |
| ACCEPTABLE RECOVERY LIMITS FOR SURROGATE (TCMX): 65% - 135% | 1.10 | 1.10 | 1.10 | 1.10 | 1.10 | 1.10 | 1.10 | 1.10 | 1.10 | 1.10 |  |  |

DF: Reported result calculated from dilution factor, multiply detection limit by dilution factor.  
 DD: Diluted out.

QA/QC DATA - LABORATORY CONTROL SPIKE ANALYSES

|                |        |
|----------------|--------|
| Spike Added    | 1.00   |
| Measured Conc. | 1.05   |
| % Recovery     | 104.9% |

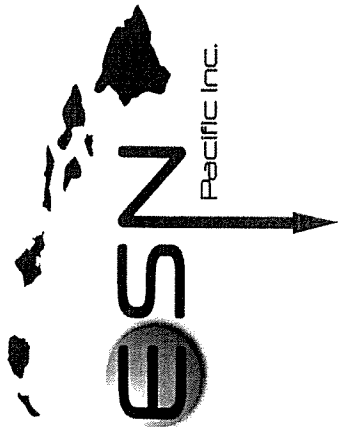
QA/QC DATA - MATRIX SPIKE ANALYSES

|                       |       |
|-----------------------|-------|
| Sample Name: RHS-DU-1 |       |
| Spike Added           | 1.00  |
| Measured Conc.        | 0.86  |
| % Recovery            | 86.3% |
| Spike Added           | 1.00  |
| Measured Conc.        | 0.85  |
| % Recovery            | 85.0% |
| RPD                   | 1.5%  |

% Recovery LIMITS: 80% TO 120%  
 RPD LIMIT: 20%

ANALYSES PERFORMED AND REVIEWED BY : K. Carvallo

2020 Kahai Street, Honolulu, Hawaii 96819



# Environmental Services Network

Bureau Veritas North America, Inc. PROJECT #17012.012148.48  
Radford High School

ESN Project #D1402050037

PCB ANALYSES OF SOILS BY EPA 8082 MODIFIED

| SAMPLE NUMBER   | DATE SAMPLED | DATE EXTRACTED | DATE ANALYZED | PCB-1016 (mg/kg) | PCB-1221 (mg/kg) | PCB-1232 (mg/kg) | PCB-1242 (mg/kg) | PCB-1248 (mg/kg) | PCB-1254 (mg/kg) | PCB-1260 (mg/kg) | SURROGATE RECOVERY(%) | FLAGS |
|-----------------|--------------|----------------|---------------|------------------|------------------|------------------|------------------|------------------|------------------|------------------|-----------------------|-------|
| Method Blank    |              | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 71%                   |       |
| RHS-DU-17       | 2/4/2014     | 2/7/2014       | 2/14/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 2.23             | DO                    | DF 2  |
| RHS-DU-17 Dup   | 2/4/2014     | 2/7/2014       | 2/14/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 2.26             | DO                    | DF 2  |
| RHS-DU-17.2     | 2/4/2014     | 2/6/2014       | 2/14/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 2.35             | DO                    | DF 2  |
| RHS-DU-17.3     | 2/4/2014     | 2/7/2014       | 2/14/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 2.24             | DO                    | DF 2  |
| RHS-DU-18       | 2/4/2014     | 2/7/2014       | 2/12/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 1.19             | 77%                   |       |
| RHS-DU-19       | 2/4/2014     | 2/7/2014       | 2/12/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 0.47             | 83%                   |       |
| RHS-DU-20       | 2/4/2014     | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 0.29             | 96%                   |       |
| RHS-DU-21       | 2/4/2014     | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 0.20             | 78%                   |       |
| RHS-DU-21 Dup   | 2/4/2014     | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 0.24             | 75%                   |       |
| RHS-DU-22       | 2/4/2014     | 2/7/2014       | 2/12/2014     | nd               | nd               | nd               | nd               | nd               | nd               | 1.46             | 86%                   |       |
| RHS-DU-23       | 2/4/2014     | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 87%                   |       |
| RHS-DU-24       | 2/4/2014     | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 82%                   |       |
| RHS-DU-25       | 2/4/2014     | 2/6/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 69%                   |       |
| RHS-DU-26       | 2/4/2014     | 2/7/2014       | 2/12/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 68%                   |       |
| RHS-DU-26.2     | 2/4/2014     | 2/7/2014       | 2/12/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 85%                   |       |
| RHS-DU-26.3     | 2/4/2014     | 2/7/2014       | 2/12/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 95%                   |       |
| RHS-DU-27       | 2/4/2014     | 2/7/2014       | 2/12/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 87%                   |       |
| RHS-DU-28       | 2/4/2014     | 2/7/2014       | 2/10/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 70%                   |       |
| RHS-DU-29       | 2/4/2014     | 2/7/2014       | 2/12/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 83%                   |       |
| RHS-DU-29 Dup   | 2/4/2014     | 2/7/2014       | 2/12/2014     | nd               | nd               | nd               | nd               | nd               | nd               | nd               | 83%                   |       |
| PQL             |              |                |               | 0.10             | 0.20             | 0.20             | 0.05             | 0.05             | 0.05             | 0.05             |                       |       |
| MDL             |              |                |               | 0.05             | 0.15             | 0.10             | 0.04             | 0.02             | 0.04             | 0.04             |                       |       |
| 2005 HI DOH EAL |              |                |               | 1.10             | 1.10             | 1.10             | 1.10             | 1.10             | 1.10             | 1.10             |                       |       |

ACCEPTABLE RECOVERY LIMITS FOR SURROGATE (TCMX): 65% - 135%

DF: Reported result calculated from dilution factor, multiply detection limit by dilution factor.  
DO: Diluted out.

QA/QC DATA - LABORATORY CONTROL SPIKE ANALYSES

|                |       |
|----------------|-------|
| Spike Added    | 1.00  |
| Measured Conc. | 0.96  |
| % Recovery     | 95.9% |

QA/QC DATA - MATRIX SPIKE ANALYSES

Sample Name: RHS-DU-21  
\*Any hits in sample spiked for MS/MSD are subtracted before reported as measured concentration.

|                |        |
|----------------|--------|
| Spike Added    | 1.00   |
| Measured Conc. | 1.02   |
| % Recovery     | 102.4% |

|                |       |
|----------------|-------|
| Spike Added    | 1.00  |
| Measured Conc. | 0.89  |
| % Recovery     | 89.2% |

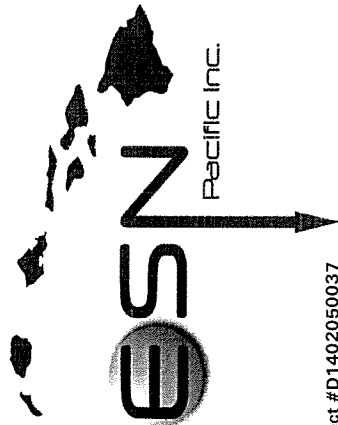
|     |       |
|-----|-------|
| RPD | 13.8% |
|-----|-------|

% Recovery LIMITS: 80% TO 120%  
RPD LIMIT: 20%

ANALYSES PERFORMED AND REVIEWED BY : K. Carvallo

2020 Kahai Street, Honolulu, Hawaii 96819





# Environmental Services Network

Bureau Veritas North America, Inc. PROJECT #17012.012148.48  
Radford High School

ESN Project #D1402050037

ORGANOCHLORINE PESTICIDES ANALYSES OF SOILS BY EPA 8081A MODIFIED

| SAMPLE NUMBER         | Method Blank | RHS-DU-1  | RHS-DU-1 Dup | RHS-DU-1.2 | RHS-DU-1.3 | RHS-DU-2  | RHS-DU-3  | RHS-DU-4  | RHS-DU-5  | RHS-DU-6  | RHS-DU-7  | RHS-DU-8  | POL     | MDL     |
|-----------------------|--------------|-----------|--------------|------------|------------|-----------|-----------|-----------|-----------|-----------|-----------|-----------|---------|---------|
| DATE SAMPLED          | 2/6/2014     | 2/4/2014  | 2/4/2014     | 2/5/2014   | 2/4/2014   | 2/4/2014  | 2/4/2014  | 2/4/2014  | 2/4/2014  | 2/4/2014  | 2/4/2014  | 2/4/2014  |         |         |
| DATE EXTRACTED        | 2/6/2014     | 2/6/2014  | 2/6/2014     | 2/6/2014   | 2/6/2014   | 2/6/2014  | 2/6/2014  | 2/6/2014  | 2/6/2014  | 2/6/2014  | 2/6/2014  | 2/6/2014  |         |         |
| DATE ANALYZED         | 2/6/2014     | 2/11/2014 | 2/11/2014    | 2/16/2014  | 2/16/2014  | 2/16/2014 | 2/16/2014 | 2/16/2014 | 2/16/2014 | 2/16/2014 | 2/16/2014 | 2/16/2014 |         |         |
| Alpha-BHC             | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.005   | 0.002   |
| Beta-BHC              | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.005   | 0.003   |
| Gamma-BHC (Lindane)   | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.005   | 0.002   |
| Delta-BHC             | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.005   | 0.002   |
| Heptachlor            | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.005   | 0.002   |
| Aldrin                | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.005   | 0.002   |
| Heptachlor epoxide    | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.005   | 0.002   |
| Gamma-Chlordane       | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.005   | 0.002   |
| Endosulfan I          | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.005   | 0.002   |
| Endosulfan II         | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.005   | 0.002   |
| Alpha-Chlordane       | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.005   | 0.002   |
| Dieldrin              | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.005   | 0.002   |
| p,p'-DDE              | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | 0.045     | 0.010   | 0.003   |
| p,p'-DDD              | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | 0.010     | 0.010   | 0.003   |
| Endosulfan II         | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.010   | 0.003   |
| p,p'-DDD              | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.010   | 0.003   |
| Endrin aldehyde       | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.010   | 0.003   |
| Endosulfan sulfate    | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.010   | 0.003   |
| Endrin ketone         | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.010   | 0.003   |
| p,p'-DDT              | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.010   | 0.006   |
| Methoxychlor          | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.010   | 0.005   |
| Chlordane (technical) | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | 0.011     | 0.010   | 0.005   |
| Toxaphene             | nd           | nd        | nd           | nd         | nd         | nd        | nd        | nd        | nd        | nd        | nd        | nd        | 0.010   | 0.005   |
|                       | (mg/kg)      | (mg/kg)   | (mg/kg)      | (mg/kg)    | (mg/kg)    | (mg/kg)   | (mg/kg)   | (mg/kg)   | (mg/kg)   | (mg/kg)   | (mg/kg)   | (mg/kg)   | (mg/kg) | (mg/kg) |

|  |            |     |     |     |     |     |     |     |     |     |     |     |
|--|------------|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|
| FLAGS  | 87%        | 80% | 84% | 86% | 88% | 86% | 90% | 88% | 87% | 96% | 89% | 87% |
| SURROGATE RECOVERY (%)                           | 87%        | 80% | 84% | 86% | 88% | 86% | 90% | 88% | 87% | 96% | 89% | 87% |
| ACCEPTABLE RECOVERY LIMITS FOR SURROGATE (TCMX): | 65% - 135% |     |     |     |     |     |     |     |     |     |     |     |

QA/QC DATA - LABORATORY CONTROL SPIKE ANALYSES

|               | Laboratory Control Spike |                        | Spike Recovery (%) |
|---------------|--------------------------|------------------------|--------------------|
|               | Spiked Conc. (mg/kg)     | Measured Conc. (mg/kg) |                    |
| Beta-BHC      | 0.050                    | 0.044                  | 87.4%              |
| p,p'-DDE      | 0.100                    | 0.085                  | 85.1%              |
| Endrin ketone | 0.100                    | 0.095                  | 94.9%              |

QA/QC DATA - MATRIX SPIKE ANALYSES

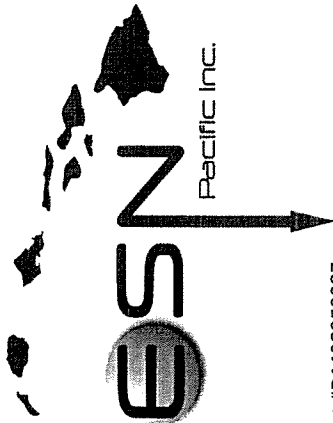
| Sample Name: RHS-DU-1 | Matrix Spike         |                        | Matrix Spike Duplicate |                        | RPD (%) | FLAGS |
|-----------------------|----------------------|------------------------|------------------------|------------------------|---------|-------|
|                       | Spiked Conc. (mg/kg) | Measured Conc. (mg/kg) | Spiked Conc. (mg/kg)   | Measured Conc. (mg/kg) |         |       |
| Beta-BHC              | 0.050                | 0.048                  | 0.050                  | 0.056                  | 112.0%  | 16.2% |
| p,p'-DDE              | 0.100                | 0.097                  | 0.100                  | 0.112                  | 111.5%  | 13.7% |
| Endrin ketone         | 0.100                | 0.095                  | 0.100                  | 0.104                  | 103.9%  | 8.9%  |

% Recovery LIMITS: 80% TO 120%

RPD LIMIT: 20%

ANALYSES PERFORMED AND REVIEWED BY : K. Carvallo

2020 Kahai Street, Honolulu, Hawaii 96819



# Environmental Services Network

Bureau Veritas North America, Inc. PROJECT #17012.012148.48  
Radford High School

ESN Project #D1402050037

ORGANOCHLORINE PESTICIDES ANALYSES OF SOILS BY EPA 8081A MODIFIED

| SAMPLE NUMBER         | Method Blank | RHS-DU-9  | RHS-DU-10 | RHS-DU-11 | RHS-DU-12 | RHS-DU-13 | RHS-DU-14 | RHS-DU-15 | RHS-DU-16 | RHS-DU-17 | RHS-DU-17 DUT | PQL     | MDL     |
|-----------------------|--------------|-----------|-----------|-----------|-----------|-----------|-----------|-----------|-----------|-----------|---------------|---------|---------|
| DATE SAMPLED          |              | 2/5/2014  | 2/5/2014  | 2/5/2014  | 2/5/2014  | 2/5/2014  | 2/4/2014  | 2/4/2014  | 2/4/2014  | 2/4/2014  | 2/4/2014      |         |         |
| DATE EXTRACTED        |              | 2/6/2014  | 2/6/2014  | 2/6/2014  | 2/6/2014  | 2/6/2014  | 2/6/2014  | 2/6/2014  | 2/6/2014  | 2/7/2014  | 2/7/2014      |         |         |
| DATE ANALYZED         |              | 2/11/2014 | 2/13/2014 | 2/13/2014 | 2/13/2014 | 2/11/2014 | 2/11/2014 | 2/11/2014 | 2/11/2014 | 2/13/2014 | 2/13/2014     |         |         |
| Alpha-BHC             | nd           | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd            | 0.005   | 0.002   |
| Beta-BHC              | nd           | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd            | 0.005   | 0.003   |
| Gamma-BHC (Lindane)   | nd           | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd            | 0.005   | 0.002   |
| Delta-BHC             | nd           | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd            | 0.005   | 0.002   |
| Heptachlor            | nd           | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd            | 0.005   | 0.002   |
| Aldrin                | nd           | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd            | 0.005   | 0.002   |
| Heptachlor epoxide    | nd           | nd        | nd        | nd        | nd        | 0.022     | nd        | nd        | nd        | nd        | nd            | 0.005   | 0.002   |
| Gamma-Chlordane       | nd           | 0.006     | 0.024     | 0.014     | 0.014     | nd        | nd        | nd        | nd        | nd        | nd            | 0.005   | 0.002   |
| Endosulfan I          | nd           | nd        | nd        | nd        | nd        | nd        | 0.006     | nd        | nd        | nd        | nd            | 0.005   | 0.002   |
| Alpha-Chlordane       | nd           | 0.006     | 0.034     | 0.014     | 0.014     | 0.014     | 0.014     | nd        | nd        | nd        | nd            | 0.010   | 0.003   |
| Dieldrin              | nd           | 0.027     | 0.034     | 0.036     | 0.036     | 0.016     | 0.012     | nd        | nd        | nd        | nd            | 0.010   | 0.003   |
| p,p'-DDE              | nd           | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd            | 0.010   | 0.003   |
| Endosulfan II         | nd           | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd            | 0.010   | 0.003   |
| p,p'-DDD              | nd           | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd            | 0.010   | 0.003   |
| Endrin aldehyde       | nd           | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd            | 0.010   | 0.003   |
| Endosulfan sulfate    | nd           | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd            | 0.010   | 0.006   |
| p,p'-DDT              | nd           | 0.021     | 0.040     | 0.174     | 0.174     | 0.051     | 0.012     | 0.021     | 0.164     | 0.164     | 0.161         | 0.010   | 0.005   |
| Endrin ketone         | nd           | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd            | 0.010   | 0.005   |
| Methoxychlor          | nd           | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd            | 0.010   | 0.005   |
| Chlordane (technical) | nd           | nd        | 0.113     | 0.127     | 0.127     | nd        | nd        | nd        | 0.050     | nd        | nd            | 0.050   | 0.020   |
| Toxaphene             | nd           | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd        | nd            | 0.050   | 0.010   |
| (mg/kg)               | (mg/kg)      | (mg/kg)   | (mg/kg)   | (mg/kg)   | (mg/kg)   | (mg/kg)   | (mg/kg)   | (mg/kg)   | (mg/kg)   | (mg/kg)   | (mg/kg)       | (mg/kg) | (mg/kg) |

| CLASS  | DF 2 | DF 2 | DF 2 | DF 2 | DF 2 | DF 2 | DF 2 | DF 2 | DF 2 | DF 2 | DF 2 | DF 2 | DF 2 |
|--|------|------|------|------|------|------|------|------|------|------|------|------|------|
| SURROGATE RECOVERY (%)                           | 101% | 79%  | 93%  | 83%  | 87%  | 84%  | 83%  | 83%  | 83%  | 83%  | 83%  | 83%  | 83%  |
| ACCEPTABLE RECOVERY LIMITS FOR SURROGATE (TCMX): | 65%  | 135% |      |      |      |      |      |      |      |      |      |      |      |

DF: Reported result calculated from dilution factor, multiply detection limit by dilution factor.  
DO: Diluted out.  
DOUT 1: RPD for duplicates outside acceptable limit (20%). Matrix may be nonhomogenous.

QA/QC DATA - LABORATORY CONTROL SPIKE ANALYSES

| Laboratory Control Spike | Spiked |         | Measured |         | Spike Recovery |       |
|--------------------------|--------|---------|----------|---------|----------------|-------|
|                          | Conc.  | (mg/kg) | Conc.    | (mg/kg) | (%)            | (%)   |
| Beta-BHC                 | 0.050  | 0.044   | 0.050    | 0.044   | 87.4%          | 87.4% |
| p,p'-DDE                 | 0.100  | 0.085   | 0.100    | 0.085   | 85.1%          | 85.1% |
| Endrin ketone            | 0.100  | 0.095   | 0.100    | 0.095   | 94.9%          | 94.9% |

% Recovery LIMITS: 80% TO 120%  
RPD LIMIT: 20%

ANALYSES PERFORMED AND REVIEWED BY : K. Carvallo



# Environmental Services Network

Bureau Veritas North America, Inc. PROJECT #17012.012148.48  
Radford High School

ESN Project #D1402050037

ORGANOCHLORINE PESTICIDES ANALYSES OF SOILS BY EPA 8081A MODIFIED

| SAMPLE NUMBER         | Method Blank | RHS-DU-17.2 | RHS-DU-17.3 | RHS-DU-18 | RHS-DU-19 | RHS-DU-20 | RHS-DU-21 | RHS-DU-21 Due | RHS-DU-22 | RHS-DU-23 | RHS-DU-24 | RHS-DU-25 | POL   | MDL   |
|-----------------------|--------------|-------------|-------------|-----------|-----------|-----------|-----------|---------------|-----------|-----------|-----------|-----------|-------|-------|
| DATE SAMPLED          | 2/6/2014     | 2/4/2014    | 2/4/2014    | 2/4/2014  | 2/4/2014  | 2/4/2014  | 2/4/2014  | 2/4/2014      | 2/4/2014  | 2/4/2014  | 2/4/2014  | 2/4/2014  |       |       |
| DATE EXTRACTED        | 2/6/2014     | 2/7/2014    | 2/7/2014    | 2/7/2014  | 2/7/2014  | 2/6/2014  | 2/6/2014  | 2/6/2014      | 2/7/2014  | 2/7/2014  | 2/6/2014  | 2/6/2014  |       |       |
| DATE ANALYZED         | 2/13/2014    | 2/13/2014   | 2/13/2014   | 2/13/2014 | 2/13/2014 | 2/11/2014 | 2/13/2014 | 2/13/2014     | 2/13/2014 | 2/13/2014 | 2/11/2014 | 2/11/2014 |       |       |
| Alpha-BHC             | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.005 | 0.002 |
| Beta-BHC              | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.005 | 0.003 |
| Gamma-BHC (Lindane)   | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.005 | 0.002 |
| Delta-BHC             | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.005 | 0.002 |
| Heptachlor            | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.005 | 0.002 |
| Aldrin                | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.005 | 0.002 |
| Heptachlor epoxide    | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.005 | 0.002 |
| Gamma-Chlordane       | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.005 | 0.002 |
| Endosulfan I          | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.005 | 0.002 |
| Alpha-Chlordane       | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.005 | 0.002 |
| Dieldrin              | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.005 | 0.002 |
| p,p'-DDE              | nd           | nd          | 0.030       | nd        | nd        | nd        | nd        | nd            | nd        | 0.011     | nd        | nd        | 0.010 | 0.003 |
| Endrin                | nd           | nd          | 0.054       | nd        | nd        | 0.013     | nd        | nd            | nd        | 0.011     | nd        | nd        | 0.010 | 0.003 |
| Endosulfan II         | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.010 | 0.003 |
| p,p'-DDD              | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.010 | 0.003 |
| Endrin aldehyde       | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.010 | 0.003 |
| Endosulfan sulfate    | nd           | 0.195       | 0.175       | 0.132     | 0.025     | 0.019     | 0.015     | 0.018         | 0.022     | 0.074     | nd        | nd        | 0.010 | 0.005 |
| p,p'-DDT              | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.010 | 0.005 |
| Endrin ketone         | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.010 | 0.005 |
| Methoxychlor          | nd           | nd          | 0.225       | nd        | nd        | nd        | nd        | nd            | nd        | 0.057     | nd        | nd        | 0.050 | 0.020 |
| Chlordane (technical) | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.050 | 0.010 |
| Toxaphene             | nd           | nd          | nd          | nd        | nd        | nd        | nd        | nd            | nd        | nd        | nd        | nd        | 0.050 | 0.010 |

| FLAGS  | DF 2 | DF 2 | 81% | 83% | 100% | 78% | 75% | 80% | 85% | 107% | 85% |
|--|------|------|-----|-----|------|-----|-----|-----|-----|------|-----|
| SURROGATE RECOVERY (%)                                     | 96%  | DO   | DO  | DO  | DO   | DO  | DO  | DO  | DO  | DO   | DO  |
| ACCEPTABLE RECOVERY LIMITS FOR SURROGATE (TCMX): 65%- 135% |      |      |     |     |      |     |     |     |     |      |     |

DF: Reported result calculated from dilution factor, multiply detection limit by dilution factor.  
DO: Diluted out.

QA/QC DATA - LABORATORY CONTROL SPIKE ANALYSES

|               | Laboratory Control Spike |              | Spike                  |              |
|---------------|--------------------------|--------------|------------------------|--------------|
|               | Spiked Conc. (mg/kg)     | Recovery (%) | Measured Conc. (mg/kg) | Recovery (%) |
| Beta-BHC      | 0.050                    | 96.6%        | 0.048                  | 96.6%        |
| p,p'-DDE      | 0.100                    | 96.7%        | 0.097                  | 96.7%        |
| Endrin ketone | 0.100                    | 99.7%        | 0.100                  | 99.7%        |

QA/QC DATA - MATRIX SPIKE ANALYSES

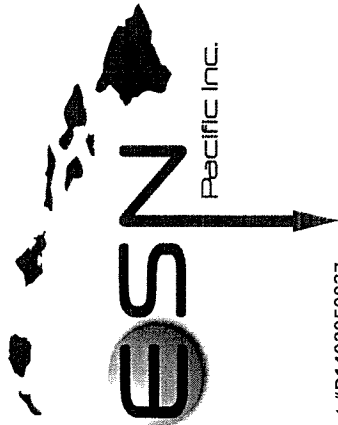
Sample Name: RHS-DU-21  
\* Any hits or matrix interference in sample spiked for MS/MSD are subtracted before reported as measured concentration.

|               | Matrix Spike         |              | Matrix Spike Duplicate |              | RPD (%) | FLAGS |
|---------------|----------------------|--------------|------------------------|--------------|---------|-------|
|               | Spiked Conc. (mg/kg) | Recovery (%) | Measured Conc. (mg/kg) | Recovery (%) |         |       |
| Beta-BHC      | 0.050                | 89.0%        | 0.045                  | 99.4%        | 11.0%   |       |
| p,p'-DDE      | 0.100                | 97.5%        | 0.098                  | 119.0%       | 19.9%   |       |
| Endrin ketone | 0.100                | 118.1%       | 0.118                  | 109.9%       | 7.2%    |       |

% Recovery LIMITS: 80% TO 120%  
RPD LIMIT: 20%

ANALYSES PERFORMED AND REVIEWED BY : K. Carvallo

2020 Kahai Street, Honolulu, Hawaii 96819



# Environmental Services Network

Bureau Veritas North America, Inc. PROJECT #17012.012148.48  
Radford High School

ESN Project #D1402050037

ORGANOCHLORINE PESTICIDES ANALYSES OF SOILS BY EPA 8081A, MODIFIED

| SAMPLE NUMBER         | Method Blank | RHS-DU-26 | RHS-DU-26.2 | RHS-DU-26.3 | RHS-DU-27 | RHS-DU-28 | RHS-DU-29 | RHS-DU-29 DUT | PQL     | MDL     |
|-----------------------|--------------|-----------|-------------|-------------|-----------|-----------|-----------|---------------|---------|---------|
| DATE SAMPLED          | 2/4/2014     | 2/4/2014  | 2/4/2014    | 2/4/2014    | 2/4/2014  | 2/4/2014  | 2/4/2014  | 2/4/2014      |         |         |
| DATE EXTRACTED        | 2/7/2014     | 2/7/2014  | 2/7/2014    | 2/7/2014    | 2/7/2014  | 2/7/2014  | 2/7/2014  | 2/7/2014      |         |         |
| DATE ANALYZED         | 2/13/2014    | 2/13/2014 | 2/13/2014   | 2/13/2014   | 2/11/2014 | 2/11/2014 | 2/11/2014 | 2/11/2014     |         |         |
| Alpha-BHC             | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.005   | 0.002   |
| Beta-BHC              | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.005   | 0.003   |
| Gamma-BHC (Lindane)   | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.005   | 0.002   |
| Delta-BHC             | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.005   | 0.002   |
| Heptachlor            | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.005   | 0.002   |
| Aldrin                | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.005   | 0.002   |
| Heptachlor epoxide    | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.005   | 0.002   |
| Gamma-Chlordane       | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.005   | 0.002   |
| Endosulfan I          | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.005   | 0.002   |
| Alpha-Chlordane       | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.005   | 0.002   |
| Dieldrin              | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.005   | 0.003   |
| p,p'-DDE              | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.010   | 0.003   |
| Endrin                | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.010   | 0.005   |
| Endosulfan II         | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.010   | 0.003   |
| p,p'-DDD              | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.010   | 0.006   |
| Endrin aldehyde       | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.010   | 0.005   |
| Endosulfan sulfate    | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.010   | 0.005   |
| p,p'-DDT              | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.010   | 0.005   |
| Endrin ketone         | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.010   | 0.009   |
| Methoxychlor          | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.050   | 0.020   |
| Chlordane (technical) | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | 0.050   | 0.010   |
| Toxaphene             | nd           | nd        | nd          | nd          | nd        | nd        | nd        | nd            | (mg/kg) | (mg/kg) |

| FLAGS  | 96% | 95%  | 76% | 94% | 96% | 94% | 85% |
|--|-----|------|-----|-----|-----|-----|-----|
| SURROGATE RECOVERY (%)                           | 96% | 95%  | 76% | 94% | 96% | 94% | 85% |
| ACCEPTABLE RECOVERY LIMITS FOR SURROGATE (TCMX): | 65% | 135% |     |     |     |     |     |

QA/QC DATA - LABORATORY CONTROL SPIKE ANALYSES

|               | Laboratory Control Spike |                       |
|---------------|--------------------------|-----------------------|
|               | Spiked Conc. (mg/kg)     | Measured Recovery (%) |
| Beta-BHC      | 0.050                    | 96.6%                 |
| p,p'-DDE      | 0.100                    | 96.7%                 |
| Endrin ketone | 0.100                    | 99.7%                 |

% Recovery LIMITS: 80% TO 120%  
RPD LIMIT: 20%

ANALYSES PERFORMED AND REVIEWED BY : K. Carvallo

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**Analysis of Diesel Range Organics & Lube Oil Range Organics in Soil  
 by Method NWTPH-Dx Extended (8015M)**

| Sample Number       | Date Prepared | Date Analyzed | Surrogate Recovery (%) | Diesel Range Organics (mg/kg) | Lube Oil Range Organics (mg/kg) |
|---------------------|---------------|---------------|------------------------|-------------------------------|---------------------------------|
| Method Blank        | 2/10/2014     | 2/11/2014     | 112                    | nd                            | nd                              |
| LCS                 | 2/10/2014     | 2/11/2014     | 100                    | 107%                          | ---                             |
| RHS-DU-1            | 2/10/2014     | 2/11/2014     | 109                    | nd                            | nd                              |
| RHS-DU-1.2          | 2/10/2014     | 2/11/2014     | 105                    | nd                            | nd                              |
| RHS-DU-1.3          | 2/10/2014     | 2/11/2014     | 104                    | nd                            | nd                              |
| RHS-DU-2            | 2/10/2014     | 2/11/2014     | 98                     | nd                            | nd                              |
| RHS-DU-3            | 2/10/2014     | 2/11/2014     | 96                     | nd                            | nd                              |
| RHS-DU-4            | 2/10/2014     | 2/11/2014     | 104                    | nd                            | nd                              |
| RHS-DU-5            | 2/10/2014     | 2/11/2014     | 106                    | nd                            | nd                              |
| RHS-DU-6            | 2/10/2014     | 2/11/2014     | 109                    | nd                            | nd                              |
| RHS-DU-7            | 2/10/2014     | 2/11/2014     | 77                     | nd                            | nd                              |
| RHS-DU-8            | 2/10/2014     | 2/11/2014     | 89                     | nd                            | nd                              |
| RHS-DU-9            | 2/10/2014     | 2/11/2014     | 114                    | nd                            | nd                              |
| RHS-DU-10           | 2/10/2014     | 2/11/2014     | 78                     | nd                            | nd                              |
| RHS-DU-11           | 2/10/2014     | 2/11/2014     | 115                    | nd                            | nd                              |
| RHS-DU-12           | 2/10/2014     | 2/11/2014     | 111                    | nd                            | nd                              |
| RHS-DU-13           | 2/12/2014     | 2/13/2014     | 89                     | nd                            | nd                              |
| RHS-DU-13 Duplicate | 2/12/2014     | 2/13/2014     | 83                     | nd                            | nd                              |
| RHS-DU-14           | 2/12/2014     | 2/13/2014     | 90                     | nd                            | nd                              |
| RHS-DU-15           | 2/12/2014     | 2/13/2014     | 82                     | nd                            | nd                              |
| RHS-DU-16           | 2/12/2014     | 2/13/2014     | 76                     | nd                            | nd                              |
| RHS-DU-16 Duplicate | 2/12/2014     | 2/13/2014     | 84                     | nd                            | nd                              |
| RHS-DU-17           | 2/10/2014     | 2/11/2014     | 109                    | nd                            | nd                              |
| RHS-DU-17.2         | 2/10/2014     | 2/11/2014     | 100                    | nd                            | nd                              |
| RHS-DU-17.3         | 2/10/2014     | 2/11/2014     | 104                    | nd                            | nd                              |
| RHS-DU-18           | 2/10/2014     | 2/11/2014     | 90                     | nd                            | nd                              |
| RHS-DU-19           | 2/10/2014     | 2/11/2014     | 78                     | nd                            | nd                              |
| RHS-DU-20           | 2/10/2014     | 2/11/2014     | 96                     | nd                            | nd                              |
| RHS-DU-21           | 2/10/2014     | 2/11/2014     | 91                     | nd                            | nd                              |
| RHS-DU-22           | 2/11/2014     | 2/12/2014     | 89                     | nd                            | nd                              |
| RHS-DU-23           | 2/11/2014     | 2/12/2014     | 87                     | nd                            | nd                              |
| RHS-DU-24           | 2/11/2014     | 2/12/2014     | 93                     | nd                            | nd                              |
| RHS-DU-25           | 2/11/2014     | 2/12/2014     | 90                     | nd                            | nd                              |
| RHS-DU-26           | 2/11/2014     | 2/12/2014     | 93                     | nd                            | nd                              |
| RHS-DU-26.2         | 2/11/2014     | 2/12/2014     | 92                     | nd                            | nd                              |
| RHS-DU-26.3         | 2/12/2014     | 2/13/2014     | 77                     | nd                            | nd                              |
| RHS-DU-27           | 2/12/2014     | 2/13/2014     | 80                     | nd                            | nd                              |
| RHS-DU-28           | 2/12/2014     | 2/13/2014     | 78                     | nd                            | nd                              |
| RHS-DU-29           | 2/12/2014     | 2/13/2014     | 77                     | nd                            | nd                              |
| RHS-DU-29 Duplicate | 2/12/2014     | 2/13/2014     | 77                     | nd                            | nd                              |
| Reporting Limits    |               |               |                        | 50                            | 100                             |

"nd" Indicates not detected at the listed detection limits.

"---" Indicates analyte not tested.

ACCEPTABLE RECOVERY LIMITS FOR SURROGATE : 50% TO 150%

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**Analysis of Semivolatile Organic Compounds in Soil by Method 8270**

Analytical Results:

|                               |           | MTH BLK  | LCS      | RHS-DU-1 | RHS-DU-1.2 | RHS-DU-1.3 | RHS-DU-2 | RHS-DU-3 |
|-------------------------------|-----------|----------|----------|----------|------------|------------|----------|----------|
| Date extracted                | Reporting | 02/10/14 | 02/10/14 | 02/10/14 | 02/10/14   | 02/10/14   | 02/10/14 | 02/10/14 |
| Date analyzed                 | Limits    | 02/11/14 | 02/11/14 | 02/11/14 | 02/11/14   | 02/11/14   | 02/11/14 | 02/11/14 |
| Moisture, %                   | (mg/kg)   |          |          | 6%       | 9%         | 12%        | 12%      | 12%      |
| Pyridine                      | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Aniline                       | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Phenol                        | 0.41*     | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 2-Chlorophenol                | 0.36*     | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Bis (2-chloroethyl) ether     | 0.40*     | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 1,3-Dichlorobenzene           | 0.34*     | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 1,4-Dichlorobenzene           | 0.26*     | nd       | 150%     | nd       | nd         | nd         | nd       | nd       |
| 1,2-Dichlorobenzene           | 0.38*     | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Benzyl alcohol                | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Hexachlorethane               | 0.39*     | nd       |          | nd       | nd         | nd         | nd       | nd       |
| N-Nitroso-di-n-propylamine    | 1.0       | nd       | 131%     | nd       | nd         | nd         | nd       | nd       |
| 3,4-Methylphenol (m,p-cresol) | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 2-Methylphenol (o-cresol)     | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Bis (2-chloroisopropyl) ether | 0.42*     | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Nitrobenzene                  | 0.41*     | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Isophorone                    | 0.41*     | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 2-Nitrophenol                 | 5.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 2,4-Dimethylphenol            | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Bis (2-chloroethoxy) methane  | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 2,4-Dichlorophenol            | 0.36*     | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 1,2,4-Trichlorobenzene        | 0.40*     | nd       | 141%     | nd       | nd         | nd         | nd       | nd       |
| Naphthalene                   | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 4-Chloroaniline               | 0.46*     | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Hexachlorobutadiene           | 0.43*     | nd       | 149%     | nd       | nd         | nd         | nd       | nd       |
| 4-Chloro-3-methylphenol       | 5.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 2-Methylnaphthalene           | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 1-Methylnaphthalene           | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Hexachlorocyclopentadiene     | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 2,4,6-Trichlorophenol         | 5.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 2,4,5-Trichlorophenol         | 5.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 2-Chloronaphthalene           | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 2-Nitroaniline                | 5.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 1,4-Dinitrobenzene            | 5.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Acenaphthylene                | 0.1       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 1,3-Dinitrobenzene            | 0.28*     | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Dimethylphthalate             | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 2,6-Dinitrotoluene            | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 1,2-Dinitrobenzene            | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Acenaphthene                  | 0.1       | nd       | 145%     | nd       | nd         | nd         | nd       | nd       |
| 3-Nitroaniline                | 5.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 2,4-Dinitrophenol             | 1.0*      | nd       | 51%      | nd       | nd         | nd         | nd       | nd       |
| 2,4-Dinitrotoluene            | 0.35*     | nd       | 145%     | nd       | nd         | nd         | nd       | nd       |
| 4-Nitrophenol                 | 5.0       | nd       | 61%      | nd       | nd         | nd         | nd       | nd       |
| Dibenzofuran                  | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 2,3,4,6-Tetrachlorophenol     | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 2,3,5,6-Tetrachlorophenol     | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Fluorene                      | 0.1       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 4-Chlorophenylphenylether     | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Diethylphthalate              | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 4-Nitroaniline                | 5.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 4,6-Dinitro-2-methylphenol    | 5.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| N-nitrosodiphenylamine        | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Azobenzene                    | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| 4-Bromophenylphenylether      | 1.0       | nd       |          | nd       | nd         | nd         | nd       | nd       |
| Hexachlorobenzene             | 0.45*     | nd       |          | nd       | nd         | nd         | nd       | nd       |



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Analysis of Semivolatile Organic Compounds in Soil by Method 8270

Analytical Results:

|                              |           | MTH BLK  | LCS      | RHS-DU-1 | RHS-DU-1.2 | RHS-DU-1.3 | RHS-DU-2    | RHS-DU-3    |
|------------------------------|-----------|----------|----------|----------|------------|------------|-------------|-------------|
| Date extracted               | Reporting | 02/10/14 | 02/10/14 | 02/10/14 | 02/10/14   | 02/10/14   | 02/10/14    | 02/10/14    |
| Date analyzed                | Limits    | 02/11/14 | 02/11/14 | 02/11/14 | 02/11/14   | 02/11/14   | 02/11/14    | 02/11/14    |
| Moisture, %                  | (mg/kg)   |          |          | 6%       | 9%         | 12%        | 12%         | 12%         |
| Pentachlorophenol            | 1.0*      | nd       | 72%      | nd       | nd         | nd         | nd          | nd          |
| Phenanthrene                 | 0.1       | nd       |          | nd       | nd         | nd         | nd          | nd          |
| Anthracene                   | 0.1       | nd       |          | nd       | nd         | nd         | nd          | nd          |
| Carbazole                    | 1.0       | nd       |          | nd       | nd         | nd         | nd          | nd          |
| Di-n-butylphthalate          | 1.0       | nd       |          | nd       | nd         | nd         | nd          | nd          |
| Fluoranthene                 | 0.1       | nd       | 128%     | nd       | nd         | nd         | <b>0.12</b> | <b>0.12</b> |
| Pyrene                       | 0.1       | nd       | 133%     | nd       | nd         | nd         | <b>0.15</b> | <b>0.14</b> |
| Butylbenzylphthalate         | 1.0       | nd       |          | nd       | nd         | nd         | nd          | nd          |
| Bis(2-ethylhexyl) adipate    | 1.0       | nd       |          | nd       | nd         | nd         | nd          | nd          |
| Benzo(a)anthracene           | 0.1       | nd       |          | nd       | nd         | nd         | nd          | nd          |
| Chrysene                     | 0.1       | nd       |          | nd       | nd         | nd         | nd          | <b>0.13</b> |
| Bis (2-ethylhexyl) phthalate | 1.0       | nd       |          | nd       | nd         | nd         | nd          | nd          |
| Di-n-octyl phthalate         | 1.0       | nd       | 138%     | nd       | nd         | nd         | nd          | nd          |
| Benzo(b)fluoranthene         | 0.1       | nd       |          | nd       | nd         | nd         | nd          | nd          |
| Benzo(k)fluoranthene         | 0.1       | nd       |          | nd       | nd         | nd         | nd          | nd          |
| Benzo(a)pyrene               | 0.1       | nd       | 69%      | nd       | nd         | nd         | nd          | nd          |
| Dibenzo(a,h)anthracene       | 0.1       | nd       |          | nd       | nd         | nd         | nd          | nd          |
| Benzo(ghi)perylene           | 0.1       | nd       |          | nd       | nd         | nd         | nd          | nd          |
| Indeno(1,2,3-cd)pyrene       | 0.1       | nd       |          | nd       | nd         | nd         | nd          | nd          |

Surrogate recoveries

|                      |      |      |      |      |      |      |      |
|----------------------|------|------|------|------|------|------|------|
| 2-Fluorophenol       | 132% | 141% | 137% | 127% | 130% | 121% | 129% |
| Phenol-d6            | 133% | 143% | 130% | 111% | 121% | 106% | 115% |
| Nitrobenzene-d5      | 120% | 152% | 114% | 111% | 116% | 109% | 117% |
| 2-Fluorobiphenyl     | 141% | 145% | 143% | 141% | 138% | 139% | 143% |
| 2,4,6-Tribromophenol | 49%  | 123% | 57%  | 54%  | 65%  | 70%  | 74%  |
| 4-Terphenyl-d14      | 115% | 123% | 112% | 108% | 106% | 101% | 99%  |

Data Qualifiers and Analytical Comments

nd - not detected at listed reporting limits

\* - method detection limit

Acceptable Recovery limits:

2-Fluorophenol: 10-135 %

Phenol - d5: 10-135 %

Nitrobenzene - d5: 20-120 %

2-Fluorobiphenyl: 50-150%

2,4,6- Tribromophenol: 29-159%

p-Terphenyl-d14: 50-150%

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**Analysis of Semivolatile Organic Compounds in Soil by Method 8270**

Analytical Results:

|                               |           | RHS-DU-4 | RHS-DU-5 | RHS-DU-6 | RHS-DU-7 | RHS-DU-8 | RHS-DU-9 | RHS-DU-10 |
|-------------------------------|-----------|----------|----------|----------|----------|----------|----------|-----------|
| Date extracted                | Reporting | 02/10/14 | 02/10/14 | 02/10/14 | 02/10/14 | 02/10/14 | 02/10/14 | 02/10/14  |
| Date analyzed                 | Limits    | 02/11/14 | 02/11/14 | 02/11/14 | 02/11/14 | 02/11/14 | 02/11/14 | 02/11/14  |
| Moisture, %                   | (mg/kg)   | 18%      | 14%      | 16%      | 22%      | 15%      | 9%       | 10%       |
| Pyridine                      | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Aniline                       | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Phenol                        | 0.41*     | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 2-Chlorophenol                | 0.36*     | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Bis (2-chloroethyl) ether     | 0.40*     | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 1,3-Dichlorobenzene           | 0.34*     | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 1,4-Dichlorobenzene           | 0.26*     | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 1,2-Dichlorobenzene           | 0.38*     | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Benzyl alcohol                | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Hexachlorethane               | 0.39*     | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| N-Nitroso-di-n-propylamine    | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 3,4-Methylphenol (m,p-cresol) | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 2-Methylphenol (o-cresol)     | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Bis (2-chloroisopropyl) ether | 0.42*     | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Nitrobenzene                  | 0.41*     | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Isophorone                    | 0.41*     | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 2-Nitrophenol                 | 5.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 2,4-Dimethylphenol            | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Bis (2-chloroethoxy) methane  | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 2,4-Dichlorophenol            | 0.36*     | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 1,2,4-Trichlorobenzene        | 0.40*     | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Naphthalene                   | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 4-Chloroaniline               | 0.46*     | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Hexachlorobutadiene           | 0.43*     | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 4-Chloro-3-methylphenol       | 5.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 2-Methylnaphthalene           | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 1-Methylnaphthalene           | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Hexachlorocyclopentadiene     | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 2,4,6-Trichlorophenol         | 5.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 2,4,5-Trichlorophenol         | 5.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 2-Chloronaphthalene           | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 2-Nitroaniline                | 5.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 1,4-Dinitrobenzene            | 5.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Acenaphthylene                | 0.1       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 1,3-Dinitrobenzene            | 0.28*     | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Dimethylphthalate             | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 2,6-Dinitrotoluene            | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 1,2-Dinitrobenzene            | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Acenaphthene                  | 0.1       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 3-Nitroaniline                | 5.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 2,4-Dinitrophenol             | 1.0*      | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 2,4-Dinitrotoluene            | 0.35*     | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 4-Nitrophenol                 | 5.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Dibenzofuran                  | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 2,3,4,6-Tetrachlorophenol     | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 2,3,5,6-Tetrachlorophenol     | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Fluorene                      | 0.1       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 4-Chlorophenylphenylether     | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Diethylphthalate              | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 4-Nitroaniline                | 5.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 4,6-Dinitro-2-methylphenol    | 5.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| N-nitrosodiphenylamine        | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Azobenzene                    | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| 4-Bromophenylphenylether      | 1.0       | nd       | nd       | nd       | nd       | nd       | nd       | nd        |
| Hexachlorobenzene             | 0.45*     | nd       | nd       | nd       | nd       | nd       | nd       | nd        |

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Analysis of Semivolatile Organic Compounds in Soil by Method 8270

Analytical Results:

|                              |           | RHS-DU-4    | RHS-DU-5 | RHS-DU-6    | RHS-DU-7    | RHS-DU-8    | RHS-DU-9    | RHS-DU-10   |
|------------------------------|-----------|-------------|----------|-------------|-------------|-------------|-------------|-------------|
| Date extracted               | Reporting | 02/10/14    | 02/10/14 | 02/10/14    | 02/10/14    | 02/10/14    | 02/10/14    | 02/10/14    |
| Date analyzed                | Limits    | 02/11/14    | 02/11/14 | 02/11/14    | 02/11/14    | 02/11/14    | 02/11/14    | 02/11/14    |
| Moisture, %                  | (mg/kg)   | 18%         | 14%      | 16%         | 22%         | 15%         | 9%          | 10%         |
| Pentachlorophenol            | 1.0*      | nd          | nd       | nd          | nd          | nd          | nd          | nd          |
| Phenanthrene                 | 0.1       | nd          | nd       | nd          | nd          | <b>0.12</b> | <b>0.13</b> | nd          |
| Anthracene                   | 0.1       | nd          | nd       | nd          | nd          | nd          | nd          | nd          |
| Carbazole                    | 1.0       | nd          | nd       | nd          | nd          | nd          | nd          | nd          |
| Di-n-butylphthalate          | 1.0       | nd          | nd       | nd          | nd          | nd          | nd          | nd          |
| Fluoranthene                 | 0.1       | nd          | nd       | <b>0.17</b> | <b>0.21</b> | <b>0.43</b> | <b>0.50</b> | <b>0.16</b> |
| Pyrene                       | 0.1       | <b>0.10</b> | nd       | <b>0.17</b> | <b>0.23</b> | <b>0.42</b> | <b>0.50</b> | <b>0.18</b> |
| Butylbenzylphthalate         | 1.0       | nd          | nd       | nd          | nd          | nd          | nd          | nd          |
| Bis(2-ethylhexyl) adipate    | 1.0       | nd          | nd       | nd          | nd          | nd          | nd          | nd          |
| Benzo(a)anthracene           | 0.1       | nd          | nd       | nd          | nd          | <b>0.11</b> | <b>0.26</b> | nd          |
| Chrysene                     | 0.1       | <b>0.12</b> | nd       | <b>0.20</b> | <b>0.19</b> | <b>0.34</b> | <b>0.58</b> | <b>0.14</b> |
| Bis (2-ethylhexyl) phthalate | 1.0       | nd          | nd       | nd          | nd          | nd          | nd          | nd          |
| Di-n-octyl phthalate         | 1.0       | nd          | nd       | nd          | nd          | nd          | nd          | nd          |
| Benzo(b)fluoranthene         | 0.1       | nd          | nd       | nd          | nd          | <b>0.12</b> | <b>0.33</b> | nd          |
| Benzo(k)fluoranthene         | 0.1       | nd          | nd       | nd          | nd          | <b>0.20</b> | <b>0.35</b> | nd          |
| Benzo(a)pyrene               | 0.1       | nd          | nd       | nd          | nd          | nd          | <b>0.17</b> | nd          |
| Dibenzo(a,h)anthracene       | 0.1       | nd          | nd       | nd          | nd          | nd          | nd          | nd          |
| Benzo(ghi)perylene           | 0.1       | nd          | nd       | nd          | nd          | nd          | <b>0.35</b> | nd          |
| Indeno(1,2,3-cd)pyrene       | 0.1       | nd          | nd       | nd          | nd          | nd          | <b>0.35</b> | nd          |
| <b>Surrogate recoveries</b>  |           |             |          |             |             |             |             |             |
| 2-Fluorophenol               |           | 135%        | 128%     | 129%        | 142%        | **156%      | 135%        | 138%        |
| Phenol-d6                    |           | 123%        | 118%     | 119%        | 138%        | 149%        | 132%        | 141%        |
| Nitrobenzene-d5              |           | 116%        | 112%     | 121%        | 61%         | 68%         | 114%        | 63%         |
| 2-Fluorobiphenyl             |           | 143%        | 140%     | 147%        | 78%         | 86%         | 150%        | 75%         |
| 2,4,6-Tribromophenol         |           | 78%         | 69%      | 63%         | 85%         | 94%         | 91%         | 92%         |
| 4-Terphenyl-d14              |           | 107%        | 109%     | 111%        | 59%         | 68%         | 117%        | 60%         |

Data Qualifiers and Analytical Comments

nd - not detected at listed reporting limits

\* - method detection limit

\*\* - The recovery for 2-Fluorophenol exceeded the control limits. No further action taken.

Acceptable Recovery limits:

2-Fluorophenol: 10-135 %

Phenol - d5: 10-135 %

Nitrobenzene - d5: 20-120 %

2-Fluorobiphenyl: 50-150%

2,4,6- Tribromophenol: 29-159%

p-Terphenyl-d14: 50-150%

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**Analysis of Semivolatile Organic Compounds in Soil by Method 8270**

Analytical Results:

|                               |           | RHS-DU-11 | RHS-DU-12 | RHS-DU-13 | RHS-DU-14 | RHS-DU-15 | RHS-DU-16 | RHS-DU-17 |
|-------------------------------|-----------|-----------|-----------|-----------|-----------|-----------|-----------|-----------|
| Date extracted                | Reporting | 02/10/14  | 02/10/14  | 02/12/14  | 02/12/14  | 02/12/14  | 02/12/14  | 02/11/14  |
| Date analyzed                 | Limits    | 02/11/14  | 02/11/14  | 02/13/14  | 02/13/14  | 02/13/14  | 02/13/14  | 02/11/14  |
| Moisture, %                   | (mg/kg)   | 15%       | 19%       | 12%       | 20%       | 14%       | 10%       | 13%       |
| Pyridine                      | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Aniline                       | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Phenol                        | 0.41*     | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 2-Chlorophenol                | 0.36*     | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Bis (2-chloroethyl) ether     | 0.40*     | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 1,3-Dichlorobenzene           | 0.34*     | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 1,4-Dichlorobenzene           | 0.26*     | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 1,2-Dichlorobenzene           | 0.38*     | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Benzyl alcohol                | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Hexachlorethane               | 0.39*     | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| N-Nitroso-di-n-propylamine    | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 3,4-Methylphenol (m,p-cresol) | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 2-Methylphenol (o-cresol)     | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Bis (2-chloroisopropyl) ether | 0.42*     | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Nitrobenzene                  | 0.41*     | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Isophorone                    | 0.41*     | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 2-Nitrophenol                 | 5.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 2,4-Dimethylphenol            | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Bis (2-chloroethoxy) methane  | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 2,4-Dichlorophenol            | 0.36*     | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 1,2,4-Trichlorobenzene        | 0.40*     | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Naphthalene                   | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 4-Chloroaniline               | 0.46*     | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Hexachlorobutadiene           | 0.43*     | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 4-Chloro-3-methylphenol       | 5.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 2-Methylnaphthalene           | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 1-Methylnaphthalene           | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Hexachlorocyclopentadiene     | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 2,4,6-Trichlorophenol         | 5.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 2,4,5-Trichlorophenol         | 5.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 2-Chloronaphthalene           | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 2-Nitroaniline                | 5.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 1,4-Dinitrobenzene            | 5.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Acenaphthylene                | 0.1       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 1,3-Dinitrobenzene            | 0.28*     | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Dimethylphthalate             | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 2,6-Dinitrotoluene            | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 1,2-Dinitrobenzene            | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Acenaphthene                  | 0.1       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 3-Nitroaniline                | 5.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 2,4-Dinitrophenol             | 1.0*      | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 2,4-Dinitrotoluene            | 0.35*     | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 4-Nitrophenol                 | 5.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Dibenzofuran                  | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 2,3,4,6-Tetrachlorophenol     | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 2,3,5,6-Tetrachlorophenol     | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Fluorene                      | 0.1       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 4-Chlorophenylphenylether     | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Diethylphthalate              | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 4-Nitroaniline                | 5.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 4,6-Dinitro-2-methylphenol    | 5.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| N-nitrosodiphenylamine        | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Azobenzene                    | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| 4-Bromophenylphenylether      | 1.0       | nd        | nd        | nd        | nd        | nd        | nd        | nd        |
| Hexachlorobenzene             | 0.45*     | nd        | nd        | nd        | nd        | nd        | nd        | nd        |

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Analysis of Semivolatile Organic Compounds in Soil by Method 8270

Analytical Results:

|                              |           | RHS-DU-11   | RHS-DU-12   | RHS-DU-13 | RHS-DU-14 | RHS-DU-15 | RHS-DU-16   | RHS-DU-17   |
|------------------------------|-----------|-------------|-------------|-----------|-----------|-----------|-------------|-------------|
| Date extracted               | Reporting | 02/10/14    | 02/10/14    | 02/12/14  | 02/12/14  | 02/12/14  | 02/12/14    | 02/11/14    |
| Date analyzed                | Limits    | 02/11/14    | 02/11/14    | 02/13/14  | 02/13/14  | 02/13/14  | 02/13/14    | 02/11/14    |
| Moisture, %                  | (mg/kg)   | 15%         | 19%         | 12%       | 20%       | 14%       | 10%         | 13%         |
| Pentachlorophenol            | 1.0*      | nd          | nd          | nd        | nd        | nd        | nd          | nd          |
| Phenanthrene                 | 0.1       | nd          | nd          | nd        | nd        | nd        | nd          | nd          |
| Anthracene                   | 0.1       | nd          | nd          | nd        | nd        | nd        | nd          | nd          |
| Carbazole                    | 1.0       | nd          | nd          | nd        | nd        | nd        | nd          | nd          |
| Di-n-butylphthalate          | 1.0       | nd          | nd          | nd        | nd        | nd        | nd          | nd          |
| Fluoranthene                 | 0.1       | <b>0.18</b> | <b>0.11</b> | nd        | nd        | nd        | <b>0.10</b> | <b>0.13</b> |
| Pyrene                       | 0.1       | <b>0.23</b> | <b>0.14</b> | nd        | nd        | nd        | <b>0.10</b> | <b>0.17</b> |
| Butylbenzylphthalate         | 1.0       | nd          | nd          | nd        | nd        | nd        | nd          | nd          |
| Bis(2-ethylhexyl) adipate    | 1.0       | nd          | nd          | nd        | nd        | nd        | nd          | nd          |
| Benzo(a)anthracene           | 0.1       | nd          | nd          | nd        | nd        | nd        | nd          | nd          |
| Chrysene                     | 0.1       | <b>0.20</b> | <b>0.13</b> | nd        | nd        | nd        | <b>0.10</b> | <b>0.16</b> |
| Bis (2-ethylhexyl) phthalate | 1.0       | nd          | nd          | nd        | nd        | nd        | nd          | nd          |
| Di-n-octyl phthalate         | 1.0       | nd          | nd          | nd        | nd        | nd        | nd          | nd          |
| Benzo(b)fluoranthene         | 0.1       | nd          | nd          | nd        | nd        | nd        | nd          | nd          |
| Benzo(k)fluoranthene         | 0.1       | nd          | nd          | nd        | nd        | nd        | nd          | nd          |
| Benzo(a)pyrene               | 0.1       | nd          | nd          | nd        | nd        | nd        | nd          | nd          |
| Dibenzo(a,h)anthracene       | 0.1       | nd          | nd          | nd        | nd        | nd        | nd          | nd          |
| Benzo(ghi)perylene           | 0.1       | nd          | nd          | nd        | nd        | nd        | nd          | nd          |
| Indeno(1,2,3-cd)pyrene       | 0.1       | nd          | nd          | nd        | nd        | nd        | nd          | nd          |

Surrogate recoveries

|                      |      |      |      |      |      |      |      |
|----------------------|------|------|------|------|------|------|------|
| 2-Fluorophenol       | 141% | 147% | 120% | 112% | 114% | 107% | 110% |
| Phenol-d6            | 140% | 147% | 114% | 92%  | 89%  | 76%  | 101% |
| Nitrobenzene-d5      | 115% | 122% | 111% | 33%  | 24%  | 91%  | 111% |
| 2-Fluorobiphenyl     | 142% | 149% | 140% | 115% | 121% | 121% | 134% |
| 2,4,6-Tribromophenol | 91%  | 97%  | 63%  | 53%  | 33%  | 28%  | 63%  |
| 4-Terphenyl-d14      | 118% | 113% | 107% | 93%  | 84%  | 78%  | 112% |

Data Qualifiers and Analytical Comments

nd - not detected at listed reporting limits

\* - method detection limit

Acceptable Recovery limits:

2-Fluorophenol: 10-135 %

Phenol - d5: 10-135 %

Nitrobenzene - d5: 20-120 %

2-Fluorobiphenyl: 50-150%

2,4,6- Tribromophenol: 29-159%

p-Terphenyl-d14: 50-150%

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Analysis of Semivolatile Organic Compounds in Soil by Method 8270

Analytical Results:

|                               |           | RHS-DU-17.2 | RHS-DU-17.3 | RHS-DU-18 | RHS-DU-19 | RHS-DU-20 | RHS-DU-21 |
|-------------------------------|-----------|-------------|-------------|-----------|-----------|-----------|-----------|
| Date extracted                | Reporting | 02/11/14    | 02/11/14    | 02/11/14  | 02/11/14  | 02/11/14  | 02/11/14  |
| Date analyzed                 | Limits    | 02/11/14    | 02/11/14    | 02/11/14  | 02/11/14  | 02/11/14  | 02/12/14  |
| Moisture, %                   | (mg/kg)   | 11%         | 11%         | 11%       | 9%        | 18%       | 14%       |
| Pyridine                      | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| Aniline                       | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| Phenol                        | 0.41*     | nd          | nd          | nd        | nd        | nd        | nd        |
| 2-Chlorophenol                | 0.36*     | nd          | nd          | nd        | nd        | nd        | nd        |
| Bis (2-chloroethyl) ether     | 0.40*     | nd          | nd          | nd        | nd        | nd        | nd        |
| 1,3-Dichlorobenzene           | 0.34*     | nd          | nd          | nd        | nd        | nd        | nd        |
| 1,4-Dichlorobenzene           | 0.26*     | nd          | nd          | nd        | nd        | nd        | nd        |
| 1,2-Dichlorobenzene           | 0.38*     | nd          | nd          | nd        | nd        | nd        | nd        |
| Benzyl alcohol                | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| Hexachlorethane               | 0.39*     | nd          | nd          | nd        | nd        | nd        | nd        |
| N-Nitroso-di-n-propylamine    | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 3,4-Methylphenol (m,p-cresol) | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 2-Methylphenol (o-cresol)     | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| Bis (2-chloroisopropyl) ether | 0.42*     | nd          | nd          | nd        | nd        | nd        | nd        |
| Nitrobenzene                  | 0.41*     | nd          | nd          | nd        | nd        | nd        | nd        |
| Isophorone                    | 0.41*     | nd          | nd          | nd        | nd        | nd        | nd        |
| 2-Nitrophenol                 | 5.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 2,4-Dimethylphenol            | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| Bis (2-chloroethoxy) methane  | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 2,4-Dichlorophenol            | 0.36*     | nd          | nd          | nd        | nd        | nd        | nd        |
| 1,2,4-Trichlorobenzene        | 0.40*     | nd          | nd          | nd        | nd        | nd        | nd        |
| Naphthalene                   | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 4-Chloroaniline               | 0.46*     | nd          | nd          | nd        | nd        | nd        | nd        |
| Hexachlorobutadiene           | 0.43*     | nd          | nd          | nd        | nd        | nd        | nd        |
| 4-Chloro-3-methylphenol       | 5.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 2-Methylnaphthalene           | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 1-Methylnaphthalene           | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| Hexachlorocyclopentadiene     | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 2,4,6-Trichlorophenol         | 5.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 2,4,5-Trichlorophenol         | 5.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 2-Chloronaphthalene           | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 2-Nitroaniline                | 5.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 1,4-Dinitrobenzene            | 5.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| Acenaphthylene                | 0.1       | nd          | nd          | nd        | nd        | nd        | nd        |
| 1,3-Dinitrobenzene            | 0.28*     | nd          | nd          | nd        | nd        | nd        | nd        |
| Dimethylphthalate             | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 2,6-Dinitrotoluene            | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 1,2-Dinitrobenzene            | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| Acenaphthene                  | 0.1       | nd          | nd          | nd        | nd        | nd        | nd        |
| 3-Nitroaniline                | 5.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 2,4-Dinitrophenol             | 1.0*      | nd          | nd          | nd        | nd        | nd        | nd        |
| 2,4-Dinitrotoluene            | 0.35*     | nd          | nd          | nd        | nd        | nd        | nd        |
| 4-Nitrophenol                 | 5.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| Dibenzofuran                  | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 2,3,4,6-Tetrachlorophenol     | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 2,3,5,6-Tetrachlorophenol     | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| Fluorene                      | 0.1       | nd          | nd          | nd        | nd        | nd        | nd        |
| 4-Chlorophenylphenylether     | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| Diethylphthalate              | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 4-Nitroaniline                | 5.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 4,6-Dinitro-2-methylphenol    | 5.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| N-nitrosodiphenylamine        | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| Azobenzene                    | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| 4-Bromophenylphenylether      | 1.0       | nd          | nd          | nd        | nd        | nd        | nd        |
| Hexachlorobenzene             | 0.45*     | nd          | nd          | nd        | nd        | nd        | nd        |



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Analysis of Semivolatile Organic Compounds in Soil by Method 8270

Analytical Results:

|                              |           | RHS-DU-17.2 | RHS-DU-17.3 | RHS-DU-18   | RHS-DU-19   | RHS-DU-20   | RHS-DU-21   |
|------------------------------|-----------|-------------|-------------|-------------|-------------|-------------|-------------|
| Date extracted               | Reporting | 02/11/14    | 02/11/14    | 02/11/14    | 02/11/14    | 02/11/14    | 02/11/14    |
| Date analyzed                | Limits    | 02/11/14    | 02/11/14    | 02/11/14    | 02/11/14    | 02/11/14    | 02/12/14    |
| Moisture, %                  | (mg/kg)   | 11%         | 11%         | 11%         | 9%          | 18%         | 14%         |
| Pentachlorophenol            | 1.0*      | nd          | nd          | nd          | nd          | nd          | nd          |
| Phenanthrene                 | 0.1       | nd          | <b>0.12</b> | nd          | nd          | nd          | nd          |
| Anthracene                   | 0.1       | nd          | nd          | nd          | nd          | nd          | nd          |
| Carbazole                    | 1.0       | nd          | nd          | nd          | nd          | nd          | nd          |
| Di-n-butylphthalate          | 1.0       | nd          | nd          | nd          | nd          | nd          | nd          |
| Fluoranthene                 | 0.1       | <b>0.11</b> | <b>0.31</b> | <b>0.11</b> | nd          | <b>0.22</b> | <b>0.15</b> |
| Pyrene                       | 0.1       | <b>0.14</b> | <b>0.33</b> | <b>0.12</b> | <b>0.12</b> | <b>0.20</b> | <b>0.16</b> |
| Butylbenzylphthalate         | 1.0       | nd          | nd          | nd          | nd          | nd          | nd          |
| Bis(2-ethylhexyl) adipate    | 1.0       | nd          | nd          | nd          | nd          | nd          | nd          |
| Benzo(a)anthracene           | 0.1       | nd          | <b>0.18</b> | nd          | nd          | nd          | nd          |
| Chrysene                     | 0.1       | <b>0.15</b> | <b>0.46</b> | <b>0.12</b> | <b>0.11</b> | <b>0.15</b> | <b>0.12</b> |
| Bis (2-ethylhexyl) phthalate | 1.0       | nd          | nd          | nd          | nd          | nd          | nd          |
| Di-n-octyl phthalate         | 1.0       | nd          | nd          | nd          | nd          | nd          | nd          |
| Benzo(b)fluoranthene         | 0.1       | nd          | <b>0.40</b> | nd          | nd          | nd          | nd          |
| Benzo(k)fluoranthene         | 0.1       | nd          | <b>0.35</b> | nd          | nd          | nd          | nd          |
| Benzo(a)pyrene               | 0.1       | nd          | <b>0.43</b> | nd          | nd          | nd          | nd          |
| Dibenzo(a,h)anthracene       | 0.1       | nd          | nd          | nd          | nd          | nd          | nd          |
| Benzo(ghi)perylene           | 0.1       | nd          | <b>0.38</b> | nd          | nd          | nd          | nd          |
| Indeno(1,2,3-cd)pyrene       | 0.1       | nd          | <b>0.40</b> | nd          | nd          | nd          | nd          |
| <b>Surrogate recoveries</b>  |           |             |             |             |             |             |             |
| 2-Fluorophenol               |           | 106%        | 120%        | 110%        | 103%        | 110%        | 109%        |
| Phenol-d6                    |           | 97%         | 114%        | 104%        | 89%         | 107%        | 104%        |
| Nitrobenzene-d5              |           | 107%        | 111%        | 95%         | 93%         | 92%         | 95%         |
| 2-Fluorobiphenyl             |           | 137%        | 140%        | 119%        | 117%        | 117%        | 119%        |
| 2,4,6-Tribromophenol         |           | 60%         | 63%         | 58%         | 47%         | 68%         | 68%         |
| 4-Terphenyl-d14              |           | 103%        | 107%        | 92%         | 80%         | 98%         | 94%         |

Data Qualifiers and Analytical Comments

nd - not detected at listed reporting limits

\* - method detection limit

Acceptable Recovery limits:

2-Fluorophenol: 10-135 %

Phenol - d5: 10-135 %

Nitrobenzene - d5: 20-120 %

2-Fluorobiphenyl: 50-150%

2,4,6- Tribromophenol: 29-159%

p-Terphenyl-d14: 50-150%

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Analysis of Semivolatile Organic Compounds in Soil by Method 8270

Analytical Results:

|                               |           | RHS-DU-22 | RHS-DU-23 | RHS-DU-24 | RHS-DU-25 | RHS-DU-26 | RHS-DU-26.2 |
|-------------------------------|-----------|-----------|-----------|-----------|-----------|-----------|-------------|
| Date extracted                | Reporting | 02/11/14  | 02/11/14  | 02/11/14  | 02/11/14  | 02/11/14  | 02/11/14    |
| Date analyzed                 | Limits    | 02/12/14  | 02/12/14  | 02/12/14  | 02/12/14  | 02/12/14  | 02/12/14    |
| Moisture, %                   | (mg/kg)   | 12%       | 13%       | 17%       | 18%       | 30%       | 20%         |
| Pyridine                      | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| Aniline                       | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| Phenol                        | 0.41*     | nd        | nd        | nd        | nd        | nd        | nd          |
| 2-Chlorophenol                | 0.36*     | nd        | nd        | nd        | nd        | nd        | nd          |
| Bis (2-chloroethyl) ether     | 0.40*     | nd        | nd        | nd        | nd        | nd        | nd          |
| 1,3-Dichlorobenzene           | 0.34*     | nd        | nd        | nd        | nd        | nd        | nd          |
| 1,4-Dichlorobenzene           | 0.26*     | nd        | nd        | nd        | nd        | nd        | nd          |
| 1,2-Dichlorobenzene           | 0.38*     | nd        | nd        | nd        | nd        | nd        | nd          |
| Benzyl alcohol                | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| Hexachlorethane               | 0.39*     | nd        | nd        | nd        | nd        | nd        | nd          |
| N-Nitroso-di-n-propylamine    | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 3,4-Methylphenol (m,p-cresol) | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 2-Methylphenol (o-cresol)     | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| Bis (2-chloroisopropyl) ether | 0.42*     | nd        | nd        | nd        | nd        | nd        | nd          |
| Nitrobenzene                  | 0.41*     | nd        | nd        | nd        | nd        | nd        | nd          |
| Isophorone                    | 0.41*     | nd        | nd        | nd        | nd        | nd        | nd          |
| 2-Nitrophenol                 | 5.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 2,4-Dimethylphenol            | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| Bis (2-chloroethoxy) methane  | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 2,4-Dichlorophenol            | 0.36*     | nd        | nd        | nd        | nd        | nd        | nd          |
| 1,2,4-Trichlorobenzene        | 0.40*     | nd        | nd        | nd        | nd        | nd        | nd          |
| Naphthalene                   | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 4-Chloroaniline               | 0.46*     | nd        | nd        | nd        | nd        | nd        | nd          |
| Hexachlorobutadiene           | 0.43*     | nd        | nd        | nd        | nd        | nd        | nd          |
| 4-Chloro-3-methylphenol       | 5.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 2-Methylnaphthalene           | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 1-Methylnaphthalene           | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| Hexachlorocyclopentadiene     | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 2,4,6-Trichlorophenol         | 5.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 2,4,5-Trichlorophenol         | 5.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 2-Chloronaphthalene           | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 2-Nitroaniline                | 5.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 1,4-Dinitrobenzene            | 5.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| Acenaphthylene                | 0.1       | nd        | nd        | nd        | nd        | nd        | nd          |
| 1,3-Dinitrobenzene            | 0.28*     | nd        | nd        | nd        | nd        | nd        | nd          |
| Dimethylphthalate             | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 2,6-Dinitrotoluene            | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 1,2-Dinitrobenzene            | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| Acenaphthene                  | 0.1       | nd        | nd        | nd        | nd        | nd        | nd          |
| 3-Nitroaniline                | 5.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 2,4-Dinitrophenol             | 1.0*      | nd        | nd        | nd        | nd        | nd        | nd          |
| 2,4-Dinitrotoluene            | 0.35*     | nd        | nd        | nd        | nd        | nd        | nd          |
| 4-Nitrophenol                 | 5.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| Dibenzofuran                  | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 2,3,4,6-Tetrachlorophenol     | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 2,3,5,6-Tetrachlorophenol     | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| Fluorene                      | 0.1       | nd        | nd        | nd        | nd        | nd        | nd          |
| 4-Chlorophenylphenylether     | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| Diethylphthalate              | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 4-Nitroaniline                | 5.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 4,6-Dinitro-2-methylphenol    | 5.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| N-nitrosodiphenylamine        | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| Azobenzene                    | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| 4-Bromophenylphenylether      | 1.0       | nd        | nd        | nd        | nd        | nd        | nd          |
| Hexachlorobenzene             | 0.45*     | nd        | nd        | nd        | nd        | nd        | nd          |

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Analysis of Semivolatile Organic Compounds in Soil by Method 8270

Analytical Results:

|                              |           | RHS-DU-22   | RHS-DU-23   | RHS-DU-24   | RHS-DU-25   | RHS-DU-26 | RHS-DU-26.2 |
|------------------------------|-----------|-------------|-------------|-------------|-------------|-----------|-------------|
| Date extracted               | Reporting | 02/11/14    | 02/11/14    | 02/11/14    | 02/11/14    | 02/11/14  | 02/11/14    |
| Date analyzed                | Limits    | 02/12/14    | 02/12/14    | 02/12/14    | 02/12/14    | 02/12/14  | 02/12/14    |
| Moisture, %                  | (mg/kg)   | 12%         | 13%         | 17%         | 18%         | 30%       | 20%         |
| Pentachlorophenol            | 1.0*      | nd          | nd          | nd          | nd          | nd        | nd          |
| Phenanthrene                 | 0.1       | <b>0.20</b> | <b>0.16</b> | nd          | nd          | nd        | nd          |
| Anthracene                   | 0.1       | nd          | nd          | nd          | nd          | nd        | nd          |
| Carbazole                    | 1.0       | nd          | nd          | nd          | nd          | nd        | nd          |
| Di-n-butylphthalate          | 1.0       | nd          | nd          | nd          | nd          | nd        | nd          |
| Fluoranthene                 | 0.1       | <b>0.49</b> | <b>0.49</b> | <b>0.13</b> | <b>0.11</b> | nd        | nd          |
| Pyrene                       | 0.1       | <b>0.41</b> | <b>0.50</b> | <b>0.14</b> | <b>0.11</b> | nd        | nd          |
| Butylbenzylphthalate         | 1.0       | nd          | nd          | nd          | nd          | nd        | nd          |
| Bis(2-ethylhexyl) adipate    | 1.0       | nd          | nd          | nd          | nd          | nd        | nd          |
| Benzo(a)anthracene           | 0.1       | nd          | <b>0.16</b> | nd          | nd          | nd        | nd          |
| Chrysene                     | 0.1       | <b>0.40</b> | <b>0.48</b> | <b>0.21</b> | <b>0.11</b> | nd        | nd          |
| Bis (2-ethylhexyl) phthalate | 1.0       | nd          | nd          | nd          | nd          | nd        | nd          |
| Di-n-octyl phthalate         | 1.0       | nd          | nd          | nd          | nd          | nd        | nd          |
| Benzo(b)fluoranthene         | 0.1       | nd          | <b>0.31</b> | nd          | nd          | nd        | nd          |
| Benzo(k)fluoranthene         | 0.1       | nd          | <b>0.33</b> | nd          | nd          | nd        | nd          |
| Benzo(a)pyrene               | 0.1       | nd          | <b>0.15</b> | nd          | nd          | nd        | nd          |
| Dibenzo(a,h)anthracene       | 0.1       | nd          | nd          | nd          | nd          | nd        | nd          |
| Benzo(ghi)perylene           | 0.1       | nd          | <b>0.44</b> | nd          | nd          | nd        | nd          |
| Indeno(1,2,3-cd)pyrene       | 0.1       | nd          | <b>0.38</b> | nd          | nd          | nd        | nd          |
| <b>Surrogate recoveries</b>  |           |             |             |             |             |           |             |
| 2-Fluorophenol               |           | 105%        | 103%        | 108%        | 112%        | 112%      | 111%        |
| Phenol-d6                    |           | 86%         | 86%         | 91%         | 92%         | 96%       | 94%         |
| Nitrobenzene-d5              |           | 32%         | 34%         | 36%         | 33%         | 35%       | 95%         |
| 2-Fluorobiphenyl             |           | 115%        | 118%        | 116%        | 115%        | 119%      | 116%        |
| 2,4,6-Tribromophenol         |           | 42%         | 49%         | 50%         | 53%         | 56%       | 60%         |
| 4-Terphenyl-d14              |           | 91%         | 89%         | 96%         | 93%         | 95%       | 94%         |

Data Qualifiers and Analytical Comments

nd - not detected at listed reporting limits

\* - method detection limit

Acceptable Recovery limits:

2-Fluorophenol: 10-135 %

Phenol - d5: 10-135 %

Nitrobenzene - d5: 20-120 %

2-Fluorobiphenyl: 50-150%

2,4,6- Tribromophenol: 29-159%

p-Terphenyl-d14: 50-150%

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Analysis of Semivolatile Organic Compounds in Soil by Method 8270

Analytical Results:

|                               |           | RHS-DU-26.3 | RHS-DU-27 | RHS-DU-28 | RHS-DU-29 | MS       | MSD      | RPD |
|-------------------------------|-----------|-------------|-----------|-----------|-----------|----------|----------|-----|
| Date extracted                | Reporting | 02/11/14    | 02/11/14  | 02/11/14  | 02/11/14  | 02/11/14 | 02/11/14 |     |
| Date analyzed                 | Limits    | 02/13/14    | 02/13/14  | 02/13/14  | 02/13/14  | 02/13/14 | 02/13/14 |     |
| Moisture, %                   | (mg/kg)   | 23%         | 27%       | 23%       | 21%       |          |          |     |
| Pyridine                      | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| Aniline                       | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| Phenol                        | 0.41*     | nd          | nd        | nd        | nd        | 104%     | 82%      | 24% |
| 2-Chlorophenol                | 0.36*     | nd          | nd        | nd        | nd        | 113%     | 92%      | 20% |
| Bis (2-chloroethyl) ether     | 0.40*     | nd          | nd        | nd        | nd        |          |          |     |
| 1,3-Dichlorobenzene           | 0.34*     | nd          | nd        | nd        | nd        |          |          |     |
| 1,4-Dichlorobenzene           | 0.26*     | nd          | nd        | nd        | nd        | 110%     | 91%      | 19% |
| 1,2-Dichlorobenzene           | 0.38*     | nd          | nd        | nd        | nd        |          |          |     |
| Benzyl alcohol                | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| Hexachloroethane              | 0.39*     | nd          | nd        | nd        | nd        |          |          |     |
| N-Nitroso-di-n-propylamine    | 1.0       | nd          | nd        | nd        | nd        | 78%      | 66%      | 17% |
| 3,4-Methylphenol (m,p-cresol) | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| 2-Methylphenol (o-cresol)     | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| Bis (2-chloroisopropyl) ether | 0.42*     | nd          | nd        | nd        | nd        |          |          |     |
| Nitrobenzene                  | 0.41*     | nd          | nd        | nd        | nd        |          |          |     |
| Isophorone                    | 0.41*     | nd          | nd        | nd        | nd        |          |          |     |
| 2-Nitrophenol                 | 5.0       | nd          | nd        | nd        | nd        |          |          |     |
| 2,4-Dimethylphenol            | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| Bis (2-chloroethoxy) methane  | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| 2,4-Dichlorophenol            | 0.36*     | nd          | nd        | nd        | nd        |          |          |     |
| 1,2,4-Trichlorobenzene        | 0.40*     | nd          | nd        | nd        | nd        | 107%     | 88%      | 19% |
| Naphthalene                   | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| 4-Chloroaniline               | 0.46*     | nd          | nd        | nd        | nd        |          |          |     |
| Hexachlorobutadiene           | 0.43*     | nd          | nd        | nd        | nd        |          |          |     |
| 4-Chloro-3-methylphenol       | 5.0       | nd          | nd        | nd        | nd        | 89%      | 76%      | 16% |
| 2-Methylnaphthalene           | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| 1-Methylnaphthalene           | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| Hexachlorocyclopentadiene     | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| 2,4,6-Trichlorophenol         | 5.0       | nd          | nd        | nd        | nd        |          |          |     |
| 2,4,5-Trichlorophenol         | 5.0       | nd          | nd        | nd        | nd        |          |          |     |
| 2-Chloronaphthalene           | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| 2-Nitroaniline                | 5.0       | nd          | nd        | nd        | nd        |          |          |     |
| 1,4-Dinitrobenzene            | 5.0       | nd          | nd        | nd        | nd        |          |          |     |
| Acenaphthylene                | 0.1       | nd          | nd        | nd        | nd        |          |          |     |
| 1,3-Dinitrobenzene            | 0.28*     | nd          | nd        | nd        | nd        |          |          |     |
| Dimethylphthalate             | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| 2,6-Dinitrotoluene            | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| 1,2-Dinitrobenzene            | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| Acenaphthene                  | 0.1       | nd          | nd        | nd        | nd        | 127%     | 102%     | 22% |
| 3-Nitroaniline                | 5.0       | nd          | nd        | nd        | nd        |          |          |     |
| 2,4-Dinitrophenol             | 1.0*      | nd          | nd        | nd        | nd        |          |          |     |
| 2,4-Dinitrotoluene            | 0.35*     | nd          | nd        | nd        | nd        | 53%      | 51%      | 4%  |
| 4-Nitrophenol                 | 5.0       | nd          | nd        | nd        | nd        |          |          |     |
| Dibenzofuran                  | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| 2,3,4,6-Tetrachlorophenol     | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| 2,3,5,6-Tetrachlorophenol     | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| Fluorene                      | 0.1       | nd          | nd        | nd        | nd        |          |          |     |
| 4-Chlorophenylphenylether     | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| Diethylphthalate              | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| 4-Nitroaniline                | 5.0       | nd          | nd        | nd        | nd        |          |          |     |
| 4,6-Dinitro-2-methylphenol    | 5.0       | nd          | nd        | nd        | nd        |          |          |     |
| N-nitrosodiphenylamine        | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| Azobenzene                    | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| 4-Bromophenylphenylether      | 1.0       | nd          | nd        | nd        | nd        |          |          |     |
| Hexachlorobenzene             | 0.45*     | nd          | nd        | nd        | nd        |          |          |     |

ESN NORTHWEST CHEMISTRY LABORATORY

Bureau Veritas  
 BV-RADFORD HIGH SCHOOL PROJECT  
 Client Project #17012-012148.48  
 ESN Project #D1402050037  
 Hawaii

ESN Northwest  
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 Olympia, WA 98501  
 (360) 459-4670 (360) 459-3432 Fax  
 lab@esnw.com

Analysis of Semivolatile Organic Compounds in Soil by Method 8270

Analytical Results:

|                |           | RHS-DU-26.3 | RHS-DU-27 | RHS-DU-28 | RHS-DU-29 | MS       | MSD      | RPD |
|----------------|-----------|-------------|-----------|-----------|-----------|----------|----------|-----|
| Date extracted | Reporting | 02/11/14    | 02/11/14  | 02/11/14  | 02/11/14  | 02/11/14 | 02/11/14 |     |
| Date analyzed  | Limits    | 02/13/14    | 02/13/14  | 02/13/14  | 02/13/14  | 02/13/14 | 02/13/14 |     |
| Moisture, %    | (mg/kg)   | 23%         | 27%       | 23%       | 21%       |          |          |     |

|                              |      |    |    |             |             |     |     |     |
|------------------------------|------|----|----|-------------|-------------|-----|-----|-----|
| Pentachlorophenol            | 1.0* | nd | nd | nd          | nd          |     |     |     |
| Phenanthrene                 | 0.1  | nd | nd | nd          | nd          |     |     |     |
| Anthracene                   | 0.1  | nd | nd | nd          | nd          |     |     |     |
| Carbazole                    | 1.0  | nd | nd | nd          | nd          |     |     |     |
| Di-n-butylphthalate          | 1.0  | nd | nd | nd          | nd          |     |     |     |
| Fluoranthene                 | 0.1  | nd | nd | <b>0.11</b> | <b>0.10</b> |     |     |     |
| Pyrene                       | 0.1  | nd | nd | <b>0.10</b> | <b>0.10</b> | 94% | 82% | 14% |
| Butylbenzylphthalate         | 1.0  | nd | nd | nd          | nd          |     |     |     |
| Bis(2-ethylhexyl) adipate    | 1.0  | nd | nd | nd          | nd          |     |     |     |
| Benzo(a)anthracene           | 0.1  | nd | nd | nd          | nd          |     |     |     |
| Chrysene                     | 0.1  | nd | nd | <b>0.10</b> | <b>0.14</b> |     |     |     |
| Bis (2-ethylhexyl) phthalate | 1.0  | nd | nd | nd          | nd          |     |     |     |
| Di-n-octyl phthalate         | 1.0  | nd | nd | nd          | nd          |     |     |     |
| Benzo(b)fluoranthene         | 0.1  | nd | nd | nd          | nd          |     |     |     |
| Benzo(k)fluoranthene         | 0.1  | nd | nd | nd          | nd          |     |     |     |
| Benzo(a)pyrene               | 0.1  | nd | nd | nd          | nd          |     |     |     |
| Dibenzo(a,h)anthracene       | 0.1  | nd | nd | nd          | nd          |     |     |     |
| Benzo(ghi)perylene           | 0.1  | nd | nd | nd          | nd          |     |     |     |
| Indeno(1,2,3-cd)pyrene       | 0.1  | nd | nd | nd          | nd          |     |     |     |

Surrogate recoveries

|                      |      |      |      |      |      |      |
|----------------------|------|------|------|------|------|------|
| 2-Fluorophenol       | 109% | 108% | 110% | 106% | 104% | 103% |
| Phenol-d6            | 87%  | 88%  | 86%  | 89%  | 97%  | 98%  |
| Nitrobenzene-d5      | 22%  | 20%  | 20%  | 22%  | 96%  | 95%  |
| 2-Fluorobiphenyl     | 114% | 110% | 117% | 112% | 109% | 107% |
| 2,4,6-Tribromophenol | 39%  | 42%  | 39%  | 43%  | 42%  | 46%  |
| 4-Terphenyl-d14      | 79%  | 82%  | 80%  | 79%  | 72%  | 74%  |

Data Qualifiers and Analytical Comments

nd - not detected at listed reporting limits

\* - method detection limit

Acceptable Recovery limits:

2-Fluorophenol: 10-135 %

Phenol - d5: 10-135 %

Nitrobenzene - d5: 20-120 %

2-Fluorobiphenyl: 50-150%

2,4,6- Tribromophenol: 29-159%

p-Terphenyl-d14: 50-150%

Acceptable RPD limit: 35%

**ESN NORTHWEST CHEMISTRY LABORATORY**

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 RADFORD HIGH SCHOOL PROJECT  
 Client Project #17012-012148.48  
 ESN Project #D1402050037  
 Hawaii

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 lab@esnnw.com

**Total Metals in Soil by EPA-6020 Series**

| Sample Number       | Date Analyzed | Lead (Pb) (mg/kg) | Cadmium (Cd) (mg/kg) | Chromium (Cr) (mg/kg) | Arsenic (As) (mg/kg) | Silver (Ag) (mg/kg) | Barium (Ba) (mg/kg) | Selenium (Se) (mg/kg) | Mercury (Hg) (mg/kg) |
|---------------------|---------------|-------------------|----------------------|-----------------------|----------------------|---------------------|---------------------|-----------------------|----------------------|
| Method Blank        | 2/10/2014     | nd                | nd                   | nd                    | nd                   | nd                  | nd                  | nd                    | nd                   |
| RHS-DU-1            | 2/11/2014     | 4800              | 14**                 | 300*                  | 35**                 | nd                  | 1200                | nd                    | 5.2                  |
| RHS-DU-1 Duplicate  | 2/11/2014     | 5200              | 12**                 | 350*                  | 33**                 | nd                  | 1100                | nd                    | 4.9                  |
| RHS-DU-1.2          | 2/11/2014     | 5800              | 14**                 | 230*                  | 38**                 | nd                  | 1100                | nd                    | 11                   |
| RHS-DU-1.3          | 2/11/2014     | 15000             | 26**                 | 310*                  | 28**                 | nd                  | 1100                | nd                    | 18                   |
| RHS-DU-2            | 2/11/2014     | 1600              | 5.9**                | 150*                  | 11**                 | nd                  | 550                 | nd                    | 15                   |
| RHS-DU-3            | 2/12/2014     | 85                | nd                   | 160                   | 6.0                  | nd                  | 410                 | nd                    | 5.7                  |
| RHS-DU-4            | 2/12/2014     | 140               | nd                   | 170                   | 7.5                  | nd                  | 460                 | nd                    | 3.5                  |
| RHS-DU-5            | 2/10/2014     | 61                | nd                   | 180                   | 6.1                  | nd                  | 280                 | nd                    | 2.9                  |
| RHS-DU-6            | 2/10/2014     | 66                | nd                   | 240                   | 7.6                  | nd                  | 200                 | nd                    | 3.3                  |
| RHS-DU-7            | 2/10/2014     | 190               | 1.0                  | 230                   | 7.3                  | nd                  | 280                 | nd                    | 6.1                  |
| RHS-DU-8            | 2/10/2014     | 260               | 1.3                  | 170                   | 6.8                  | nd                  | 320                 | nd                    | 21                   |
| RHS-DU-9            | 2/12/2014     | 660               | 1.6**                | 110*                  | 8.7**                | nd                  | 320                 | nd                    | 18                   |
| RHS-DU-10           | 2/10/2014     | 490               | 1.7                  | 180                   | 12                   | nd                  | 370                 | nd                    | 23                   |
| RHS-DU-10 Duplicate | 2/10/2014     | 420               | 2.2                  | 180                   | 11                   | nd                  | 320                 | nd                    | 22                   |
| RHS-DU-11           | 2/10/2014     | 1000              | 3.7                  | 200                   | 19                   | nd                  | 522                 | nd                    | 68                   |
| RHS-DU-12           | 2/10/2014     | 1300              | 7.1                  | 210                   | 23                   | nd                  | 670                 | nd                    | 57                   |
| RHS-DU-13           | 2/10/2014     | 2800              | 6.8                  | 240                   | 19                   | nd                  | 740                 | nd                    | 23                   |
| RHS-DU-14           | 2/10/2014     | 240               | nd                   | 220                   | 12                   | nd                  | 260                 | nd                    | 8.2                  |
| RHS-DU-15           | 2/10/2014     | 360               | 1.4                  | 190                   | 15                   | nd                  | 270                 | nd                    | 14                   |
| RHS-DU-16           | 2/10/2014     | 95                | nd                   | 150                   | 5.4                  | nd                  | 220                 | nd                    | 5.0                  |
| RHS-DU-17           | 2/12/2014     | 1100              | 4.5**                | 210*                  | 23**                 | nd                  | 590                 | nd                    | 73                   |
| RHS-DU-17.2         | 2/10/2014     | 1600              | 4.6                  | 220                   | 24                   | nd                  | 720                 | nd                    | 66                   |
| RHS-DU-17.3         | 2/12/2014     | 1600              | 6.4**                | 150*                  | 26**                 | nd                  | 670                 | nd                    | 83                   |
| RHS-DU-18           | 2/12/2014     | 6200              | 11**                 | 170*                  | 26**                 | nd                  | 710                 | nd                    | 45                   |
| RHS-DU-19           | 2/12/2014     | 2300              | 7.8**                | 160*                  | 26**                 | nd                  | 610                 | nd                    | 45                   |
| RHS-DU-20           | 2/10/2014     | 120               | nd                   | 190                   | 6.4                  | nd                  | 290                 | nd                    | 18                   |
| RHS-DU-21           | 2/10/2014     | 2300              | 12                   | 240                   | 18                   | nd                  | 810                 | nd                    | 20                   |
| RHS-DU-22           | 2/12/2014     | 110               | nd                   | 130                   | 6.9                  | nd                  | 320                 | nd                    | 18                   |
| RHS-DU-23           | 2/12/2014     | 500               | 1.9                  | 150                   | 18                   | nd                  | 340                 | nd                    | 200                  |
| RHS-DU-24           | 2/10/2014     | 110               | nd                   | 140                   | 5.1                  | nd                  | 270                 | nd                    | 4.8                  |
| RHS-DU-25           | 2/10/2014     | 120               | nd                   | 190                   | 5.8                  | nd                  | 280                 | nd                    | 4.4                  |
| RHS-DU-26           | 2/13/2001     | 89                | nd                   | 250*                  | 5.3                  | nd                  | 100                 | nd                    | 3.5                  |
| RHS-DU-26.2         | 2/13/2014     | 50                | nd                   | 220*                  | nd                   | nd                  | 69                  | nd                    | 2.7                  |
| RHS-DU-26.3         | 2/13/2014     | 78                | nd                   | 190*                  | 6.2                  | nd                  | 85                  | nd                    | 2.9                  |
| RHS-DU-27           | 2/13/2014     | 59                | nd                   | 230*                  | 6.3                  | nd                  | 85                  | nd                    | 2.9                  |
| RHS-DU-27 Duplicate | 2/13/2014     | 39                | nd                   | 220*                  | 5.5                  | nd                  | 78                  | nd                    | 2.8                  |
| RHS-DU-28           | 2/13/2014     | 44                | nd                   | 250*                  | 4.6                  | nd                  | 91                  | nd                    | 3.1                  |
| RHS-DU-29           | 2/13/2014     | 68                | nd                   | 270*                  | 5.7                  | nd                  | 99                  | nd                    | 5.1                  |
| Reporting Limits    |               | 5.0               | 1.0                  | 5.0                   | 5.0                  | 20                  | 50                  | 20                    | 0.5                  |

nd - not detected at listed reporting limits

\* The Continuing Calibration standard failed low. The samples were reanalyzed with similar results.

\*\* The Continuing Calibration standard failed high. The samples were reanalyzed with similar results.

**QA/QC Data - Total Metals EPA-6020**

| Sample Number: RHS-DU-27 |                      |                        |                    |                        |                        |                    |      |
|--------------------------|----------------------|------------------------|--------------------|------------------------|------------------------|--------------------|------|
|                          | Matrix Spike         |                        |                    | Matrix Spike Duplicate |                        |                    | RPD  |
|                          | Spiked Conc. (mg/kg) | Measured Conc. (mg/kg) | Spike Recovery (%) | Spiked Conc. (mg/kg)   | Measured Conc. (mg/kg) | Spike Recovery (%) |      |
| Lead                     | 70.2                 | 67.1                   | 95.6               | 71.2                   | 77.6                   | 109                | 13.1 |
| Cadmium                  | 70.2                 | 54.5                   | 77.6M              | 71.2                   | 58.3                   | 81.9               | 5.32 |
| Chromium                 | 70.2                 | 109                    | 155M               | 71.2                   | 101                    | 142M               | 9.0  |
| Arsenic                  | 70.2                 | 58.3                   | 83.0               | 71.2                   | 62.6                   | 87.9               | 5.70 |
| Silver                   | 70.2                 | 45.4                   | 64.7M              | 71.2                   | 48.9                   | 68.7M              | 6.01 |
| Barium                   | 70.2                 | 84.9                   | 121M               | 71.2                   | 108                    | 152M               | 22.6 |
| Selenium                 | 70.2                 | 52.2                   | 74.4M              | 71.2                   | 55.0                   | 77.2M              | 3.81 |
| Mercury                  | 7.02                 | 6.06                   | 86.3               | 7.12                   | 6.97                   | 97.9               | 12.6 |

| Laboratory Control Sample |                      |                        |                    |
|---------------------------|----------------------|------------------------|--------------------|
|                           | Spiked Conc. (mg/kg) | Measured Conc. (mg/kg) | Spike Recovery (%) |
| Lead                      | 100                  | 101                    | 101                |
| Cadmium                   | 100                  | 97.5                   | 97.5               |
| Chromium                  | 100                  | 102                    | 102                |
| Arsenic                   | 100                  | 103                    | 103                |
| Silver                    | 100                  | 85.8                   | 85.8               |
| Barium                    | 100                  | 99.1                   | 99.1               |
| Selenium                  | 100                  | 101                    | 101                |
| Mercury                   | 10                   | 10.6                   | 106                |

ACCEPTABLE RECOVERY LIMITS FOR MATRIX SPIKES: 80%-120%

ACCEPTABLE RPD IS 35%

M - Matrix Spike recovery failed due to matrix interference.

**Report Prepared for:**

Karen Carvallo  
ESN Pacific  
2020-B Kahai Street  
Honolulu HI 96819

**REPORT OF  
LABORATORY  
ANALYSIS FOR  
PCDD/PCDF**

**Report Information:**

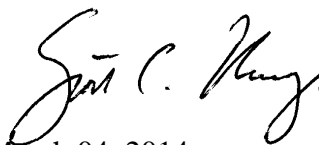
**Pace Project #: 10257346**  
**Sample Receipt Date: 02/11/2014**  
**Client Project #: D1402050037**  
**Client Sub PO #: N/A**  
**State Cert #: SLD**

**Invoicing & Reporting Options:**

The report provided has been invoiced as a Level 2 PCDD/PCDF Report. If an upgrade of this report package is requested, an additional charge may be applied.

Please review the attached invoice for accuracy and forward any questions to Scott Unze, your Pace Project Manager.

**This report has been reviewed by:**



March 04, 2014

Scott Unze, Project Manager  
(612) 607-6383  
(612) 607-6444 (fax)  
scott.unze@pacelabs.com

**Report Prepared Date:**

March 3, 2014



**Report of Laboratory Analysis**

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The results relate only to the samples included in this report.





## **DISCUSSION**

This report presents the results from the analyses performed on eighteen samples submitted by a representative of ESN Pacific. The samples were analyzed for the presence or absence of polychlorodibenzo-p-dioxins (PCDDs) and polychlorodibenzofurans (PCDFs) using a modified version of USEPA Method 8290. The reporting limits were based on signal-to-noise measurements.

The results provided in this report include lower-bound, mid-bound, and upper-bound toxic equivalence (TEQ) results. Lower bound TEQ results include only contributions from positive values in the sample. Mid-bound TEQ results include one-half of the reporting limit in cases where an analyte was not detected, along with contributions from positives in the sample. Upper-bound TEQ results include the full reporting limit in cases where an analyte was not detected, along with contributions from positives in the sample.

Second column confirmation analyses of 2,3,7,8-TCDF values obtained from the primary (DB5-MS) column are performed only when specifically requested for a project and only when the values are above the concentration of the lowest calibration standard. Typical resolution for this isomer using the DB5-MS column ranges from 25-30%.

The recoveries of the isotopically-labeled PCDD/PCDF internal standards in the sample extracts ranged from 40-107%. All of the labeled standard recoveries obtained for this project were within the 40-135% target range specified in Method 8290. Also, since the quantification of the native 2,3,7,8-substituted congeners was based on isotope dilution, the data were automatically corrected for variation in recovery and accurate values were obtained.

In some cases, interfering substances impacted the determinations of PCDD or PCDF congeners; the affected values were flagged "I" where incorrect isotope ratios were obtained or "P" where polychlorinated diphenyl ethers were present. Concentrations below the calibration range were flagged "J" and should be regarded as estimates. Concentrations above the calibration range were flagged "E" and should also be regarded as estimates.

A laboratory method blank was prepared and analyzed with each sample batch as part of our routine quality control procedures. The results show the blanks to contain trace levels of selected congeners. These levels were below the calibration range of the method. The levels reported for the affected congeners in the field samples were higher than the corresponding blank levels by one or more orders of magnitude. These results indicate that the sample processing steps did not contribute significantly to the levels reported for the field samples.

Laboratory and matrix spike samples were also prepared with the sample batches using clean sand or sample matrix that had been fortified with native standard materials. The results show that the spiked native compounds in the laboratory spike samples were recovered at 89-128% with relative percent differences (RPDs) of 0.0-7.5%. These results indicate high degrees of accuracy and precision for these determinations. Several recovery values obtained for the native congeners in the matrix spike samples were outside of the 70-130% target range due to the levels of these congeners in the sample material. The RPD values obtained for the matrix spike analyses ranged from 0.2-5.1%. Matrix spikes were not prepared with the 02/24/2014 sample batch.

The responses obtained for selected native and labeled congeners in calibration standard analyses U140226B\_20 and P140301A\_15 were outside the target ranges. As specified in the method, the averages of the daily response factors for these compounds were used in the calculations for the samples

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**DISCUSSION**

from these runshifts. The affected values were flagged "Y" on the results tables.

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## Minnesota Laboratory Certifications

| Authority       | Certificate # | Authority       | Certificate # |
|-----------------|---------------|-----------------|---------------|
| A2LA            | 2926.01       | Minnesota       | 027-053-137   |
| Alabama         | 40770         | Mississippi     | MN00064       |
| Alaska          | MN00064       | Montana         | 92            |
| Arizona         | AZ0014        | Nebraska        |               |
| Arkansas        | 88-0680       | Nevada          | MN_00064_200  |
| California      | 01155CA       | New Jersey (NE) | MN002         |
| Colorado        | MN00064       | New York (NEL)  | 11647         |
| Connecticut     | PH-0256       | North Carolina  | 27700         |
| EPA Region 8    | 8TMS-Q        | North Dakota    | R-036         |
| Florida (NELAP) | E87605        | Ohio            | 4150          |
| Georgia (DNR)   | 959           | Oklahoma        | D9922         |
| Guam            | 959           | Oregon (ELAP)   | MN200001-005  |
| Hawaii          | SLD           | Oregon (OREL)   | MN300001-001  |
| Idaho           | MN00064       | Pennsylvania    | 68-00563      |
| Illinois        | 200012        | Puerto Rico     | MN00064       |
| Indiana         | C-MN-01       | Saipan          | MP0003        |
| Indiana         | C-MN-01       | South Carolina  | 74003001      |
| Iowa            | 368           | Texas           | T104704192-08 |
| Kansas          | E-10167       | Utah (NELAP)    | MN00064       |
| Kentucky        | 90062         | Virginia        | 00251         |
| Louisiana       | 03086         | Washington      | C755          |
| Maine           | 2007029       | West Virginia   | 9952C         |
| Maryland        | 322           | Wisconsin       | 999407970     |
| Michigan        | 9909          | Wyoming         | 8TMS-Q        |

## REPORT OF LABORATORY ANALYSIS

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## Appendix A

### Sample Management

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10257346

# ESN PACIFIC'S CHAIN-OF-CUSTODY RECORD

CLIENT: ESN Pacific  
 ADDRESS: 2020-B Kahai St Honolulu, HI 96819  
 PHONE: 8088470067  
 EMAIL: esn@esnpacific.com  
 TAT: 10-day  
 DATE: 2-7-14  
 ESN PROJECT #: D1402050037  
 LOCATION/PROJECT NAME: BV - Radford High School  
 COLLECTOR: DATE COLLECTED: 2/4-2/5  
 Project Manager: K. Carvallo  
 Dioxins/Furans 8290

PAGE 1 OF 2

| Sample ID#   | Date | Time | Sample Type | Container Type | Comments | # of Containers |
|--------------|------|------|-------------|----------------|----------|-----------------|
| 1 RHS-DU-1   | 2/4  | 1525 | soil        | X              |          | 1               |
| 2 RHS-DU-1.2 | 2/5  | 933  |             | X              |          | 1               |
| 3 RHS-DU-1.3 | 2/5  | 1030 |             | X              |          | 1               |
| 4 RHS-DU-2   | 2/4  | 1500 |             | X              |          | 1               |
| 5 RHS-DU-3   | 2/4  | 1403 |             | X              |          | 1               |
| 6 RHS-DU-4   | 2/4  | 1145 |             | X              |          | 1               |
| 7 RHS-DU-5   | 2/4  | 1220 |             | X              |          | 1               |
| 8 RHS-DU-6   | 2/4  | 1402 |             | X              |          | 1               |
| 9 RHS-DU-7   | 2/4  | 1131 |             | X              |          | 1               |
| 10 RHS-DU-8  | 2/4  | 1424 |             | X              |          | 1               |
| 11 RHS-DU-9  | 2/4  | 1455 |             | X              |          | 1               |
| 12 RHS-DU-10 | 2/5  | 915  |             | X              |          | 1               |
| 13 RHS-DU-11 | 2/5  | 900  |             | X              |          | 1               |
| 14 RHS-DU-12 | 2/5  | 945  |             | X              |          | 1               |
| 15 RHS-DU-13 | 2/5  | 945  |             | X              |          | 1               |
| 16 RHS-DU-14 | 2/4  | 1252 |             | X              |          | 1               |
| 17 RHS-DU-15 | 2/4  | 1345 |             | X              |          | 1               |
| 18 RHS-DU-16 | 2/4  | 1411 |             | X              |          | 1               |

Project Manager: K. Carvallo  
 Dioxins/Furans 8290

RECEIVED BY (Signature): *K. Carvallo* DATE/TIME: 2/7/14 1500  
 RECEIVED BY (Signature): *Carroll Pace* DATE/TIME: 2/4/14 907  
 RECEIVED BY (Signature): \_\_\_\_\_ DATE/TIME: \_\_\_\_\_

SAMPLE RECEIPT:  
 TOTAL # OF CONTAINERS: 18 (of 35)  
 COC SEALS Y / N / NA  
 SEALS INTACT Y / N / NA  
 RECEIVED TEMP: \_\_\_\_\_

LABORATORY NOTES:  
 Samples MI prepped at ESN Pacific

**Sample Condition Upon Receipt**

Client Name: ESN Pacifics

Project #: **WO#: 10257346**



Courier:  Fed Ex  UPS  USPS  Client  
 Commercial  Pace  Other: \_\_\_\_\_

Tracking Number: 18 VZ3 U3F B3 9904 0403

Custody Seal on Cooler/Box Present?  Yes  No      Seals Intact?  Yes  No      Optional: Proj. Due Date: \_\_\_\_\_ Proj. Name: \_\_\_\_\_

Packing Material:  Bubble Wrap  Bubble Bags  None  Other: \_\_\_\_\_      Temp Blank?  Yes  No

Thermom. Used:  80512447  72337080  B88A912167504  B88A9132521491      Type of Ice:  Wet  Blue  None  Samples on ice, cooling process has begun

Cooler Temp Read (°C): 24      Cooler Temp Corrected (°C): 24      Biological Tissue Frozen?  Yes  No  N/A  
 Temp should be above freezing to 6°C      Correction Factor: TRUE      Date and Initials of Person Examining Contents: CMB 2/11/14

**Comments:**

|   |  |  |
|---|--|--|
| Chain of Custody Present?   | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A | 1.   |
| Chain of Custody Filled Out?  | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A | 2.   |
| Chain of Custody Relinquished?  | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A | 3.   |
| Sampler Name and/or Signature on COC?   | <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <input type="checkbox"/> N/A | 4.   |
| Samples Arrived within Hold Time?   | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A | 5.   |
| Short Hold Time Analysis (<72 hr)?  | <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <input type="checkbox"/> N/A | 6.   |
| Rush Turn Around Time Requested?  | <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <input type="checkbox"/> N/A | 7.   |
| Sufficient Volume?  | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A | 8.   |
| Correct Containers Used?  | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A | 9.   |
| -Pace Containers Used?  | <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <input type="checkbox"/> N/A |  |
| Containers Intact?  | <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> N/A | 10.  |
| Filtered Volume Received for Dissolved Tests?   | <input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A | 11.  |
| Sample Labels Match COC?  | <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No <input type="checkbox"/> N/A | 12. <u>No date/time on samples</u>   |
| -Includes Date/Time/ID/Analysis Matrix: <u>SL</u>   |  |  |
| All containers needing acid/base preservation have been checked? Noncompliances are noted in 13.  | <input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A | 13. <input type="checkbox"/> HNO <sub>3</sub> <input type="checkbox"/> H <sub>2</sub> SO <sub>4</sub> <input type="checkbox"/> NaOH <input type="checkbox"/> HCl |
| All containers needing preservation are found to be in compliance with EPA recommendation? (HNO <sub>3</sub> , H <sub>2</sub> SO <sub>4</sub> , HCl<2; NaOH>12) | <input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A | Sample #   |
| Exceptions: VOA, Coliform, TOC, Oil and Grease, WI-DRO (water) DOC  | <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No                              | Initial when completed: _____ Lot # of added preservative: _____   |
| Headspace in VOA Vials (>6mm)?  | <input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A | 14.  |
| Trip Blank Present?   | <input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A | 15.  |
| Trip Blank Custody Seals Present?   | <input type="checkbox"/> Yes <input type="checkbox"/> No <input checked="" type="checkbox"/> N/A |  |
| Pace Trip Blank Lot # (if purchased): _____   |  |  |

**CLIENT NOTIFICATION/RESOLUTION**

Field Data Required?  Yes  No

Person Contacted: \_\_\_\_\_ Date/Time: \_\_\_\_\_

Comments/Resolution: \_\_\_\_\_

**Project Manager Review:**

Date: 02/11/14

Note: Whenever there is a discrepancy affecting North Carolina compliance samples, a copy of this form will be sent to the North Carolina DEHR Certification Office (i.e. out of hold, incorrect preservative, out of temp, incorrect containers)

## Reporting Flags

- A = Reporting Limit based on signal to noise
- B = Less than 10x higher than method blank level
- C = Result obtained from confirmation analysis
- D = Result obtained from analysis of diluted sample
- E = Exceeds calibration range
- I = Interference present
- J = Estimated value
- Nn = Value obtained from additional analysis
- P = PCDE Interference
- R = Recovery outside target range
- S = Peak saturated
- U = Analyte not detected
- V = Result verified by confirmation analysis
- X = %D Exceeds limits
- Y = Calculated using average of daily RFs
- \* = See Discussion

### REPORT OF LABORATORY ANALYSIS



## Appendix B

### Sample Analysis Summary

CarrollCox.com



### Method 8290 Sample Analysis Results

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-1                  |           |                  |  |
| Lab Sample ID          | 10257346001-S             |           |                  |  |
| Filename               | U140226B_14               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.1 g                    | Matrix    | Soil             |  |
| % Moisture             | 5.2                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.57 g                    | Collected | 02/04/2014 15:25 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140226B_06 & U140226B_20 | Extracted | 02/24/2014 21:00 |  |
| Method Blank ID        | BLANK-39446               | Analyzed  | 02/27/2014 03:40 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg |   | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|---|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 460        | ----       | 0.91     | E | 2,3,7,8-TCDF-13C         | 2.00       | 88 Y             |
| Total TCDF          | 13000      | ----       | 0.91     | E | 2,3,7,8-TCDD-13C         | 2.00       | 89               |
|                     |            |            |          |   | 1,2,3,7,8-PeCDF-13C      | 2.00       | 95 Y             |
| 2,3,7,8-TCDD        | 29         | ----       | 0.16     | Y | 2,3,4,7,8-PeCDF-13C      | 2.00       | 94 Y             |
| Total TCDD          | 1500       | ----       | 0.16     | Y | 1,2,3,7,8-PeCDD-13C      | 2.00       | 96 Y             |
|                     |            |            |          |   | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 97               |
| 1,2,3,7,8-PeCDF     | 360        | ----       | 0.86     |   | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 97               |
| 2,3,4,7,8-PeCDF     | ----       | 800        | 0.60     | P | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 99               |
| Total PeCDF         | 9100       | ----       | 0.73     |   | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 91               |
|                     |            |            |          |   | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 81               |
| 1,2,3,7,8-PeCDD     | 110        | ----       | 0.46     |   | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 76               |
| Total PeCDD         | 1800       | ----       | 0.46     |   | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 73               |
|                     |            |            |          |   | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 67               |
| 1,2,3,4,7,8-HxCDF   | 610        | ----       | 0.56     |   | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 63               |
| 1,2,3,6,7,8-HxCDF   | ----       | 540        | 0.53     | P | OCDD-13C                 | 4.00       | 70               |
| 2,3,4,6,7,8-HxCDF   | 750        | ----       | 0.59     |   |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 100        | ----       | 0.56     |   | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 6100       | ----       | 0.56     | E | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |   |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 91         | ----       | 0.84     |   | 2,3,7,8-TCDD-37Cl4       | 0.20       | 96               |
| 1,2,3,6,7,8-HxCDD   | 170        | ----       | 0.45     |   |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 130        | ----       | 0.40     |   |                          |            |                  |
| Total HxCDD         | 2200       | ----       | 0.56     |   |                          |            |                  |
|                     |            |            |          |   |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 2500       | ----       | 0.46     | E | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 130        | ----       | 0.38     |   | Equivalence: 710 ng/Kg   |            |                  |
| Total HpCDF         | 3100       | ----       | 0.42     | E | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |   |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 840        | ----       | 0.53     |   |                          |            |                  |
| Total HpCDD         | 1700       | ----       | 0.53     |   |                          |            |                  |
|                     |            |            |          |   |                          |            |                  |
| OCDF                | 910        | ----       | 0.35     | Y |                          |            |                  |
| OCDD                | 2400       | ----       | 0.21     |   |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
P = PCDE Interference  
E = Exceeds calibration range  
Y = Calculated using average of daily RFs

## REPORT OF LABORATORY ANALYSIS

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-1                  |           |                  |  |
| Lab Sample ID          | 10257346001-S             |           |                  |  |
| Filename               | U140226B_14               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.1 g                    | Matrix    | Soil             |  |
| % Moisture             | 5.2                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.57 g                    | Collected | 02/04/2014 15:25 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140226B_06 & U140226B_20 | Extracted | 02/24/2014 21:00 |  |
| Method Blank ID        | BLANK-39446               | Analyzed  | 02/27/2014 03:40 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB       | MB       | UB       |
|---------------------|---------------|-------------|---------|----------|----------|----------|
| 2,3,7,8-TCDF        | 460           | 0.91        | 0.10000 | 46.1804  | 46.1804  | 46.1804  |
| Total TCDF          | 13000         | 0.91        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 2,3,7,8-TCDD        | 29            | 0.16        | 1.00000 | 29.2299  | 29.2299  | 29.2299  |
| Total TCDD          | 1500          | 0.16        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,7,8-PeCDF     | 360           | 0.86        | 0.03000 | 10.8854  | 10.8854  | 10.8854  |
| 2,3,4,7,8-PeCDF     | ND            | 0.60        | 0.30000 | 239.3679 | 239.3679 | 239.3679 |
| Total PeCDF         | 9100          | 0.73        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,7,8-PeCDD     | 110           | 0.46        | 1.00000 | 113.0036 | 113.0036 | 113.0036 |
| Total PeCDD         | 1800          | 0.46        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,7,8-HxCDF   | 610           | 0.56        | 0.10000 | 61.0951  | 61.0951  | 61.0951  |
| 1,2,3,6,7,8-HxCDF   | ND            | 0.53        | 0.10000 | 54.4034  | 54.4034  | 54.4034  |
| 2,3,4,6,7,8-HxCDF   | 750           | 0.59        | 0.10000 | 74.5342  | 74.5342  | 74.5342  |
| 1,2,3,7,8,9-HxCDF   | 100           | 0.56        | 0.10000 | 10.1314  | 10.1314  | 10.1314  |
| Total HxCDF         | 6100          | 0.56        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,7,8-HxCDD   | 91            | 0.84        | 0.10000 | 9.1338   | 9.1338   | 9.1338   |
| 1,2,3,6,7,8-HxCDD   | 170           | 0.45        | 0.10000 | 17.2408  | 17.2408  | 17.2408  |
| 1,2,3,7,8,9-HxCDD   | 130           | 0.40        | 0.10000 | 13.3970  | 13.3970  | 13.3970  |
| Total HxCDD         | 2200          | 0.56        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,6,7,8-HpCDF | 2500          | 0.46        | 0.01000 | 25.1529  | 25.1529  | 25.1529  |
| 1,2,3,4,7,8,9-HpCDF | 130           | 0.38        | 0.01000 | 1.2781   | 1.2781   | 1.2781   |
| Total HpCDF         | 3100          | 0.42        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,6,7,8-HpCDD | 840           | 0.53        | 0.01000 | 8.4056   | 8.4056   | 8.4056   |
| Total HpCDD         | 1700          | 0.53        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| OCDF                | 910           | 0.35        | 0.00030 | 0.2722   | 0.2722   | 0.2722   |
| OCDD                | 2400          | 0.21        | 0.00030 | 0.7093   | 0.7093   | 0.7093   |

**710 ng/Kg      710 ng/Kg      710 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

**REPORT OF LABORATORY ANALYSIS**

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### Method 8290 Sample Analysis Results

Client - ESN Pacific

|                        |                           |           |                  |  |  |
|------------------------|---------------------------|-----------|------------------|--|--|
| Client's Sample ID     | RHS-DU-1.2                |           |                  |  |  |
| Lab Sample ID          | 10257346002-S             |           |                  |  |  |
| Filename               | P140301A_06               |           |                  |  |  |
| Injected By            | BAL                       |           |                  |  |  |
| Total Amount Extracted | 10.2 g                    | Matrix    | Soil             |  |  |
| % Moisture             | 8.5                       | Dilution  | NA               |  |  |
| Dry Weight Extracted   | 9.33 g                    | Collected | 02/05/2014 09:33 |  |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |  |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Extracted | 02/26/2014 21:30 |  |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 10:46 |  |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg |   | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|---|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 630        | ----       | 0.15     | E | 2,3,7,8-TCDF-13C         | 2.00       | 73               |
| Total TCDF          | 17000      | ----       | 0.15     | E | 2,3,7,8-TCDD-13C         | 2.00       | 77               |
|                     |            |            |          |   | 1,2,3,7,8-PeCDF-13C      | 2.00       | 71               |
| 2,3,7,8-TCDD        | 38         | ----       | 0.14     |   | 2,3,4,7,8-PeCDF-13C      | 2.00       | 69               |
| Total TCDD          | 1700       | ----       | 0.14     |   | 1,2,3,7,8-PeCDD-13C      | 2.00       | 71               |
|                     |            |            |          |   | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 85               |
| 1,2,3,7,8-PeCDF     | 470        | ----       | 0.43     |   | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 84               |
| 2,3,4,7,8-PeCDF     | 970        | ----       | 1.40     |   | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 81               |
| Total PeCDF         | 11000      | ----       | 0.92     |   | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 79               |
|                     |            |            |          |   | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 78               |
| 1,2,3,7,8-PeCDD     | 120        | ----       | 0.44     |   | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 67               |
| Total PeCDD         | 2000       | ----       | 0.44     |   | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 74               |
|                     |            |            |          |   | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 71               |
| 1,2,3,4,7,8-HxCDF   | 870        | ----       | 1.10     |   | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 76               |
| 1,2,3,6,7,8-HxCDF   | 740        | ----       | 1.40     |   | OCDD-13C                 | 4.00       | 79 Y             |
| 2,3,4,6,7,8-HxCDF   | 1000       | ----       | 0.82     |   |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 120        | ----       | 0.68     |   | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 8700       | ----       | 1.00     | E | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |   |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 110        | ----       | 0.54     |   | 2,3,7,8-TCDD-37Cl4       | 0.20       | 79               |
| 1,2,3,6,7,8-HxCDD   | 200        | ----       | 0.72     |   |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 150        | ----       | 0.83     |   |                          |            |                  |
| Total HxCDD         | 2700       | ----       | 0.70     |   |                          |            |                  |
|                     |            |            |          |   |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 3200       | ----       | 0.66     | E | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 100        | ----       | 0.81     |   | Equivalence: 890 ng/Kg   |            |                  |
| Total HpCDF         | 3900       | ----       | 0.73     | E | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |   |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 850        | ----       | 0.22     |   |                          |            |                  |
| Total HpCDD         | 1800       | ----       | 0.22     |   |                          |            |                  |
|                     |            |            |          |   |                          |            |                  |
| OCDF                | 560        | ----       | 0.27     |   |                          |            |                  |
| OCDD                | 1900       | ----       | 0.23     |   |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
E = Exceeds calibration range  
Y = Calculated using average of daily RfFs

## REPORT OF LABORATORY ANALYSIS

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-1.2                |           |                  |  |
| Lab Sample ID          | 10257346002-S             |           |                  |  |
| Filename               | P140301A_06               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.2 g                    | Matrix    | Soil             |  |
| % Moisture             | 8.5                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.33 g                    | Collected | 02/05/2014 09:33 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 10:46 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB       | MB       | UB       |
|---------------------|---------------|-------------|---------|----------|----------|----------|
| 2,3,7,8-TCDF        | 630           | 0.15        | 0.10000 | 63.0003  | 63.0003  | 63.0003  |
| Total TCDF          | 17000         | 0.15        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 2,3,7,8-TCDD        | 38            | 0.14        | 1.00000 | 38.2979  | 38.2979  | 38.2979  |
| Total TCDD          | 1700          | 0.14        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,7,8-PeCDF     | 470           | 0.43        | 0.03000 | 14.1460  | 14.1460  | 14.1460  |
| 2,3,4,7,8-PeCDF     | 970           | 1.4         | 0.30000 | 290.7745 | 290.7745 | 290.7745 |
| Total PeCDF         | 11000         | 0.92        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,7,8-PeCDD     | 120           | 0.44        | 1.00000 | 117.8814 | 117.8814 | 117.8814 |
| Total PeCDD         | 2000          | 0.44        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,7,8-HxCDF   | 870           | 1.1         | 0.10000 | 87.2673  | 87.2673  | 87.2673  |
| 1,2,3,6,7,8-HxCDF   | 740           | 1.4         | 0.10000 | 74.0624  | 74.0624  | 74.0624  |
| 2,3,4,6,7,8-HxCDF   | 1000          | 0.82        | 0.10000 | 101.9369 | 101.9369 | 101.9369 |
| 1,2,3,7,8,9-HxCDF   | 120           | 0.68        | 0.10000 | 12.1000  | 12.1000  | 12.1000  |
| Total HxCDF         | 8700          | 1.0         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,7,8-HxCDD   | 110           | 0.54        | 0.10000 | 10.6051  | 10.6051  | 10.6051  |
| 1,2,3,6,7,8-HxCDD   | 200           | 0.72        | 0.10000 | 19.9110  | 19.9110  | 19.9110  |
| 1,2,3,7,8,9-HxCDD   | 150           | 0.83        | 0.10000 | 15.4994  | 15.4994  | 15.4994  |
| Total HxCDD         | 2700          | 0.70        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,6,7,8-HpCDF | 3200          | 0.66        | 0.01000 | 32.3356  | 32.3356  | 32.3356  |
| 1,2,3,4,7,8,9-HpCDF | 100           | 0.81        | 0.01000 | 1.0486   | 1.0486   | 1.0486   |
| Total HpCDF         | 3900          | 0.73        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,6,7,8-HpCDD | 850           | 0.22        | 0.01000 | 8.4969   | 8.4969   | 8.4969   |
| Total HpCDD         | 1800          | 0.22        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| OCDF                | 560           | 0.27        | 0.00030 | 0.1688   | 0.1688   | 0.1688   |
| OCDD                | 1900          | 0.23        | 0.00030 | 0.5720   | 0.5720   | 0.5720   |

**890 ng/Kg      890 ng/Kg      890 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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### Method 8290 Sample Analysis Results

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-1.3                |           |                  |  |
| Lab Sample ID          | 10257346003-S             |           |                  |  |
| Filename               | P140228B_04               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.4 g                    | Matrix    | Soil             |  |
| % Moisture             | 9.6                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.40 g                    | Collected | 02/05/2014 10:30 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 02/28/2014 19:47 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg |    | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|----|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | -----      | 2700       | 0.40     | PE | 2,3,7,8-TCDF-13C         | 2.00       | 67               |
| Total TCDF          | 57000      | -----      | 0.40     | E  | 2,3,7,8-TCDD-13C         | 2.00       | 75               |
|                     |            |            |          |    | 1,2,3,7,8-PeCDF-13C      | 2.00       | 68               |
| 2,3,7,8-TCDD        | 140        | -----      | 0.66     |    | 2,3,4,7,8-PeCDF-13C      | 2.00       | 66               |
| Total TCDD          | 5700       | -----      | 0.66     | E  | 1,2,3,7,8-PeCDD-13C      | 2.00       | 72               |
|                     |            |            |          |    | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 78               |
| 1,2,3,7,8-PeCDF     | 1700       | -----      | 0.75     |    | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 77               |
| 2,3,4,7,8-PeCDF     | 3800       | -----      | 2.80     | E  | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 75               |
| Total PeCDF         | 38000      | -----      | 1.80     | E  | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 73               |
|                     |            |            |          |    | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 71               |
| 1,2,3,7,8-PeCDD     | 340        | -----      | 1.10     |    | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 65               |
| Total PeCDD         | 5500       | -----      | 1.10     |    | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 73               |
|                     |            |            |          |    | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 72               |
| 1,2,3,4,7,8-HxCDF   | 2300       | -----      | 1.10     | E  | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 77               |
| 1,2,3,6,7,8-HxCDF   | 2000       | -----      | 2.60     |    | OCDD-13C                 | 4.00       | 58               |
| 2,3,4,6,7,8-HxCDF   | 2600       | -----      | 1.80     | E  |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 310        | -----      | 2.00     |    | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 23000      | -----      | 1.90     | E  | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 260        | -----      | 1.40     |    | 2,3,7,8-TCDD-37Cl4       | 0.20       | 80               |
| 1,2,3,6,7,8-HxCDD   | 410        | -----      | 1.50     |    |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 370        | -----      | 1.50     |    |                          |            |                  |
| Total HxCDD         | 6200       | -----      | 1.50     | E  |                          |            |                  |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 7700       | -----      | 1.20     | E  | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 270        | -----      | 1.60     |    | Equivalence: 2900 ng/Kg  |            |                  |
| Total HpCDF         | 9400       | -----      | 1.40     | E  | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 2100       | -----      | 0.50     | E  |                          |            |                  |
| Total HpCDD         | 4600       | -----      | 0.50     | E  |                          |            |                  |
|                     |            |            |          |    |                          |            |                  |
| OCDF                | 1500       | -----      | 0.52     |    |                          |            |                  |
| OCDD                | 5100       | -----      | 0.42     | E  |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
P = PCDE Interference  
E = Exceeds calibration range

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-1.3                |           |                  |  |
| Lab Sample ID          | 10257346003-S             |           |                  |  |
| Filename               | P140228B_04               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.4 g                    | Matrix    | Soil             |  |
| % Moisture             | 9.6                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.40 g                    | Collected | 02/05/2014 10:30 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 02/28/2014 19:47 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB                | MB                | UB                |
|---------------------|---------------|-------------|---------|-------------------|-------------------|-------------------|
| 2,3,7,8-TCDF        | ND            | 0.40        | 0.10000 | 266.4271          | 266.4271          | 266.4271          |
| Total TCDF          | 57000         | 0.40        | 0.00000 | 0.0000            | 0.0000            | 0.0000            |
| 2,3,7,8-TCDD        | 140           | 0.66        | 1.00000 | 140.4906          | 140.4906          | 140.4906          |
| Total TCDD          | 5700          | 0.66        | 0.00000 | 0.0000            | 0.0000            | 0.0000            |
| 1,2,3,7,8-PeCDF     | 1700          | 0.75        | 0.03000 | 51.8696           | 51.8696           | 51.8696           |
| 2,3,4,7,8-PeCDF     | 3800          | 2.8         | 0.30000 | 1150.0767         | 1150.0767         | 1150.0767         |
| Total PeCDF         | 38000         | 1.8         | 0.00000 | 0.0000            | 0.0000            | 0.0000            |
| 1,2,3,7,8-PeCDD     | 340           | 1.1         | 1.00000 | 344.9645          | 344.9645          | 344.9645          |
| Total PeCDD         | 5500          | 1.1         | 0.00000 | 0.0000            | 0.0000            | 0.0000            |
| 1,2,3,4,7,8-HxCDF   | 2300          | 1.1         | 0.10000 | 225.4365          | 225.4365          | 225.4365          |
| 1,2,3,6,7,8-HxCDF   | 2000          | 2.6         | 0.10000 | 196.5160          | 196.5160          | 196.5160          |
| 2,3,4,6,7,8-HxCDF   | 2600          | 1.8         | 0.10000 | 264.4692          | 264.4692          | 264.4692          |
| 1,2,3,7,8,9-HxCDF   | 310           | 2.0         | 0.10000 | 30.7299           | 30.7299           | 30.7299           |
| Total HxCDF         | 23000         | 1.9         | 0.00000 | 0.0000            | 0.0000            | 0.0000            |
| 1,2,3,4,7,8-HxCDD   | 260           | 1.4         | 0.10000 | 25.8999           | 25.8999           | 25.8999           |
| 1,2,3,6,7,8-HxCDD   | 410           | 1.5         | 0.10000 | 41.2923           | 41.2923           | 41.2923           |
| 1,2,3,7,8,9-HxCDD   | 370           | 1.5         | 0.10000 | 36.6176           | 36.6176           | 36.6176           |
| Total HxCDD         | 6200          | 1.5         | 0.00000 | 0.0000            | 0.0000            | 0.0000            |
| 1,2,3,4,6,7,8-HpCDF | 7700          | 1.2         | 0.01000 | 77.4586           | 77.4586           | 77.4586           |
| 1,2,3,4,7,8,9-HpCDF | 270           | 1.6         | 0.01000 | 2.7123            | 2.7123            | 2.7123            |
| Total HpCDF         | 9400          | 1.4         | 0.00000 | 0.0000            | 0.0000            | 0.0000            |
| 1,2,3,4,6,7,8-HpCDD | 2100          | 0.50        | 0.01000 | 21.3697           | 21.3697           | 21.3697           |
| Total HpCDD         | 4600          | 0.50        | 0.00000 | 0.0000            | 0.0000            | 0.0000            |
| OCDF                | 1500          | 0.52        | 0.00030 | 0.4615            | 0.4615            | 0.4615            |
| OCDD                | 5100          | 0.42        | 0.00030 | 1.5290            | 1.5290            | 1.5290            |
|                     |               |             |         | <b>2900 ng/Kg</b> | <b>2900 ng/Kg</b> | <b>2900 ng/Kg</b> |

Final values are valid to only 2 significant figures  
 TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
 LB = Lower Bound, Where "ND", TEQ Conc = 0  
 MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
 UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
 RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |
|------------------------|---------------------------|-----------|------------------|
| Client's Sample ID     | RHS-DU-2                  |           |                  |
| Lab Sample ID          | 10257346004-S             |           |                  |
| Filename               | P140228B_05               |           |                  |
| Injected By            | BAL                       |           |                  |
| Total Amount Extracted | 10.5 g                    | Matrix    | Soil             |
| % Moisture             | 9.2                       | Dilution  | NA               |
| Dry Weight Extracted   | 9.53 g                    | Collected | 02/04/2014 15:00 |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |
| Method Blank ID        | BLANK-39465               | Analyzed  | 02/28/2014 20:30 |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 230        | ----       | 1.20     | 2,3,7,8-TCDF-13C         | 2.00       | 71               |
| Total TCDF          | 5800       | ----       | 1.20 E   | 2,3,7,8-TCDD-13C         | 2.00       | 80               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 68               |
| 2,3,7,8-TCDD        | 16         | ----       | 0.70     | 2,3,4,7,8-PeCDF-13C      | 2.00       | 66               |
| Total TCDD          | 480        | ----       | 0.70     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 71               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 84               |
| 1,2,3,7,8-PeCDF     | 170        | ----       | 3.30     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 84               |
| 2,3,4,7,8-PeCDF     | 330        | ----       | 3.10     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 79               |
| Total PeCDF         | 3600       | ----       | 3.20     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 73               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 82               |
| 1,2,3,7,8-PeCDD     | 43         | ----       | 1.40     | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 67               |
| Total PeCDD         | 540        | ----       | 1.40     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 70               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 66               |
| 1,2,3,4,7,8-HxCDF   | 220        | ----       | 2.70     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 72               |
| 1,2,3,6,7,8-HxCDF   | 210        | ----       | 1.50     | OCDD-13C                 | 4.00       | 52               |
| 2,3,4,6,7,8-HxCDF   | 270        | ----       | 1.70     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 44         | ----       | 2.70     | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 2400       | ----       | 2.20 E   | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 34         | ----       | 0.98     | 2,3,7,8-TCDD-37Cl4       | 0.20       | 78               |
| 1,2,3,6,7,8-HxCDD   | 68         | ----       | 1.10     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 53         | ----       | 0.90     |                          |            |                  |
| Total HxCDD         | 820        | ----       | 0.98     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 710        | ----       | 0.96     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 39         | ----       | 1.60     | Equivalence: 290 ng/Kg   |            |                  |
| Total HpCDF         | 1000       | ----       | 1.30     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 590        | ----       | 2.40     |                          |            |                  |
| Total HpCDD         | 1200       | ----       | 2.40     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 330        | ----       | 2.30     |                          |            |                  |
| OCDD                | 3700       | ----       | 4.50     |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
E = Exceeds calibration range

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-2                  |           |                  |  |
| Lab Sample ID          | 10257346004-S             |           |                  |  |
| Filename               | P140228B_05               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.5 g                    | Matrix    | Soil             |  |
| % Moisture             | 9.2                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.53 g                    | Collected | 02/04/2014 15:00 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 02/28/2014 20:30 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB      | MB      | UB      |
|---------------------|---------------|-------------|---------|---------|---------|---------|
| 2,3,7,8-TCDF        | 230           | 1.2         | 0.10000 | 22.7176 | 22.7176 | 22.7176 |
| Total TCDF          | 5800          | 1.2         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 2,3,7,8-TCDD        | 16            | 0.70        | 1.00000 | 16.4296 | 16.4296 | 16.4296 |
| Total TCDD          | 480           | 0.70        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDF     | 170           | 3.3         | 0.03000 | 5.1067  | 5.1067  | 5.1067  |
| 2,3,4,7,8-PeCDF     | 330           | 3.1         | 0.30000 | 97.9930 | 97.9930 | 97.9930 |
| Total PeCDF         | 3600          | 3.2         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDD     | 43            | 1.4         | 1.00000 | 42.7471 | 42.7471 | 42.7471 |
| Total PeCDD         | 540           | 1.4         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDF   | 220           | 2.7         | 0.10000 | 21.9228 | 21.9228 | 21.9228 |
| 1,2,3,6,7,8-HxCDF   | 210           | 1.5         | 0.10000 | 20.8312 | 20.8312 | 20.8312 |
| 2,3,4,6,7,8-HxCDF   | 270           | 1.7         | 0.10000 | 27.3254 | 27.3254 | 27.3254 |
| 1,2,3,7,8,9-HxCDF   | 44            | 2.7         | 0.10000 | 4.4244  | 4.4244  | 4.4244  |
| Total HxCDF         | 2400          | 2.2         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDD   | 34            | 0.98        | 0.10000 | 3.3965  | 3.3965  | 3.3965  |
| 1,2,3,6,7,8-HxCDD   | 68            | 1.1         | 0.10000 | 6.7560  | 6.7560  | 6.7560  |
| 1,2,3,7,8,9-HxCDD   | 53            | 0.90        | 0.10000 | 5.3329  | 5.3329  | 5.3329  |
| Total HxCDD         | 820           | 0.98        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDF | 710           | 0.96        | 0.01000 | 7.1142  | 7.1142  | 7.1142  |
| 1,2,3,4,7,8,9-HpCDF | 39            | 1.6         | 0.01000 | 0.3860  | 0.3860  | 0.3860  |
| Total HpCDF         | 1000          | 1.3         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDD | 590           | 2.4         | 0.01000 | 5.8672  | 5.8672  | 5.8672  |
| Total HpCDD         | 1200          | 2.4         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| OCDF                | 330           | 2.3         | 0.00030 | 0.0998  | 0.0998  | 0.0998  |
| OCDD                | 3700          | 4.5         | 0.00030 | 1.1246  | 1.1246  | 1.1246  |

**290 ng/Kg      290 ng/Kg      290 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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### Method 8290 Sample Analysis Results

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-3                  |           |                  |  |
| Lab Sample ID          | 10257346005-S             |           |                  |  |
| Filename               | P140228B_06               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.3 g                    | Matrix    | Soil             |  |
| % Moisture             | 10.8                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.19 g                    | Collected | 02/04/2014 14:03 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 02/28/2014 21:13 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 7.90       | ----       | 1.10     | 2,3,7,8-TCDF-13C         | 2.00       | 75               |
| Total TCDF          | 180.00     | ----       | 1.10     | 2,3,7,8-TCDD-13C         | 2.00       | 83               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 70               |
| 2,3,7,8-TCDD        | 0.75       | ----       | 0.63 J   | 2,3,4,7,8-PeCDF-13C      | 2.00       | 68               |
| Total TCDD          | 15.00      | ----       | 0.63     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 71               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 87               |
| 1,2,3,7,8-PeCDF     | 7.60       | ----       | 0.87     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 84               |
| 2,3,4,7,8-PeCDF     | 21.00      | ----       | 0.80     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 82               |
| Total PeCDF         | 230.00     | ----       | 0.83     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 75               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 83               |
| 1,2,3,7,8-PeCDD     | 2.50       | ----       | 0.90 J   | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 72               |
| Total PeCDD         | 22.00      | ----       | 0.90     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 72               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 69               |
| 1,2,3,4,7,8-HxCDF   | 15.00      | ----       | 0.80     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 73               |
| 1,2,3,6,7,8-HxCDF   | ----       | 26         | 0.71 P   | OCDD-13C                 | 4.00       | 54               |
| 2,3,4,6,7,8-HxCDF   | 20.00      | ----       | 0.75     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 3.20       | ----       | 0.99 J   | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 240.00     | ----       | 0.82     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 4.90       | ----       | 0.91 J   | 2,3,7,8-TCDD-37Cl4       | 0.20       | 82               |
| 1,2,3,6,7,8-HxCDD   | 12.00      | ----       | 1.00     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 7.80       | ----       | 0.94     |                          |            |                  |
| Total HxCDD         | 110.00     | ----       | 0.96     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 110.00     | ----       | 0.83     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 5.70       | ----       | 1.10     | Equivalence: 25 ng/Kg    |            |                  |
| Total HpCDF         | 310.00     | ----       | 0.98     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 290.00     | ----       | 1.70     |                          |            |                  |
| Total HpCDD         | 570.00     | ----       | 1.70     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 250.00     | ----       | 0.89     |                          |            |                  |
| OCDD                | 4100.00    | ----       | 1.20     |                          |            |                  |

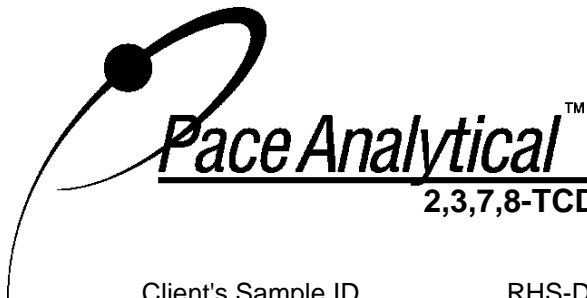
Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
J = Estimated value  
P = PCDE Interference

## REPORT OF LABORATORY ANALYSIS

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-3                  |           |                  |  |
| Lab Sample ID          | 10257346005-S             |           |                  |  |
| Filename               | P140228B_06               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.3 g                    | Matrix    | Soil             |  |
| % Moisture             | 10.8                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.19 g                    | Collected | 02/04/2014 14:03 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 02/28/2014 21:13 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | 7.90          | 1.1         | 0.10000 | 0.7872 | 0.7872 | 0.7872 |
| Total TCDF          | 180.00        | 1.1         | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | 0.75          | 0.63        | 1.00000 | 0.7474 | 0.7474 | 0.7474 |
| Total TCDD          | 15.00         | 0.63        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | 7.60          | 0.87        | 0.03000 | 0.2293 | 0.2293 | 0.2293 |
| 2,3,4,7,8-PeCDF     | 21.00         | 0.80        | 0.30000 | 6.2023 | 6.2023 | 6.2023 |
| Total PeCDF         | 230.00        | 0.83        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | 2.50          | 0.90        | 1.00000 | 2.4729 | 2.4729 | 2.4729 |
| Total PeCDD         | 22.00         | 0.90        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | 15.00         | 0.80        | 0.10000 | 1.4750 | 1.4750 | 1.4750 |
| 1,2,3,6,7,8-HxCDF   | ND            | 0.71        | 0.10000 | 2.6405 | 2.6405 | 2.6405 |
| 2,3,4,6,7,8-HxCDF   | 20.00         | 0.75        | 0.10000 | 1.9850 | 1.9850 | 1.9850 |
| 1,2,3,7,8,9-HxCDF   | 3.20          | 0.99        | 0.10000 | 0.3170 | 0.3170 | 0.3170 |
| Total HxCDF         | 240.00        | 0.82        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | 4.90          | 0.91        | 0.10000 | 0.4950 | 0.4950 | 0.4950 |
| 1,2,3,6,7,8-HxCDD   | 12.00         | 1.0         | 0.10000 | 1.2189 | 1.2189 | 1.2189 |
| 1,2,3,7,8,9-HxCDD   | 7.80          | 0.94        | 0.10000 | 0.7800 | 0.7800 | 0.7800 |
| Total HxCDD         | 110.00        | 0.96        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | 110.00        | 0.83        | 0.01000 | 1.1335 | 1.1335 | 1.1335 |
| 1,2,3,4,7,8,9-HpCDF | 5.70          | 1.1         | 0.01000 | 0.0565 | 0.0565 | 0.0565 |
| Total HpCDF         | 310.00        | 0.98        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | 290.00        | 1.7         | 0.01000 | 2.8638 | 2.8638 | 2.8638 |
| Total HpCDD         | 570.00        | 1.7         | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | 250.00        | 0.89        | 0.00030 | 0.0764 | 0.0764 | 0.0764 |
| OCDD                | 4100.00       | 1.2         | 0.00030 | 1.2196 | 1.2196 | 1.2196 |

**25 ng/Kg      25 ng/Kg      25 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |
|------------------------|---------------------------|-----------|------------------|
| Client's Sample ID     | RHS-DU-4                  |           |                  |
| Lab Sample ID          | 10257346006-S             |           |                  |
| Filename               | P140228B_07               |           |                  |
| Injected By            | BAL                       |           |                  |
| Total Amount Extracted | 10.4 g                    | Matrix    | Soil             |
| % Moisture             | 16.0                      | Dilution  | NA               |
| Dry Weight Extracted   | 8.74 g                    | Collected | 02/04/2014 11:45 |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |
| Method Blank ID        | BLANK-39465               | Analyzed  | 02/28/2014 21:56 |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 16.00      | ----       | 0.15     | 2,3,7,8-TCDF-13C         | 2.00       | 70               |
| Total TCDF          | 410.00     | ----       | 0.15     | 2,3,7,8-TCDD-13C         | 2.00       | 78               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 66               |
| 2,3,7,8-TCDD        | 0.99       | ----       | 0.19 J   | 2,3,4,7,8-PeCDF-13C      | 2.00       | 64               |
| Total TCDD          | 37.00      | ----       | 0.19     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 69               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 81               |
| 1,2,3,7,8-PeCDF     | 13.00      | ----       | 0.45     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 80               |
| 2,3,4,7,8-PeCDF     | 26.00      | ----       | 0.33     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 76               |
| Total PeCDF         | 310.00     | ----       | 0.39     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 67               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 76               |
| 1,2,3,7,8-PeCDD     | 4.30       | ----       | 0.46 J   | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 67               |
| Total PeCDD         | 57.00      | ----       | 0.46     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 67               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 66               |
| 1,2,3,4,7,8-HxCDF   | 22.00      | ----       | 0.25     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 71               |
| 1,2,3,6,7,8-HxCDF   | 20.00      | ----       | 0.21     | OCDD-13C                 | 4.00       | 54               |
| 2,3,4,6,7,8-HxCDF   | 29.00      | ----       | 0.22     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 3.90       | ----       | 0.28 J   | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 310.00     | ----       | 0.24     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 5.40       | ----       | 0.28 J   | 2,3,7,8-TCDD-37Cl4       | 0.20       | 78               |
| 1,2,3,6,7,8-HxCDD   | 31.00      | ----       | 0.28     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 17.00      | ----       | 0.24     |                          |            |                  |
| Total HxCDD         | 240.00     | ----       | 0.27     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 170.00     | ----       | 0.26     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 7.50       | ----       | 0.35     | Equivalence: 35 ng/Kg    |            |                  |
| Total HpCDF         | 370.00     | ----       | 0.31     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 360.00     | ----       | 0.83     |                          |            |                  |
| Total HpCDD         | 700.00     | ----       | 0.83     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 310.00     | ----       | 0.54     |                          |            |                  |
| OCDD                | 3400.00    | ----       | 1.70     |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
J = Estimated value

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-4                  |           |                  |  |
| Lab Sample ID          | 10257346006-S             |           |                  |  |
| Filename               | P140228B_07               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.4 g                    | Matrix    | Soil             |  |
| % Moisture             | 16.0                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 8.74 g                    | Collected | 02/04/2014 11:45 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 02/28/2014 21:56 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | 16.00         | 0.15        | 0.10000 | 1.6233 | 1.6233 | 1.6233 |
| Total TCDF          | 410.00        | 0.15        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | 0.99          | 0.19        | 1.00000 | 0.9907 | 0.9907 | 0.9907 |
| Total TCDD          | 37.00         | 0.19        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | 13.00         | 0.45        | 0.03000 | 0.3785 | 0.3785 | 0.3785 |
| 2,3,4,7,8-PeCDF     | 26.00         | 0.33        | 0.30000 | 7.9467 | 7.9467 | 7.9467 |
| Total PeCDF         | 310.00        | 0.39        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | 4.30          | 0.46        | 1.00000 | 4.3193 | 4.3193 | 4.3193 |
| Total PeCDD         | 57.00         | 0.46        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | 22.00         | 0.25        | 0.10000 | 2.2258 | 2.2258 | 2.2258 |
| 1,2,3,6,7,8-HxCDF   | 20.00         | 0.21        | 0.10000 | 2.0104 | 2.0104 | 2.0104 |
| 2,3,4,6,7,8-HxCDF   | 29.00         | 0.22        | 0.10000 | 2.8826 | 2.8826 | 2.8826 |
| 1,2,3,7,8,9-HxCDF   | 3.90          | 0.28        | 0.10000 | 0.3949 | 0.3949 | 0.3949 |
| Total HxCDF         | 310.00        | 0.24        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | 5.40          | 0.28        | 0.10000 | 0.5421 | 0.5421 | 0.5421 |
| 1,2,3,6,7,8-HxCDD   | 31.00         | 0.28        | 0.10000 | 3.1137 | 3.1137 | 3.1137 |
| 1,2,3,7,8,9-HxCDD   | 17.00         | 0.24        | 0.10000 | 1.7111 | 1.7111 | 1.7111 |
| Total HxCDD         | 240.00        | 0.27        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | 170.00        | 0.26        | 0.01000 | 1.6818 | 1.6818 | 1.6818 |
| 1,2,3,4,7,8,9-HpCDF | 7.50          | 0.35        | 0.01000 | 0.0753 | 0.0753 | 0.0753 |
| Total HpCDF         | 370.00        | 0.31        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | 360.00        | 0.83        | 0.01000 | 3.5904 | 3.5904 | 3.5904 |
| Total HpCDD         | 700.00        | 0.83        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | 310.00        | 0.54        | 0.00030 | 0.0922 | 0.0922 | 0.0922 |
| OCDD                | 3400.00       | 1.7         | 0.00030 | 1.0149 | 1.0149 | 1.0149 |

**35 ng/Kg      35 ng/Kg      35 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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### Method 8290 Sample Analysis Results

Client - ESN Pacific

|                        |                           |           |                  |
|------------------------|---------------------------|-----------|------------------|
| Client's Sample ID     | RHS-DU-5                  |           |                  |
| Lab Sample ID          | 10257346007-S             |           |                  |
| Filename               | P140228B_08               |           |                  |
| Injected By            | BAL                       |           |                  |
| Total Amount Extracted | 10.3 g                    | Matrix    | Soil             |
| % Moisture             | 13.3                      | Dilution  | NA               |
| Dry Weight Extracted   | 8.93 g                    | Collected | 02/04/2014 12:20 |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |
| Method Blank ID        | BLANK-39465               | Analyzed  | 02/28/2014 22:38 |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 10.00      | ----       | 0.11     | 2,3,7,8-TCDF-13C         | 2.00       | 75               |
| Total TCDF          | 210.00     | ----       | 0.11     | 2,3,7,8-TCDD-13C         | 2.00       | 83               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 72               |
| 2,3,7,8-TCDD        | 0.78       | ----       | 0.15 J   | 2,3,4,7,8-PeCDF-13C      | 2.00       | 69               |
| Total TCDD          | 24.00      | ----       | 0.15     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 75               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 88               |
| 1,2,3,7,8-PeCDF     | 10.00      | ----       | 0.18     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 87               |
| 2,3,4,7,8-PeCDF     | 22.00      | ----       | 0.19     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 81               |
| Total PeCDF         | 250.00     | ----       | 0.18     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 65               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 82               |
| 1,2,3,7,8-PeCDD     | 4.10       | ----       | 0.14 J   | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 67               |
| Total PeCDD         | 51.00      | ----       | 0.14     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 71               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 68               |
| 1,2,3,4,7,8-HxCDF   | 22.00      | ----       | 0.29     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 72               |
| 1,2,3,6,7,8-HxCDF   | 25.00      | ----       | 0.26     | OCDD-13C                 | 4.00       | 54               |
| 2,3,4,6,7,8-HxCDF   | 29.00      | ----       | 0.24     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 3.80       | ----       | 0.37 J   | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 350.00     | ----       | 0.29     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 6.50       | ----       | 0.26     | 2,3,7,8-TCDD-37Cl4       | 0.20       | 86               |
| 1,2,3,6,7,8-HxCDD   | 22.00      | ----       | 0.28     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 14.00      | ----       | 0.22     |                          |            |                  |
| Total HxCDD         | 190.00     | ----       | 0.25     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 210.00     | ----       | 0.26     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 8.90       | ----       | 0.33     | Equivalence: 33 ng/Kg    |            |                  |
| Total HpCDF         | 510.00     | ----       | 0.29     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 460.00     | ----       | 0.77     |                          |            |                  |
| Total HpCDD         | 890.00     | ----       | 0.77     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 430.00     | ----       | 0.32     |                          |            |                  |
| OCDD                | 4200.00    | ----       | 0.27     |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
J = Estimated value

## REPORT OF LABORATORY ANALYSIS

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-5                  |           |                  |  |
| Lab Sample ID          | 10257346007-S             |           |                  |  |
| Filename               | P140228B_08               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.3 g                    | Matrix    | Soil             |  |
| % Moisture             | 13.3                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 8.93 g                    | Collected | 02/04/2014 12:20 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 02/28/2014 22:38 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB              | MB              | UB              |
|---------------------|---------------|-------------|---------|-----------------|-----------------|-----------------|
| 2,3,7,8-TCDF        | 10.00         | 0.11        | 0.10000 | 1.0134          | 1.0134          | 1.0134          |
| Total TCDF          | 210.00        | 0.11        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 2,3,7,8-TCDD        | 0.78          | 0.15        | 1.00000 | 0.7805          | 0.7805          | 0.7805          |
| Total TCDD          | 24.00         | 0.15        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,7,8-PeCDF     | 10.00         | 0.18        | 0.03000 | 0.3104          | 0.3104          | 0.3104          |
| 2,3,4,7,8-PeCDF     | 22.00         | 0.19        | 0.30000 | 6.6764          | 6.6764          | 6.6764          |
| Total PeCDF         | 250.00        | 0.18        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,7,8-PeCDD     | 4.10          | 0.14        | 1.00000 | 4.0811          | 4.0811          | 4.0811          |
| Total PeCDD         | 51.00         | 0.14        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,4,7,8-HxCDF   | 22.00         | 0.29        | 0.10000 | 2.1995          | 2.1995          | 2.1995          |
| 1,2,3,6,7,8-HxCDF   | 25.00         | 0.26        | 0.10000 | 2.5369          | 2.5369          | 2.5369          |
| 2,3,4,6,7,8-HxCDF   | 29.00         | 0.24        | 0.10000 | 2.9204          | 2.9204          | 2.9204          |
| 1,2,3,7,8,9-HxCDF   | 3.80          | 0.37        | 0.10000 | 0.3819          | 0.3819          | 0.3819          |
| Total HxCDF         | 350.00        | 0.29        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,4,7,8-HxCDD   | 6.50          | 0.26        | 0.10000 | 0.6510          | 0.6510          | 0.6510          |
| 1,2,3,6,7,8-HxCDD   | 22.00         | 0.28        | 0.10000 | 2.1999          | 2.1999          | 2.1999          |
| 1,2,3,7,8,9-HxCDD   | 14.00         | 0.22        | 0.10000 | 1.3581          | 1.3581          | 1.3581          |
| Total HxCDD         | 190.00        | 0.25        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,4,6,7,8-HpCDF | 210.00        | 0.26        | 0.01000 | 2.0881          | 2.0881          | 2.0881          |
| 1,2,3,4,7,8,9-HpCDF | 8.90          | 0.33        | 0.01000 | 0.0887          | 0.0887          | 0.0887          |
| Total HpCDF         | 510.00        | 0.29        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,4,6,7,8-HpCDD | 460.00        | 0.77        | 0.01000 | 4.5638          | 4.5638          | 4.5638          |
| Total HpCDD         | 890.00        | 0.77        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| OCDF                | 430.00        | 0.32        | 0.00030 | 0.1277          | 0.1277          | 0.1277          |
| OCDD                | 4200.00       | 0.27        | 0.00030 | 1.2722          | 1.2722          | 1.2722          |
|                     |               |             |         | <b>33 ng/Kg</b> | <b>33 ng/Kg</b> | <b>33 ng/Kg</b> |

Final values are valid to only 2 significant figures  
 TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
 LB = Lower Bound, Where "ND", TEQ Conc = 0  
 MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
 UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
 RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-6                  |           |                  |  |
| Lab Sample ID          | 10257346008-S             |           |                  |  |
| Filename               | U140226B_15               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.1 g                    | Matrix    | Soil             |  |
| % Moisture             | 14.3                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 8.66 g                    | Collected | 02/04/2014 14:02 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140226B_06 & U140226B_20 | Extracted | 02/24/2014 21:00 |  |
| Method Blank ID        | BLANK-39446               | Analyzed  | 02/27/2014 04:27 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 22.0       | ----       | 0.560    | 2,3,7,8-TCDF-13C         | 2.00       | 82 Y             |
| Total TCDF          | 600.0      | ----       | 0.560    | 2,3,7,8-TCDD-13C         | 2.00       | 88               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 89 Y             |
| 2,3,7,8-TCDD        | 1.4        | ----       | 0.098 Y  | 2,3,4,7,8-PeCDF-13C      | 2.00       | 86 Y             |
| Total TCDD          | 55.0       | ----       | 0.098 Y  | 1,2,3,7,8-PeCDD-13C      | 2.00       | 87 Y             |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 100              |
| 1,2,3,7,8-PeCDF     | 22.0       | ----       | 0.270    | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 103              |
| 2,3,4,7,8-PeCDF     | 46.0       | ----       | 0.230    | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 97               |
| Total PeCDF         | 570.0      | ----       | 0.250    | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 91               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 81               |
| 1,2,3,7,8-PeCDD     | 9.3        | ----       | 0.240    | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 75               |
| Total PeCDD         | 120.0      | ----       | 0.240    | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 64               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 60               |
| 1,2,3,4,7,8-HxCDF   | 42.0       | ----       | 0.280    | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 59               |
| 1,2,3,6,7,8-HxCDF   | ----       | 43         | 0.200 P  | OCDD-13C                 | 4.00       | 67               |
| 2,3,4,6,7,8-HxCDF   | 56.0       | ----       | 0.240    |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 8.8        | ----       | 0.440    | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 500.0      | ----       | 0.290    | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 13.0       | ----       | 0.460    | 2,3,7,8-TCDD-37Cl4       | 0.20       | 96               |
| 1,2,3,6,7,8-HxCDD   | 42.0       | ----       | 0.580    |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 27.0       | ----       | 0.600    |                          |            |                  |
| Total HxCDD         | 390.0      | ----       | 0.550    |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 240.0      | ----       | 0.200    | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 16.0       | ----       | 0.400    | Equivalence: 63 ng/Kg    |            |                  |
| Total HpCDF         | 630.0      | ----       | 0.300    | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 740.0      | ----       | 0.140    |                          |            |                  |
| Total HpCDD         | 1400.0     | ----       | 0.140    |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 570.0      | ----       | 0.460 Y  |                          |            |                  |
| OCDD                | 6400.0     | ----       | 0.440    |                          |            |                  |

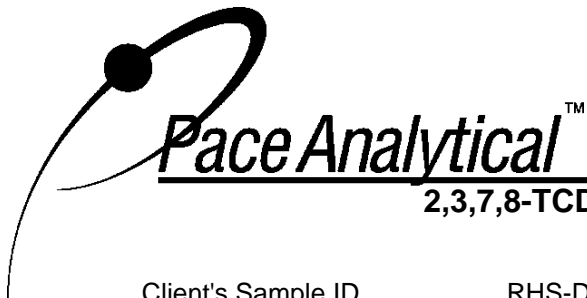
Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
P = PCDE Interference  
Y = Calculated using average of daily RfFs

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-6                  |           |                  |  |
| Lab Sample ID          | 10257346008-S             |           |                  |  |
| Filename               | U140226B_15               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.1 g                    | Matrix    | Soil             |  |
| % Moisture             | 14.3                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 8.66 g                    | Collected | 02/04/2014 14:02 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140226B_06 & U140226B_20 | Extracted | 02/24/2014 21:00 |  |
| Method Blank ID        | BLANK-39446               | Analyzed  | 02/27/2014 04:27 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB      | MB      | UB      |
|---------------------|---------------|-------------|---------|---------|---------|---------|
| 2,3,7,8-TCDF        | 22.0          | 0.56        | 0.10000 | 2.2016  | 2.2016  | 2.2016  |
| Total TCDF          | 600.0         | 0.56        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 2,3,7,8-TCDD        | 1.4           | 0.098       | 1.00000 | 1.4168  | 1.4168  | 1.4168  |
| Total TCDD          | 55.0          | 0.098       | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDF     | 22.0          | 0.27        | 0.03000 | 0.6533  | 0.6533  | 0.6533  |
| 2,3,4,7,8-PeCDF     | 46.0          | 0.23        | 0.30000 | 13.8748 | 13.8748 | 13.8748 |
| Total PeCDF         | 570.0         | 0.25        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDD     | 9.3           | 0.24        | 1.00000 | 9.3190  | 9.3190  | 9.3190  |
| Total PeCDD         | 120.0         | 0.24        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDF   | 42.0          | 0.28        | 0.10000 | 4.1906  | 4.1906  | 4.1906  |
| 1,2,3,6,7,8-HxCDF   | ND            | 0.20        | 0.10000 | 4.2820  | 4.2820  | 4.2820  |
| 2,3,4,6,7,8-HxCDF   | 56.0          | 0.24        | 0.10000 | 5.5569  | 5.5569  | 5.5569  |
| 1,2,3,7,8,9-HxCDF   | 8.8           | 0.44        | 0.10000 | 0.8820  | 0.8820  | 0.8820  |
| Total HxCDF         | 500.0         | 0.29        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDD   | 13.0          | 0.46        | 0.10000 | 1.2841  | 1.2841  | 1.2841  |
| 1,2,3,6,7,8-HxCDD   | 42.0          | 0.58        | 0.10000 | 4.2335  | 4.2335  | 4.2335  |
| 1,2,3,7,8,9-HxCDD   | 27.0          | 0.60        | 0.10000 | 2.6887  | 2.6887  | 2.6887  |
| Total HxCDD         | 390.0         | 0.55        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDF | 240.0         | 0.20        | 0.01000 | 2.4353  | 2.4353  | 2.4353  |
| 1,2,3,4,7,8,9-HpCDF | 16.0          | 0.40        | 0.01000 | 0.1649  | 0.1649  | 0.1649  |
| Total HpCDF         | 630.0         | 0.30        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDD | 740.0         | 0.14        | 0.01000 | 7.4428  | 7.4428  | 7.4428  |
| Total HpCDD         | 1400.0        | 0.14        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| OCDF                | 570.0         | 0.46        | 0.00030 | 0.1702  | 0.1702  | 0.1702  |
| OCDD                | 6400.0        | 0.44        | 0.00030 | 1.9249  | 1.9249  | 1.9249  |

**63 ng/Kg      63 ng/Kg      63 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |
|------------------------|---------------------------|-----------|------------------|
| Client's Sample ID     | RHS-DU-7                  |           |                  |
| Lab Sample ID          | 10257346009-S             |           |                  |
| Filename               | P140228B_09               |           |                  |
| Injected By            | BAL                       |           |                  |
| Total Amount Extracted | 10.2 g                    | Matrix    | Soil             |
| % Moisture             | 10.2                      | Dilution  | NA               |
| Dry Weight Extracted   | 9.16 g                    | Collected | 02/04/2014 11:31 |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |
| Method Blank ID        | BLANK-39465               | Analyzed  | 02/28/2014 23:21 |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 42.0       | ----       | 0.17     | 2,3,7,8-TCDF-13C         | 2.00       | 72               |
| Total TCDF          | 1100.0     | ----       | 0.17     | 2,3,7,8-TCDD-13C         | 2.00       | 80               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 70               |
| 2,3,7,8-TCDD        | 2.6        | ----       | 0.21     | 2,3,4,7,8-PeCDF-13C      | 2.00       | 67               |
| Total TCDD          | 100.0      | ----       | 0.21     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 72               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 85               |
| 1,2,3,7,8-PeCDF     | 46.0       | ----       | 0.26     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 81               |
| 2,3,4,7,8-PeCDF     | 98.0       | ----       | 0.55     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 78               |
| Total PeCDF         | 1100.0     | ----       | 0.40     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 49               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 80               |
| 1,2,3,7,8-PeCDD     | 15.0       | ----       | 0.25     | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 64               |
| Total PeCDD         | 200.0      | ----       | 0.25     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 68               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 65               |
| 1,2,3,4,7,8-HxCDF   | 100.0      | ----       | 0.50     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 71               |
| 1,2,3,6,7,8-HxCDF   | 88.0       | ----       | 0.54     | OCDD-13C                 | 4.00       | 55               |
| 2,3,4,6,7,8-HxCDF   | 130.0      | ----       | 0.46     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 18.0       | ----       | 0.47     | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 1300.0     | ----       | 0.49     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 19.0       | ----       | 0.44     | 2,3,7,8-TCDD-37Cl4       | 0.20       | 86               |
| 1,2,3,6,7,8-HxCDD   | 55.0       | ----       | 0.47     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 36.0       | ----       | 0.45     |                          |            |                  |
| Total HxCDD         | 600.0      | ----       | 0.45     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 600.0      | ----       | 0.42     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 25.0       | ----       | 0.45     | Equivalence: 110 ng/Kg   |            |                  |
| Total HpCDF         | 1000.0     | ----       | 0.43     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 800.0      | ----       | 0.96     |                          |            |                  |
| Total HpCDD         | 1700.0     | ----       | 0.96     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 570.0      | ----       | 0.46     |                          |            |                  |
| OCDD                | 6300.0     | ----       | 2.10     |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-7                  |           |                  |  |
| Lab Sample ID          | 10257346009-S             |           |                  |  |
| Filename               | P140228B_09               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.2 g                    | Matrix    | Soil             |  |
| % Moisture             | 10.2                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.16 g                    | Collected | 02/04/2014 11:31 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 02/28/2014 23:21 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB      | MB      | UB      |
|---------------------|---------------|-------------|---------|---------|---------|---------|
| 2,3,7,8-TCDF        | 42.0          | 0.17        | 0.10000 | 4.2120  | 4.2120  | 4.2120  |
| Total TCDF          | 1100.0        | 0.17        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 2,3,7,8-TCDD        | 2.6           | 0.21        | 1.00000 | 2.5779  | 2.5779  | 2.5779  |
| Total TCDD          | 100.0         | 0.21        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDF     | 46.0          | 0.26        | 0.03000 | 1.3924  | 1.3924  | 1.3924  |
| 2,3,4,7,8-PeCDF     | 98.0          | 0.55        | 0.30000 | 29.4800 | 29.4800 | 29.4800 |
| Total PeCDF         | 1100.0        | 0.40        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDD     | 15.0          | 0.25        | 1.00000 | 14.6872 | 14.6872 | 14.6872 |
| Total PeCDD         | 200.0         | 0.25        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDF   | 100.0         | 0.50        | 0.10000 | 10.2795 | 10.2795 | 10.2795 |
| 1,2,3,6,7,8-HxCDF   | 88.0          | 0.54        | 0.10000 | 8.7543  | 8.7543  | 8.7543  |
| 2,3,4,6,7,8-HxCDF   | 130.0         | 0.46        | 0.10000 | 13.1480 | 13.1480 | 13.1480 |
| 1,2,3,7,8,9-HxCDF   | 18.0          | 0.47        | 0.10000 | 1.8134  | 1.8134  | 1.8134  |
| Total HxCDF         | 1300.0        | 0.49        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDD   | 19.0          | 0.44        | 0.10000 | 1.9415  | 1.9415  | 1.9415  |
| 1,2,3,6,7,8-HxCDD   | 55.0          | 0.47        | 0.10000 | 5.5172  | 5.5172  | 5.5172  |
| 1,2,3,7,8,9-HxCDD   | 36.0          | 0.45        | 0.10000 | 3.5872  | 3.5872  | 3.5872  |
| Total HxCDD         | 600.0         | 0.45        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDF | 600.0         | 0.42        | 0.01000 | 6.0332  | 6.0332  | 6.0332  |
| 1,2,3,4,7,8,9-HpCDF | 25.0          | 0.45        | 0.01000 | 0.2528  | 0.2528  | 0.2528  |
| Total HpCDF         | 1000.0        | 0.43        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDD | 800.0         | 0.96        | 0.01000 | 8.0212  | 8.0212  | 8.0212  |
| Total HpCDD         | 1700.0        | 0.96        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| OCDF                | 570.0         | 0.46        | 0.00030 | 0.1720  | 0.1720  | 0.1720  |
| OCDD                | 6300.0        | 2.1         | 0.00030 | 1.8777  | 1.8777  | 1.8777  |

**110 ng/Kg      110 ng/Kg      110 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |
|------------------------|---------------------------|-----------|------------------|
| Client's Sample ID     | RHS-DU-8                  |           |                  |
| Lab Sample ID          | 10257346010-S             |           |                  |
| Filename               | P140228B_10               |           |                  |
| Injected By            | BAL                       |           |                  |
| Total Amount Extracted | 10.7 g                    | Matrix    | Soil             |
| % Moisture             | 12.6                      | Dilution  | NA               |
| Dry Weight Extracted   | 9.35 g                    | Collected | 02/04/2014 14:24 |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 00:04 |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 73.0       | ----       | 0.18     | 2,3,7,8-TCDF-13C         | 2.00       | 75               |
| Total TCDF          | 1800.0     | ----       | 0.18     | 2,3,7,8-TCDD-13C         | 2.00       | 82               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 76               |
| 2,3,7,8-TCDD        | 3.9        | ----       | 0.19     | 2,3,4,7,8-PeCDF-13C      | 2.00       | 71               |
| Total TCDD          | 170.0      | ----       | 0.19     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 77               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 107              |
| 1,2,3,7,8-PeCDF     | 73.0       | ----       | 0.21     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 90               |
| 2,3,4,7,8-PeCDF     | 160.0      | ----       | 0.23     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 84               |
| Total PeCDF         | 1600.0     | ----       | 0.22     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 45               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 88               |
| 1,2,3,7,8-PeCDD     | 20.0       | ----       | 0.24     | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 70               |
| Total PeCDD         | 290.0      | ----       | 0.24     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 73               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 64               |
| 1,2,3,4,7,8-HxCDF   | 130.0      | ----       | 0.49     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 70               |
| 1,2,3,6,7,8-HxCDF   | 130.0      | ----       | 0.39     | OCDD-13C                 | 4.00       | 45               |
| 2,3,4,6,7,8-HxCDF   | 200.0      | ----       | 0.44     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 27.0       | ----       | 0.53     | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 1700.0     | ----       | 0.46     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 26.0       | ----       | 0.52     | 2,3,7,8-TCDD-37Cl4       | 0.20       | 88               |
| 1,2,3,6,7,8-HxCDD   | 61.0       | ----       | 0.42     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 42.0       | ----       | 0.31     |                          |            |                  |
| Total HxCDD         | 770.0      | ----       | 0.42     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 690.0      | ----       | 0.43     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 32.0       | ----       | 0.25     | Equivalence: 160 ng/Kg   |            |                  |
| Total HpCDF         | 1000.0     | ----       | 0.34     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 640.0      | ----       | 0.88     |                          |            |                  |
| Total HpCDD         | 1300.0     | ----       | 0.88     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 370.0      | ----       | 0.50     |                          |            |                  |
| OCDD                | 4100.0     | ----       | 1.50     |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

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## 2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations

ESN Pacific

|                        |                           |           |                  |
|------------------------|---------------------------|-----------|------------------|
| Client's Sample ID     | RHS-DU-8                  |           |                  |
| Lab Sample ID          | 10257346010-S             |           |                  |
| Filename               | P140228B_10               |           |                  |
| Injected By            | BAL                       |           |                  |
| Total Amount Extracted | 10.7 g                    | Matrix    | Soil             |
| % Moisture             | 12.6                      | Dilution  | NA               |
| Dry Weight Extracted   | 9.35 g                    | Collected | 02/04/2014 14:24 |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 00:04 |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB      | MB      | UB      |
|---------------------|---------------|-------------|---------|---------|---------|---------|
| 2,3,7,8-TCDF        | 73.0          | 0.18        | 0.10000 | 7.3131  | 7.3131  | 7.3131  |
| Total TCDF          | 1800.0        | 0.18        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 2,3,7,8-TCDD        | 3.9           | 0.19        | 1.00000 | 3.8969  | 3.8969  | 3.8969  |
| Total TCDD          | 170.0         | 0.19        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDF     | 73.0          | 0.21        | 0.03000 | 2.2024  | 2.2024  | 2.2024  |
| 2,3,4,7,8-PeCDF     | 160.0         | 0.23        | 0.30000 | 46.9959 | 46.9959 | 46.9959 |
| Total PeCDF         | 1600.0        | 0.22        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDD     | 20.0          | 0.24        | 1.00000 | 20.2905 | 20.2905 | 20.2905 |
| Total PeCDD         | 290.0         | 0.24        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDF   | 130.0         | 0.49        | 0.10000 | 12.5722 | 12.5722 | 12.5722 |
| 1,2,3,6,7,8-HxCDF   | 130.0         | 0.39        | 0.10000 | 13.0742 | 13.0742 | 13.0742 |
| 2,3,4,6,7,8-HxCDF   | 200.0         | 0.44        | 0.10000 | 19.7418 | 19.7418 | 19.7418 |
| 1,2,3,7,8,9-HxCDF   | 27.0          | 0.53        | 0.10000 | 2.7032  | 2.7032  | 2.7032  |
| Total HxCDF         | 1700.0        | 0.46        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDD   | 26.0          | 0.52        | 0.10000 | 2.5672  | 2.5672  | 2.5672  |
| 1,2,3,6,7,8-HxCDD   | 61.0          | 0.42        | 0.10000 | 6.0860  | 6.0860  | 6.0860  |
| 1,2,3,7,8,9-HxCDD   | 42.0          | 0.31        | 0.10000 | 4.2065  | 4.2065  | 4.2065  |
| Total HxCDD         | 770.0         | 0.42        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDF | 690.0         | 0.43        | 0.01000 | 6.8828  | 6.8828  | 6.8828  |
| 1,2,3,4,7,8,9-HpCDF | 32.0          | 0.25        | 0.01000 | 0.3223  | 0.3223  | 0.3223  |
| Total HpCDF         | 1000.0        | 0.34        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDD | 640.0         | 0.88        | 0.01000 | 6.3833  | 6.3833  | 6.3833  |
| Total HpCDD         | 1300.0        | 0.88        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| OCDF                | 370.0         | 0.50        | 0.00030 | 0.1096  | 0.1096  | 0.1096  |
| OCDD                | 4100.0        | 1.5         | 0.00030 | 1.2413  | 1.2413  | 1.2413  |

**160 ng/Kg      160 ng/Kg      160 ng/Kg**

Final values are valid to only 2 significant figures  
 TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
 LB = Lower Bound, Where "ND", TEQ Conc = 0  
 MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
 UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
 RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-9                  |           |                  |  |
| Lab Sample ID          | 10257346011-S             |           |                  |  |
| Filename               | P140228B_11               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.2 g                    | Matrix    | Soil             |  |
| % Moisture             | 7.7                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.41 g                    | Collected | 02/04/2014 14:55 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 00:47 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 130.0      | ----       | 0.29     | 2,3,7,8-TCDF-13C         | 2.00       | 73               |
| Total TCDF          | 3400.0     | ----       | 0.29 E   | 2,3,7,8-TCDD-13C         | 2.00       | 81               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 71               |
| 2,3,7,8-TCDD        | 8.7        | ----       | 0.30     | 2,3,4,7,8-PeCDF-13C      | 2.00       | 66               |
| Total TCDD          | 340.0      | ----       | 0.30     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 70               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 97               |
| 1,2,3,7,8-PeCDF     | 130.0      | ----       | 0.41     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 92               |
| 2,3,4,7,8-PeCDF     | 290.0      | ----       | 0.52     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 84               |
| Total PeCDF         | 3200.0     | ----       | 0.46     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 57               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 81               |
| 1,2,3,7,8-PeCDD     | 36.0       | ----       | 0.52     | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 71               |
| Total PeCDD         | 620.0      | ----       | 0.52     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 67               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 65               |
| 1,2,3,4,7,8-HxCDF   | 290.0      | ----       | 0.94     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 70               |
| 1,2,3,6,7,8-HxCDF   | 210.0      | ----       | 0.78     | OCDD-13C                 | 4.00       | 48               |
| 2,3,4,6,7,8-HxCDF   | 390.0      | ----       | 1.30     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 56.0       | ----       | 1.20     | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 3200.0     | ----       | 1.10     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 47.0       | ----       | 0.80     | 2,3,7,8-TCDD-37Cl4       | 0.20       | 83               |
| 1,2,3,6,7,8-HxCDD   | 100.0      | ----       | 0.89     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 75.0       | ----       | 0.88     |                          |            |                  |
| Total HxCDD         | 1400.0     | ----       | 0.86     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 1400.0     | ----       | 0.69     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 61.0       | ----       | 0.71     | Equivalence: 290 ng/Kg   |            |                  |
| Total HpCDF         | 1900.0     | ----       | 0.70     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 920.0      | ----       | 1.10     |                          |            |                  |
| Total HpCDD         | 1900.0     | ----       | 1.10     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 490.0      | ----       | 0.63     |                          |            |                  |
| OCDD                | 4500.0     | ----       | 1.40     |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
E = Exceeds calibration range

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-9                  |           |                  |  |
| Lab Sample ID          | 10257346011-S             |           |                  |  |
| Filename               | P140228B_11               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.2 g                    | Matrix    | Soil             |  |
| % Moisture             | 7.7                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.41 g                    | Collected | 02/04/2014 14:55 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 00:47 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB      | MB      | UB      |
|---------------------|---------------|-------------|---------|---------|---------|---------|
| 2,3,7,8-TCDF        | 130.0         | 0.29        | 0.10000 | 12.7166 | 12.7166 | 12.7166 |
| Total TCDF          | 3400.0        | 0.29        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 2,3,7,8-TCDD        | 8.7           | 0.30        | 1.00000 | 8.6543  | 8.6543  | 8.6543  |
| Total TCDD          | 340.0         | 0.30        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDF     | 130.0         | 0.41        | 0.03000 | 4.0015  | 4.0015  | 4.0015  |
| 2,3,4,7,8-PeCDF     | 290.0         | 0.52        | 0.30000 | 87.8277 | 87.8277 | 87.8277 |
| Total PeCDF         | 3200.0        | 0.46        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDD     | 36.0          | 0.52        | 1.00000 | 36.0399 | 36.0399 | 36.0399 |
| Total PeCDD         | 620.0         | 0.52        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDF   | 290.0         | 0.94        | 0.10000 | 28.6286 | 28.6286 | 28.6286 |
| 1,2,3,6,7,8-HxCDF   | 210.0         | 0.78        | 0.10000 | 21.0516 | 21.0516 | 21.0516 |
| 2,3,4,6,7,8-HxCDF   | 390.0         | 1.3         | 0.10000 | 38.6524 | 38.6524 | 38.6524 |
| 1,2,3,7,8,9-HxCDF   | 56.0          | 1.2         | 0.10000 | 5.5926  | 5.5926  | 5.5926  |
| Total HxCDF         | 3200.0        | 1.1         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDD   | 47.0          | 0.80        | 0.10000 | 4.6681  | 4.6681  | 4.6681  |
| 1,2,3,6,7,8-HxCDD   | 100.0         | 0.89        | 0.10000 | 10.3913 | 10.3913 | 10.3913 |
| 1,2,3,7,8,9-HxCDD   | 75.0          | 0.88        | 0.10000 | 7.5209  | 7.5209  | 7.5209  |
| Total HxCDD         | 1400.0        | 0.86        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDF | 1400.0        | 0.69        | 0.01000 | 14.2588 | 14.2588 | 14.2588 |
| 1,2,3,4,7,8,9-HpCDF | 61.0          | 0.71        | 0.01000 | 0.6129  | 0.6129  | 0.6129  |
| Total HpCDF         | 1900.0        | 0.70        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDD | 920.0         | 1.1         | 0.01000 | 9.2390  | 9.2390  | 9.2390  |
| Total HpCDD         | 1900.0        | 1.1         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| OCDF                | 490.0         | 0.63        | 0.00030 | 0.1458  | 0.1458  | 0.1458  |
| OCDD                | 4500.0        | 1.4         | 0.00030 | 1.3427  | 1.3427  | 1.3427  |

**290 ng/Kg      290 ng/Kg      290 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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### Method 8290 Sample Analysis Results

Client - ESN Pacific

|                        |                           |           |                  |
|------------------------|---------------------------|-----------|------------------|
| Client's Sample ID     | RHS-DU-10                 |           |                  |
| Lab Sample ID          | 10257346012-S             |           |                  |
| Filename               | P140228B_12               |           |                  |
| Injected By            | BAL                       |           |                  |
| Total Amount Extracted | 10.8 g                    | Matrix    | Soil             |
| % Moisture             | 8.3                       | Dilution  | NA               |
| Dry Weight Extracted   | 9.90 g                    | Collected | 02/05/2014 09:15 |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 01:30 |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 54.0       | ----       | 0.21     | 2,3,7,8-TCDF-13C         | 2.00       | 68               |
| Total TCDF          | 1300.0     | ----       | 0.21     | 2,3,7,8-TCDD-13C         | 2.00       | 72               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 69               |
| 2,3,7,8-TCDD        | 3.0        | ----       | 0.21     | 2,3,4,7,8-PeCDF-13C      | 2.00       | 65               |
| Total TCDD          | 120.0      | ----       | 0.21     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 70               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 86               |
| 1,2,3,7,8-PeCDF     | 57.0       | ----       | 0.31     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 81               |
| 2,3,4,7,8-PeCDF     | 110.0      | ----       | 0.46     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 75               |
| Total PeCDF         | 1200.0     | ----       | 0.38     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 51               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 76               |
| 1,2,3,7,8-PeCDD     | 14.0       | ----       | 0.32     | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 64               |
| Total PeCDD         | 190.0      | ----       | 0.32     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 63               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 55               |
| 1,2,3,4,7,8-HxCDF   | 93.0       | ----       | 0.41     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 61               |
| 1,2,3,6,7,8-HxCDF   | 79.0       | ----       | 0.42     | OCDD-13C                 | 4.00       | 40               |
| 2,3,4,6,7,8-HxCDF   | 120.0      | ----       | 0.45     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 17.0       | ----       | 0.54     | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 1100.0     | ----       | 0.46     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 19.0       | ----       | 0.26     | 2,3,7,8-TCDD-37Cl4       | 0.20       | 76               |
| 1,2,3,6,7,8-HxCDD   | 49.0       | ----       | 0.36     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 37.0       | ----       | 0.61     |                          |            |                  |
| Total HxCDD         | 550.0      | ----       | 0.41     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 440.0      | ----       | 0.32     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 20.0       | ----       | 0.41     | Equivalence: 110 ng/Kg   |            |                  |
| Total HpCDF         | 720.0      | ----       | 0.37     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 920.0      | ----       | 0.97     |                          |            |                  |
| Total HpCDD         | 1900.0     | ----       | 0.97     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 430.0      | ----       | 0.46     |                          |            |                  |
| OCDD                | 5700.0     | ----       | 0.48     |                          |            |                  |

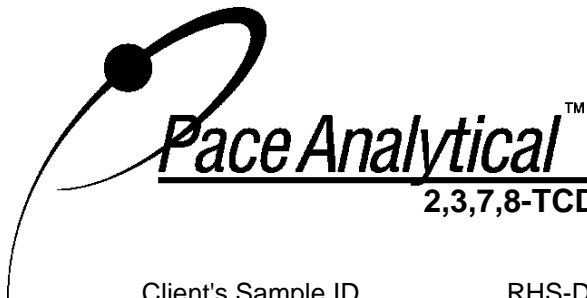
Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

## REPORT OF LABORATORY ANALYSIS

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-10                 |           |                  |  |
| Lab Sample ID          | 10257346012-S             |           |                  |  |
| Filename               | P140228B_12               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.8 g                    | Matrix    | Soil             |  |
| % Moisture             | 8.3                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.90 g                    | Collected | 02/05/2014 09:15 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 01:30 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB      | MB      | UB      |
|---------------------|---------------|-------------|---------|---------|---------|---------|
| 2,3,7,8-TCDF        | 54.0          | 0.21        | 0.10000 | 5.3519  | 5.3519  | 5.3519  |
| Total TCDF          | 1300.0        | 0.21        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 2,3,7,8-TCDD        | 3.0           | 0.21        | 1.00000 | 2.9992  | 2.9992  | 2.9992  |
| Total TCDD          | 120.0         | 0.21        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDF     | 57.0          | 0.31        | 0.03000 | 1.6989  | 1.6989  | 1.6989  |
| 2,3,4,7,8-PeCDF     | 110.0         | 0.46        | 0.30000 | 31.8489 | 31.8489 | 31.8489 |
| Total PeCDF         | 1200.0        | 0.38        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDD     | 14.0          | 0.32        | 1.00000 | 14.4344 | 14.4344 | 14.4344 |
| Total PeCDD         | 190.0         | 0.32        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDF   | 93.0          | 0.41        | 0.10000 | 9.2541  | 9.2541  | 9.2541  |
| 1,2,3,6,7,8-HxCDF   | 79.0          | 0.42        | 0.10000 | 7.8939  | 7.8939  | 7.8939  |
| 2,3,4,6,7,8-HxCDF   | 120.0         | 0.45        | 0.10000 | 11.5691 | 11.5691 | 11.5691 |
| 1,2,3,7,8,9-HxCDF   | 17.0          | 0.54        | 0.10000 | 1.6940  | 1.6940  | 1.6940  |
| Total HxCDF         | 1100.0        | 0.46        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDD   | 19.0          | 0.26        | 0.10000 | 1.8932  | 1.8932  | 1.8932  |
| 1,2,3,6,7,8-HxCDD   | 49.0          | 0.36        | 0.10000 | 4.8812  | 4.8812  | 4.8812  |
| 1,2,3,7,8,9-HxCDD   | 37.0          | 0.61        | 0.10000 | 3.6766  | 3.6766  | 3.6766  |
| Total HxCDD         | 550.0         | 0.41        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDF | 440.0         | 0.32        | 0.01000 | 4.4471  | 4.4471  | 4.4471  |
| 1,2,3,4,7,8,9-HpCDF | 20.0          | 0.41        | 0.01000 | 0.1967  | 0.1967  | 0.1967  |
| Total HpCDF         | 720.0         | 0.37        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDD | 920.0         | 0.97        | 0.01000 | 9.2161  | 9.2161  | 9.2161  |
| Total HpCDD         | 1900.0        | 0.97        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| OCDF                | 430.0         | 0.46        | 0.00030 | 0.1299  | 0.1299  | 0.1299  |
| OCDD                | 5700.0        | 0.48        | 0.00030 | 1.7155  | 1.7155  | 1.7155  |

**110 ng/Kg      110 ng/Kg      110 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-11                 |           |                  |  |
| Lab Sample ID          | 10257346013-S             |           |                  |  |
| Filename               | P140228B_13               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.6 g                    | Matrix    | Soil             |  |
| % Moisture             | 12.2                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.31 g                    | Collected | 02/05/2014 09:00 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 02:12 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 140        | ----       | 0.31     | 2,3,7,8-TCDF-13C         | 2.00       | 76               |
| Total TCDF          | 3200       | ----       | 0.31 E   | 2,3,7,8-TCDD-13C         | 2.00       | 84               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 72               |
| 2,3,7,8-TCDD        | -----      | 7.2        | 0.24 I   | 2,3,4,7,8-PeCDF-13C      | 2.00       | 70               |
| Total TCDD          | 250        | ----       | 0.24     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 74               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 101              |
| 1,2,3,7,8-PeCDF     | 120        | ----       | 0.64     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 94               |
| 2,3,4,7,8-PeCDF     | 240        | ----       | 0.46     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 85               |
| Total PeCDF         | 2800       | ----       | 0.55     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 59               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 88               |
| 1,2,3,7,8-PeCDD     | 26         | ----       | 0.36     | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 73               |
| Total PeCDD         | 390        | ----       | 0.36     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 69               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 62               |
| 1,2,3,4,7,8-HxCDF   | 200        | ----       | 0.63     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 68               |
| 1,2,3,6,7,8-HxCDF   | 160        | ----       | 0.77     | OCDD-13C                 | 4.00       | 43               |
| 2,3,4,6,7,8-HxCDF   | 250        | ----       | 0.83     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 34         | ----       | 0.77     | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 2200       | ----       | 0.75     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 28         | ----       | 0.53     | 2,3,7,8-TCDD-37Cl4       | 0.20       | 117              |
| 1,2,3,6,7,8-HxCDD   | 62         | ----       | 0.60     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 48         | ----       | 0.54     |                          |            |                  |
| Total HxCDD         | 810        | ----       | 0.56     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 870        | ----       | 0.55     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 37         | ----       | 0.96     | Equivalence: 220 ng/Kg   |            |                  |
| Total HpCDF         | 1300       | ----       | 0.75     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 920        | ----       | 1.90     |                          |            |                  |
| Total HpCDD         | 1900       | ----       | 1.90     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 630        | ----       | 1.00     |                          |            |                  |
| OCDD                | 6000       | ----       | 0.92     |                          |            |                  |

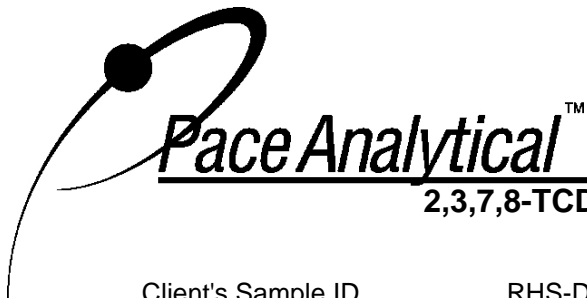
Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
E = Exceeds calibration range  
I = Interference present

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-11                 |           |                  |  |
| Lab Sample ID          | 10257346013-S             |           |                  |  |
| Filename               | P140228B_13               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.6 g                    | Matrix    | Soil             |  |
| % Moisture             | 12.2                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.31 g                    | Collected | 02/05/2014 09:00 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 02:12 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB      | MB      | UB      |
|---------------------|---------------|-------------|---------|---------|---------|---------|
| 2,3,7,8-TCDF        | 140           | 0.31        | 0.10000 | 13.8611 | 13.8611 | 13.8611 |
| Total TCDF          | 3200          | 0.31        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 2,3,7,8-TCDD        | ND            | 0.24        | 1.00000 | 7.2140  | 7.2140  | 7.2140  |
| Total TCDD          | 250           | 0.24        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDF     | 120           | 0.64        | 0.03000 | 3.6829  | 3.6829  | 3.6829  |
| 2,3,4,7,8-PeCDF     | 240           | 0.46        | 0.30000 | 72.1345 | 72.1345 | 72.1345 |
| Total PeCDF         | 2800          | 0.55        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDD     | 26            | 0.36        | 1.00000 | 26.2290 | 26.2290 | 26.2290 |
| Total PeCDD         | 390           | 0.36        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDF   | 200           | 0.63        | 0.10000 | 20.1051 | 20.1051 | 20.1051 |
| 1,2,3,6,7,8-HxCDF   | 160           | 0.77        | 0.10000 | 15.8041 | 15.8041 | 15.8041 |
| 2,3,4,6,7,8-HxCDF   | 250           | 0.83        | 0.10000 | 24.8891 | 24.8891 | 24.8891 |
| 1,2,3,7,8,9-HxCDF   | 34            | 0.77        | 0.10000 | 3.4094  | 3.4094  | 3.4094  |
| Total HxCDF         | 2200          | 0.75        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDD   | 28            | 0.53        | 0.10000 | 2.8169  | 2.8169  | 2.8169  |
| 1,2,3,6,7,8-HxCDD   | 62            | 0.60        | 0.10000 | 6.2413  | 6.2413  | 6.2413  |
| 1,2,3,7,8,9-HxCDD   | 48            | 0.54        | 0.10000 | 4.7967  | 4.7967  | 4.7967  |
| Total HxCDD         | 810           | 0.56        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDF | 870           | 0.55        | 0.01000 | 8.6815  | 8.6815  | 8.6815  |
| 1,2,3,4,7,8,9-HpCDF | 37            | 0.96        | 0.01000 | 0.3717  | 0.3717  | 0.3717  |
| Total HpCDF         | 1300          | 0.75        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDD | 920           | 1.9         | 0.01000 | 9.1633  | 9.1633  | 9.1633  |
| Total HpCDD         | 1900          | 1.9         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| OCDF                | 630           | 1.0         | 0.00030 | 0.1884  | 0.1884  | 0.1884  |
| OCDD                | 6000          | 0.92        | 0.00030 | 1.8081  | 1.8081  | 1.8081  |

**220 ng/Kg      220 ng/Kg      220 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-12                 |           |                  |  |
| Lab Sample ID          | 10257346014-S             |           |                  |  |
| Filename               | P140228B_14               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.2 g                    | Matrix    | Soil             |  |
| % Moisture             | 16.9                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 8.48 g                    | Collected | 02/05/2014 09:45 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 02:55 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 160        | ----       | 0.96     | 2,3,7,8-TCDF-13C         | 2.00       | 67               |
| Total TCDF          | 3700       | ----       | 0.96 E   | 2,3,7,8-TCDD-13C         | 2.00       | 73               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 64               |
| 2,3,7,8-TCDD        | -----      | 7.8        | 0.92 I   | 2,3,4,7,8-PeCDF-13C      | 2.00       | 61               |
| Total TCDD          | 210        | ----       | 0.92     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 63               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 85               |
| 1,2,3,7,8-PeCDF     | 140        | ----       | 2.20     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 83               |
| 2,3,4,7,8-PeCDF     | 250        | ----       | 2.00     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 77               |
| Total PeCDF         | 2800       | ----       | 2.10     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 70               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 77               |
| 1,2,3,7,8-PeCDD     | 26         | ----       | 1.30     | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 66               |
| Total PeCDD         | 340        | ----       | 1.30     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 63               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 59               |
| 1,2,3,4,7,8-HxCDF   | 200        | ----       | 1.40     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 65               |
| 1,2,3,6,7,8-HxCDF   | 200        | ----       | 1.70     | OCDD-13C                 | 4.00       | 42               |
| 2,3,4,6,7,8-HxCDF   | 250        | ----       | 1.70     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 41         | ----       | 1.90     | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 2200       | ----       | 1.70     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 28         | ----       | 2.10     | 2,3,7,8-TCDD-37Cl4       | 0.20       | 87               |
| 1,2,3,6,7,8-HxCDD   | 66         | ----       | 2.60     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 52         | ----       | 2.70     |                          |            |                  |
| Total HxCDD         | 740        | ----       | 2.50     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 810        | ----       | 1.10     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 45         | ----       | 1.20     | Equivalence: 230 ng/Kg   |            |                  |
| Total HpCDF         | 1200       | ----       | 1.20     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 1100       | ----       | 2.70     |                          |            |                  |
| Total HpCDD         | 2100       | ----       | 2.70     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 660        | ----       | 3.00     |                          |            |                  |
| OCDD                | 7000       | ----       | 2.80     |                          |            |                  |

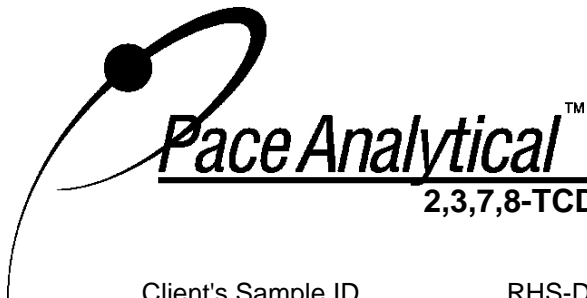
Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
E = Exceeds calibration range  
I = Interference present

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-12                 |           |                  |  |
| Lab Sample ID          | 10257346014-S             |           |                  |  |
| Filename               | P140228B_14               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.2 g                    | Matrix    | Soil             |  |
| % Moisture             | 16.9                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 8.48 g                    | Collected | 02/05/2014 09:45 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 02:55 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB      | MB      | UB      |
|---------------------|---------------|-------------|---------|---------|---------|---------|
| 2,3,7,8-TCDF        | 160           | 0.96        | 0.10000 | 16.4758 | 16.4758 | 16.4758 |
| Total TCDF          | 3700          | 0.96        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 2,3,7,8-TCDD        | ND            | 0.92        | 1.00000 | 7.7820  | 7.7820  | 7.7820  |
| Total TCDD          | 210           | 0.92        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDF     | 140           | 2.2         | 0.03000 | 4.1408  | 4.1408  | 4.1408  |
| 2,3,4,7,8-PeCDF     | 250           | 2.0         | 0.30000 | 74.6559 | 74.6559 | 74.6559 |
| Total PeCDF         | 2800          | 2.1         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDD     | 26            | 1.3         | 1.00000 | 25.9939 | 25.9939 | 25.9939 |
| Total PeCDD         | 340           | 1.3         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDF   | 200           | 1.4         | 0.10000 | 19.7951 | 19.7951 | 19.7951 |
| 1,2,3,6,7,8-HxCDF   | 200           | 1.7         | 0.10000 | 20.2713 | 20.2713 | 20.2713 |
| 2,3,4,6,7,8-HxCDF   | 250           | 1.7         | 0.10000 | 24.6069 | 24.6069 | 24.6069 |
| 1,2,3,7,8,9-HxCDF   | 41            | 1.9         | 0.10000 | 4.0567  | 4.0567  | 4.0567  |
| Total HxCDF         | 2200          | 1.7         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDD   | 28            | 2.1         | 0.10000 | 2.7781  | 2.7781  | 2.7781  |
| 1,2,3,6,7,8-HxCDD   | 66            | 2.6         | 0.10000 | 6.5790  | 6.5790  | 6.5790  |
| 1,2,3,7,8,9-HxCDD   | 52            | 2.7         | 0.10000 | 5.2103  | 5.2103  | 5.2103  |
| Total HxCDD         | 740           | 2.5         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDF | 810           | 1.1         | 0.01000 | 8.1337  | 8.1337  | 8.1337  |
| 1,2,3,4,7,8,9-HpCDF | 45            | 1.2         | 0.01000 | 0.4486  | 0.4486  | 0.4486  |
| Total HpCDF         | 1200          | 1.2         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDD | 1100          | 2.7         | 0.01000 | 10.7723 | 10.7723 | 10.7723 |
| Total HpCDD         | 2100          | 2.7         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| OCDF                | 660           | 3.0         | 0.00030 | 0.1974  | 0.1974  | 0.1974  |
| OCDD                | 7000          | 2.8         | 0.00030 | 2.0878  | 2.0878  | 2.0878  |

**230 ng/Kg      230 ng/Kg      230 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-13                 |           |                  |  |
| Lab Sample ID          | 10257346015-S             |           |                  |  |
| Filename               | P140228B_15               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.3 g                    | Matrix    | Soil             |  |
| % Moisture             | 9.4                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.33 g                    | Collected | 02/05/2014 09:45 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 03:38 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 330        | ----       | 0.71     | 2,3,7,8-TCDF-13C         | 2.00       | 78               |
| Total TCDF          | 7100       | ----       | 0.71 E   | 2,3,7,8-TCDD-13C         | 2.00       | 87               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 75               |
| 2,3,7,8-TCDD        | 17         | ----       | 0.96     | 2,3,4,7,8-PeCDF-13C      | 2.00       | 73               |
| Total TCDD          | 550        | ----       | 0.96     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 73               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 94               |
| 1,2,3,7,8-PeCDF     | 230        | ----       | 2.40     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 93               |
| 2,3,4,7,8-PeCDF     | 420        | ----       | 2.00     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 87               |
| Total PeCDF         | 4700       | ----       | 2.20     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 83               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 86               |
| 1,2,3,7,8-PeCDD     | 47         | ----       | 1.10     | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 74               |
| Total PeCDD         | 680        | ----       | 1.10     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 74               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 67               |
| 1,2,3,4,7,8-HxCDF   | 320        | ----       | 1.80     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 72               |
| 1,2,3,6,7,8-HxCDF   | 280        | ----       | 1.40     | OCDD-13C                 | 4.00       | 47               |
| 2,3,4,6,7,8-HxCDF   | 370        | ----       | 2.10     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 55         | ----       | 3.30     | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 3300       | ----       | 2.10     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 44         | ----       | 1.40     | 2,3,7,8-TCDD-37Cl4       | 0.20       | 106              |
| 1,2,3,6,7,8-HxCDD   | 83         | ----       | 1.90     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 74         | ----       | 1.50     |                          |            |                  |
| Total HxCDD         | 1000       | ----       | 1.60     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 1100       | ----       | 0.95     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 49         | ----       | 2.00     | Equivalence: 380 ng/Kg   |            |                  |
| Total HpCDF         | 1500       | ----       | 1.50     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 860        | ----       | 2.00     |                          |            |                  |
| Total HpCDD         | 1800       | ----       | 2.00     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 410        | ----       | 2.10     |                          |            |                  |
| OCDD                | 5000       | ----       | 2.90     |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
E = Exceeds calibration range

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## 2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations

ESN Pacific

|                        |                           |           |                  |  |  |
|------------------------|---------------------------|-----------|------------------|--|--|
| Client's Sample ID     | RHS-DU-13                 |           |                  |  |  |
| Lab Sample ID          | 10257346015-S             |           |                  |  |  |
| Filename               | P140228B_15               |           |                  |  |  |
| Injected By            | BAL                       |           |                  |  |  |
| Total Amount Extracted | 10.3 g                    | Matrix    | Soil             |  |  |
| % Moisture             | 9.4                       | Dilution  | NA               |  |  |
| Dry Weight Extracted   | 9.33 g                    | Collected | 02/05/2014 09:45 |  |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 03:38 |  |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB       | MB       | UB       |
|---------------------|---------------|-------------|---------|----------|----------|----------|
| 2,3,7,8-TCDF        | 330           | 0.71        | 0.10000 | 32.7150  | 32.7150  | 32.7150  |
| Total TCDF          | 7100          | 0.71        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 2,3,7,8-TCDD        | 17            | 0.96        | 1.00000 | 17.0473  | 17.0473  | 17.0473  |
| Total TCDD          | 550           | 0.96        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,7,8-PeCDF     | 230           | 2.4         | 0.03000 | 6.8601   | 6.8601   | 6.8601   |
| 2,3,4,7,8-PeCDF     | 420           | 2.0         | 0.30000 | 126.6528 | 126.6528 | 126.6528 |
| Total PeCDF         | 4700          | 2.2         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,7,8-PeCDD     | 47            | 1.1         | 1.00000 | 47.3344  | 47.3344  | 47.3344  |
| Total PeCDD         | 680           | 1.1         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,7,8-HxCDF   | 320           | 1.8         | 0.10000 | 31.9249  | 31.9249  | 31.9249  |
| 1,2,3,6,7,8-HxCDF   | 280           | 1.4         | 0.10000 | 28.1113  | 28.1113  | 28.1113  |
| 2,3,4,6,7,8-HxCDF   | 370           | 2.1         | 0.10000 | 37.3267  | 37.3267  | 37.3267  |
| 1,2,3,7,8,9-HxCDF   | 55            | 3.3         | 0.10000 | 5.5191   | 5.5191   | 5.5191   |
| Total HxCDF         | 3300          | 2.1         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,7,8-HxCDD   | 44            | 1.4         | 0.10000 | 4.3691   | 4.3691   | 4.3691   |
| 1,2,3,6,7,8-HxCDD   | 83            | 1.9         | 0.10000 | 8.3001   | 8.3001   | 8.3001   |
| 1,2,3,7,8,9-HxCDD   | 74            | 1.5         | 0.10000 | 7.3762   | 7.3762   | 7.3762   |
| Total HxCDD         | 1000          | 1.6         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,6,7,8-HpCDF | 1100          | 0.95        | 0.01000 | 11.2948  | 11.2948  | 11.2948  |
| 1,2,3,4,7,8,9-HpCDF | 49            | 2.0         | 0.01000 | 0.4879   | 0.4879   | 0.4879   |
| Total HpCDF         | 1500          | 1.5         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,6,7,8-HpCDD | 860           | 2.0         | 0.01000 | 8.5669   | 8.5669   | 8.5669   |
| Total HpCDD         | 1800          | 2.0         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| OCDF                | 410           | 2.1         | 0.00030 | 0.1232   | 0.1232   | 0.1232   |
| OCDD                | 5000          | 2.9         | 0.00030 | 1.4992   | 1.4992   | 1.4992   |

**380 ng/Kg      380 ng/Kg      380 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-14                 |           |                  |  |
| Lab Sample ID          | 10257346016-S             |           |                  |  |
| Filename               | P140228B_16               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.2 g                    | Matrix    | Soil             |  |
| % Moisture             | 16.8                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 8.49 g                    | Collected | 02/04/2014 12:52 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 04:21 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 29.0       | ----       | 0.46     | 2,3,7,8-TCDF-13C         | 2.00       | 75               |
| Total TCDF          | 690.0      | ----       | 0.46     | 2,3,7,8-TCDD-13C         | 2.00       | 81               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 70               |
| 2,3,7,8-TCDD        | -----      | 1.0        | 0.46     | 2,3,4,7,8-PeCDF-13C      | 2.00       | 66               |
| Total TCDD          | 41.0       | ----       | 0.46     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 68               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 91               |
| 1,2,3,7,8-PeCDF     | 26.0       | ----       | 0.60     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 88               |
| 2,3,4,7,8-PeCDF     | 50.0       | ----       | 0.58     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 82               |
| Total PeCDF         | 600.0      | ----       | 0.59     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 82               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 84               |
| 1,2,3,7,8-PeCDD     | 6.7        | ----       | 0.60     | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 72               |
| Total PeCDD         | 110.0      | ----       | 0.60     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 72               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 69               |
| 1,2,3,4,7,8-HxCDF   | 46.0       | ----       | 0.73     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 74               |
| 1,2,3,6,7,8-HxCDF   | 50.0       | ----       | 0.90     | OCDD-13C                 | 4.00       | 51               |
| 2,3,4,6,7,8-HxCDF   | 61.0       | ----       | 0.72     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 9.5        | ----       | 0.70     | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 610.0      | ----       | 0.76     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 12.0       | ----       | 1.20     | 2,3,7,8-TCDD-37Cl4       | 0.20       | 83               |
| 1,2,3,6,7,8-HxCDD   | 32.0       | ----       | 1.10     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 24.0       | ----       | 0.98     |                          |            |                  |
| Total HxCDD         | 320.0      | ----       | 1.10     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 280.0      | ----       | 1.20     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 13.0       | ----       | 1.20     | Equivalence: 61 ng/Kg    |            |                  |
| Total HpCDF         | 540.0      | ----       | 1.20     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 680.0      | ----       | 2.50     |                          |            |                  |
| Total HpCDD         | 1400.0     | ----       | 2.50     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 340.0      | ----       | 1.20     |                          |            |                  |
| OCDD                | 5300.0     | ----       | 1.80     |                          |            |                  |

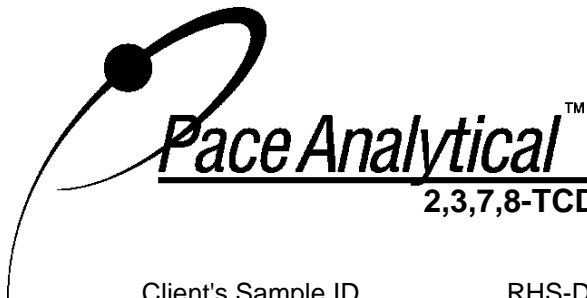
Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
J = Estimated value  
I = Interference present

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-14                 |           |                  |  |
| Lab Sample ID          | 10257346016-S             |           |                  |  |
| Filename               | P140228B_16               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.2 g                    | Matrix    | Soil             |  |
| % Moisture             | 16.8                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 8.49 g                    | Collected | 02/04/2014 12:52 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 04:21 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB      | MB      | UB      |
|---------------------|---------------|-------------|---------|---------|---------|---------|
| 2,3,7,8-TCDF        | 29.0          | 0.46        | 0.10000 | 2.8982  | 2.8982  | 2.8982  |
| Total TCDF          | 690.0         | 0.46        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 2,3,7,8-TCDD        | ND            | 0.46        | 1.00000 | 1.0046  | 1.0046  | 1.0046  |
| Total TCDD          | 41.0          | 0.46        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDF     | 26.0          | 0.60        | 0.03000 | 0.7718  | 0.7718  | 0.7718  |
| 2,3,4,7,8-PeCDF     | 50.0          | 0.58        | 0.30000 | 14.9708 | 14.9708 | 14.9708 |
| Total PeCDF         | 600.0         | 0.59        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDD     | 6.7           | 0.60        | 1.00000 | 6.6760  | 6.6760  | 6.6760  |
| Total PeCDD         | 110.0         | 0.60        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDF   | 46.0          | 0.73        | 0.10000 | 4.5985  | 4.5985  | 4.5985  |
| 1,2,3,6,7,8-HxCDF   | 50.0          | 0.90        | 0.10000 | 5.0122  | 5.0122  | 5.0122  |
| 2,3,4,6,7,8-HxCDF   | 61.0          | 0.72        | 0.10000 | 6.0587  | 6.0587  | 6.0587  |
| 1,2,3,7,8,9-HxCDF   | 9.5           | 0.70        | 0.10000 | 0.9543  | 0.9543  | 0.9543  |
| Total HxCDF         | 610.0         | 0.76        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDD   | 12.0          | 1.2         | 0.10000 | 1.1630  | 1.1630  | 1.1630  |
| 1,2,3,6,7,8-HxCDD   | 32.0          | 1.1         | 0.10000 | 3.2300  | 3.2300  | 3.2300  |
| 1,2,3,7,8,9-HxCDD   | 24.0          | 0.98        | 0.10000 | 2.3563  | 2.3563  | 2.3563  |
| Total HxCDD         | 320.0         | 1.1         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDF | 280.0         | 1.2         | 0.01000 | 2.8133  | 2.8133  | 2.8133  |
| 1,2,3,4,7,8,9-HpCDF | 13.0          | 1.2         | 0.01000 | 0.1286  | 0.1286  | 0.1286  |
| Total HpCDF         | 540.0         | 1.2         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDD | 680.0         | 2.5         | 0.01000 | 6.8214  | 6.8214  | 6.8214  |
| Total HpCDD         | 1400.0        | 2.5         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| OCDF                | 340.0         | 1.2         | 0.00030 | 0.1033  | 0.1033  | 0.1033  |
| OCDD                | 5300.0        | 1.8         | 0.00030 | 1.5833  | 1.5833  | 1.5833  |

**61 ng/Kg      61 ng/Kg      61 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-15                 |           |                  |  |
| Lab Sample ID          | 10257346017-S             |           |                  |  |
| Filename               | P140228B_17               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.4 g                    | Matrix    | Soil             |  |
| % Moisture             | 11.3                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.22 g                    | Collected | 02/04/2014 13:45 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 05:04 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 46         | ----       | 0.43     | 2,3,7,8-TCDF-13C         | 2.00       | 70               |
| Total TCDF          | 1100       | ----       | 0.43     | 2,3,7,8-TCDD-13C         | 2.00       | 77               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 66               |
| 2,3,7,8-TCDD        | -----      | 2.4        | 0.36 I   | 2,3,4,7,8-PeCDF-13C      | 2.00       | 62               |
| Total TCDD          | 87         | ----       | 0.36     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 65               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 85               |
| 1,2,3,7,8-PeCDF     | 46         | ----       | 0.62     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 84               |
| 2,3,4,7,8-PeCDF     | 97         | ----       | 0.65     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 79               |
| Total PeCDF         | 1100       | ----       | 0.64     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 74               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 76               |
| 1,2,3,7,8-PeCDD     | 14         | ----       | 0.69     | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 70               |
| Total PeCDD         | 170        | ----       | 0.69     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 68               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 65               |
| 1,2,3,4,7,8-HxCDF   | 89         | ----       | 1.70     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 70               |
| 1,2,3,6,7,8-HxCDF   | 82         | ----       | 0.89     | OCDD-13C                 | 4.00       | 50               |
| 2,3,4,6,7,8-HxCDF   | 110        | ----       | 1.20     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 17         | ----       | 0.70     | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 1000       | ----       | 1.10     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 18         | ----       | 1.60     | 2,3,7,8-TCDD-37Cl4       | 0.20       | 84               |
| 1,2,3,6,7,8-HxCDD   | 43         | ----       | 1.50     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 35         | ----       | 1.20     |                          |            |                  |
| Total HxCDD         | 490        | ----       | 1.50     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 490        | ----       | 0.53     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 24         | ----       | 0.88     | Equivalence: 110 ng/Kg   |            |                  |
| Total HpCDF         | 760        | ----       | 0.71     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 840        | ----       | 0.99     |                          |            |                  |
| Total HpCDD         | 1800       | ----       | 0.99     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 660        | ----       | 1.40     |                          |            |                  |
| OCDD                | 5700       | ----       | 2.00     |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
I = Interference present

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-15                 |           |                  |  |
| Lab Sample ID          | 10257346017-S             |           |                  |  |
| Filename               | P140228B_17               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.4 g                    | Matrix    | Soil             |  |
| % Moisture             | 11.3                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.22 g                    | Collected | 02/04/2014 13:45 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_01 & P140228B_19 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 05:04 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB      | MB      | UB      |
|---------------------|---------------|-------------|---------|---------|---------|---------|
| 2,3,7,8-TCDF        | 46            | 0.43        | 0.10000 | 4.6215  | 4.6215  | 4.6215  |
| Total TCDF          | 1100          | 0.43        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 2,3,7,8-TCDD        | ND            | 0.36        | 1.00000 | 2.3909  | 2.3909  | 2.3909  |
| Total TCDD          | 87            | 0.36        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDF     | 46            | 0.62        | 0.03000 | 1.3870  | 1.3870  | 1.3870  |
| 2,3,4,7,8-PeCDF     | 97            | 0.65        | 0.30000 | 29.1231 | 29.1231 | 29.1231 |
| Total PeCDF         | 1100          | 0.64        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDD     | 14            | 0.69        | 1.00000 | 14.2288 | 14.2288 | 14.2288 |
| Total PeCDD         | 170           | 0.69        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDF   | 89            | 1.7         | 0.10000 | 8.9167  | 8.9167  | 8.9167  |
| 1,2,3,6,7,8-HxCDF   | 82            | 0.89        | 0.10000 | 8.2004  | 8.2004  | 8.2004  |
| 2,3,4,6,7,8-HxCDF   | 110           | 1.2         | 0.10000 | 11.0519 | 11.0519 | 11.0519 |
| 1,2,3,7,8,9-HxCDF   | 17            | 0.70        | 0.10000 | 1.7175  | 1.7175  | 1.7175  |
| Total HxCDF         | 1000          | 1.1         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDD   | 18            | 1.6         | 0.10000 | 1.8308  | 1.8308  | 1.8308  |
| 1,2,3,6,7,8-HxCDD   | 43            | 1.5         | 0.10000 | 4.3374  | 4.3374  | 4.3374  |
| 1,2,3,7,8,9-HxCDD   | 35            | 1.2         | 0.10000 | 3.5304  | 3.5304  | 3.5304  |
| Total HxCDD         | 490           | 1.5         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDF | 490           | 0.53        | 0.01000 | 4.8664  | 4.8664  | 4.8664  |
| 1,2,3,4,7,8,9-HpCDF | 24            | 0.88        | 0.01000 | 0.2371  | 0.2371  | 0.2371  |
| Total HpCDF         | 760           | 0.71        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDD | 840           | 0.99        | 0.01000 | 8.4395  | 8.4395  | 8.4395  |
| Total HpCDD         | 1800          | 0.99        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| OCDF                | 660           | 1.4         | 0.00030 | 0.1972  | 0.1972  | 0.1972  |
| OCDD                | 5700          | 2.0         | 0.00030 | 1.7233  | 1.7233  | 1.7233  |

**110 ng/Kg      110 ng/Kg      110 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-16                 |           |                  |  |
| Lab Sample ID          | 10257346018-S             |           |                  |  |
| Filename               | P140301A_14               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.6 g                    | Matrix    | Soil             |  |
| % Moisture             | 9.1                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.64 g                    | Collected | 02/04/2014 14:11 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 16:29 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg |    | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|----|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 13.0       | ----       | 0.130    |    | 2,3,7,8-TCDF-13C         | 2.00       | 71               |
| Total TCDF          | 240.0      | ----       | 0.130    |    | 2,3,7,8-TCDD-13C         | 2.00       | 77               |
|                     |            |            |          |    | 1,2,3,7,8-PeCDF-13C      | 2.00       | 68               |
| 2,3,7,8-TCDD        | -----      | 0.79       | 0.110    | IJ | 2,3,4,7,8-PeCDF-13C      | 2.00       | 65               |
| Total TCDD          | 19.0       | ----       | 0.110    |    | 1,2,3,7,8-PeCDD-13C      | 2.00       | 67               |
|                     |            |            |          |    | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 78               |
| 1,2,3,7,8-PeCDF     | 13.0       | ----       | 0.130    |    | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 79               |
| 2,3,4,7,8-PeCDF     | 27.0       | ----       | 0.100    |    | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 74               |
| Total PeCDF         | 290.0      | ----       | 0.120    |    | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 54               |
|                     |            |            |          |    | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 73               |
| 1,2,3,7,8-PeCDD     | 6.1        | ----       | 0.130    |    | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 64               |
| Total PeCDD         | 51.0       | ----       | 0.130    |    | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 66               |
|                     |            |            |          |    | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 62               |
| 1,2,3,4,7,8-HxCDF   | 23.0       | ----       | 0.220    |    | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 67               |
| 1,2,3,6,7,8-HxCDF   | 16.0       | ----       | 0.370    |    | OCDD-13C                 | 4.00       | 68 Y             |
| 2,3,4,6,7,8-HxCDF   | 31.0       | ----       | 0.190    |    |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 4.2        | ----       | 0.230    | J  | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 340.0      | ----       | 0.250    |    | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 14.0       | ----       | 0.500    |    | 2,3,7,8-TCDD-37Cl4       | 0.20       | 80               |
| 1,2,3,6,7,8-HxCDD   | 36.0       | ----       | 0.460    |    |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 27.0       | ----       | 0.360    |    |                          |            |                  |
| Total HxCDD         | 370.0      | ----       | 0.440    |    |                          |            |                  |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 180.0      | ----       | 0.240    |    | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 7.0        | ----       | 0.270    |    | Equivalence: 44 ng/Kg    |            |                  |
| Total HpCDF         | 370.0      | ----       | 0.250    |    | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 850.0      | ----       | 0.280    |    |                          |            |                  |
| Total HpCDD         | 1900.0     | ----       | 0.280    |    |                          |            |                  |
|                     |            |            |          |    |                          |            |                  |
| OCDF                | 230.0      | ----       | 0.200    |    |                          |            |                  |
| OCDD                | 5800.0     | ----       | 0.310    |    |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

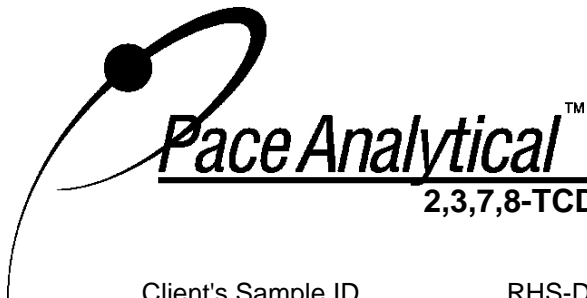
J = Estimated value

I = Interference present

Y = Calculated using average of daily RFs

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-16                 |           |                  |  |
| Lab Sample ID          | 10257346018-S             |           |                  |  |
| Filename               | P140301A_14               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.6 g                    | Matrix    | Soil             |  |
| % Moisture             | 9.1                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.64 g                    | Collected | 02/04/2014 14:11 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 16:29 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | 13.0          | 0.13        | 0.10000 | 1.2999 | 1.2999 | 1.2999 |
| Total TCDF          | 240.0         | 0.13        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | ND            | 0.11        | 1.00000 | 0.7901 | 0.7901 | 0.7901 |
| Total TCDD          | 19.0          | 0.11        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | 13.0          | 0.13        | 0.03000 | 0.3792 | 0.3792 | 0.3792 |
| 2,3,4,7,8-PeCDF     | 27.0          | 0.100       | 0.30000 | 7.9660 | 7.9660 | 7.9660 |
| Total PeCDF         | 290.0         | 0.12        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | 6.1           | 0.13        | 1.00000 | 6.1323 | 6.1323 | 6.1323 |
| Total PeCDD         | 51.0          | 0.13        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | 23.0          | 0.22        | 0.10000 | 2.2521 | 2.2521 | 2.2521 |
| 1,2,3,6,7,8-HxCDF   | 16.0          | 0.37        | 0.10000 | 1.5828 | 1.5828 | 1.5828 |
| 2,3,4,6,7,8-HxCDF   | 31.0          | 0.19        | 0.10000 | 3.0979 | 3.0979 | 3.0979 |
| 1,2,3,7,8,9-HxCDF   | 4.2           | 0.23        | 0.10000 | 0.4189 | 0.4189 | 0.4189 |
| Total HxCDF         | 340.0         | 0.25        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | 14.0          | 0.50        | 0.10000 | 1.3776 | 1.3776 | 1.3776 |
| 1,2,3,6,7,8-HxCDD   | 36.0          | 0.46        | 0.10000 | 3.5695 | 3.5695 | 3.5695 |
| 1,2,3,7,8,9-HxCDD   | 27.0          | 0.36        | 0.10000 | 2.7198 | 2.7198 | 2.7198 |
| Total HxCDD         | 370.0         | 0.44        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | 180.0         | 0.24        | 0.01000 | 1.8242 | 1.8242 | 1.8242 |
| 1,2,3,4,7,8,9-HpCDF | 7.0           | 0.27        | 0.01000 | 0.0699 | 0.0699 | 0.0699 |
| Total HpCDF         | 370.0         | 0.25        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | 850.0         | 0.28        | 0.01000 | 8.5450 | 8.5450 | 8.5450 |
| Total HpCDD         | 1900.0        | 0.28        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | 230.0         | 0.20        | 0.00030 | 0.0682 | 0.0682 | 0.0682 |
| OCDD                | 5800.0        | 0.31        | 0.00030 | 1.7485 | 1.7485 | 1.7485 |

**44 ng/Kg      44 ng/Kg      44 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Blank Analysis Results**

|                        |                           |             |                  |
|------------------------|---------------------------|-------------|------------------|
| Lab Sample ID          | BLANK-39446               | Matrix      | Solid            |
| Filename               | F140226B_09               | Dilution    | NA               |
| Total Amount Extracted | 10.5 g                    | Extracted   | 02/24/2014 21:00 |
| ICAL ID                | F131125                   | Analyzed    | 02/27/2014 02:08 |
| CCal Filename(s)       | F140226B_01 & F140226B_17 | Injected By | BAL              |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg |    | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|----|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | -----      | 0.140      | 0.035    | IJ | 2,3,7,8-TCDF-13C         | 2.00       | 74               |
| Total TCDF          | 0.430      | -----      | 0.035    | J  | 2,3,7,8-TCDD-13C         | 2.00       | 81               |
|                     |            |            |          |    | 1,2,3,7,8-PeCDF-13C      | 2.00       | 81               |
| 2,3,7,8-TCDD        | -----      | 0.060      | 0.057    | IJ | 2,3,4,7,8-PeCDF-13C      | 2.00       | 82               |
| Total TCDD          | ND         | -----      | 0.057    |    | 1,2,3,7,8-PeCDD-13C      | 2.00       | 88               |
|                     |            |            |          |    | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 78               |
| 1,2,3,7,8-PeCDF     | -----      | 0.150      | 0.062    | IJ | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 91               |
| 2,3,4,7,8-PeCDF     | ND         | -----      | 0.037    |    | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 88               |
| Total PeCDF         | ND         | -----      | 0.049    |    | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 82               |
|                     |            |            |          |    | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 73               |
| 1,2,3,7,8-PeCDD     | -----      | 0.086      | 0.039    | IJ | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 77               |
| Total PeCDD         | ND         | -----      | 0.039    |    | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 74               |
|                     |            |            |          |    | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 73               |
| 1,2,3,4,7,8-HxCDF   | ND         | -----      | 0.048    |    | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 79               |
| 1,2,3,6,7,8-HxCDF   | 0.094      | -----      | 0.035    | J  | OCDD-13C                 | 4.00       | 70               |
| 2,3,4,6,7,8-HxCDF   | -----      | 0.028      | 0.028    | IJ |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | ND         | -----      | 0.042    |    | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 0.094      | -----      | 0.038    | J  | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | ND         | -----      | 0.050    |    | 2,3,7,8-TCDD-37Cl4       | 0.20       | 87               |
| 1,2,3,6,7,8-HxCDD   | -----      | 0.081      | 0.052    | IJ |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | ND         | -----      | 0.048    |    |                          |            |                  |
| Total HxCDD         | ND         | -----      | 0.050    |    |                          |            |                  |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | -----      | 0.096      | 0.056    | IJ | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | ND         | -----      | 0.047    |    | Equivalence: 0.19 ng/Kg  |            |                  |
| Total HpCDF         | ND         | -----      | 0.051    |    | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | -----      | 0.068      | 0.054    | IJ |                          |            |                  |
| Total HpCDD         | 0.100      | -----      | 0.054    | J  |                          |            |                  |
|                     |            |            |          |    |                          |            |                  |
| OCDF                | 0.200      | -----      | 0.120    | J  |                          |            |                  |
| OCDD                | -----      | 0.370      | 0.150    | IJ |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

J = Estimated value

I = Interference present

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |  |
|------------------------|---------------------------|-----------|------------------|--|--|
| Client's Sample ID     |                           |           |                  |  |  |
| Lab Sample ID          | BLANK-39446               |           |                  |  |  |
| Filename               | F140226B_09               |           |                  |  |  |
| Injected By            | BAL                       |           |                  |  |  |
| Total Amount Extracted | 10.5 g                    | Matrix    | Solid            |  |  |
| % Moisture             | 0.0                       | Dilution  | NA               |  |  |
| Dry Weight Extracted   | 10.5 g                    | Collected | 02/24/2014 14:08 |  |  |
| ICAL ID                | F131125                   | Received  | 02/24/2014 14:08 |  |  |
| CCal Filename(s)       | F140226B_01 & F140226B_17 | Extracted | 02/24/2014 21:00 |  |  |
| Method Blank ID        |                           | Analyzed  | 02/27/2014 02:08 |  |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | ND            | 0.035       | 0.10000 | 0.0136 | 0.0136 | 0.0136 |
| Total TCDF          | 0.430         | 0.035       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | ND            | 0.057       | 1.00000 | 0.0601 | 0.0601 | 0.0601 |
| Total TCDD          | ND            | 0.057       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | ND            | 0.062       | 0.03000 | 0.0044 | 0.0044 | 0.0044 |
| 2,3,4,7,8-PeCDF     | ND            | 0.037       | 0.30000 | 0.0000 | 0.0056 | 0.0112 |
| Total PeCDF         | ND            | 0.049       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | ND            | 0.039       | 1.00000 | 0.0857 | 0.0857 | 0.0857 |
| Total PeCDD         | ND            | 0.039       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | ND            | 0.048       | 0.10000 | 0.0000 | 0.0024 | 0.0048 |
| 1,2,3,6,7,8-HxCDF   | 0.094         | 0.035       | 0.10000 | 0.0094 | 0.0094 | 0.0094 |
| 2,3,4,6,7,8-HxCDF   | ND            | 0.028       | 0.10000 | 0.0028 | 0.0028 | 0.0028 |
| 1,2,3,7,8,9-HxCDF   | ND            | 0.042       | 0.10000 | 0.0000 | 0.0021 | 0.0042 |
| Total HxCDF         | 0.094         | 0.038       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | ND            | 0.050       | 0.10000 | 0.0000 | 0.0025 | 0.0050 |
| 1,2,3,6,7,8-HxCDD   | ND            | 0.052       | 0.10000 | 0.0081 | 0.0081 | 0.0081 |
| 1,2,3,7,8,9-HxCDD   | ND            | 0.048       | 0.10000 | 0.0000 | 0.0024 | 0.0048 |
| Total HxCDD         | ND            | 0.050       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | ND            | 0.056       | 0.01000 | 0.0010 | 0.0010 | 0.0010 |
| 1,2,3,4,7,8,9-HpCDF | ND            | 0.047       | 0.01000 | 0.0000 | 0.0002 | 0.0005 |
| Total HpCDF         | ND            | 0.051       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | ND            | 0.054       | 0.01000 | 0.0007 | 0.0007 | 0.0007 |
| Total HpCDD         | 0.100         | 0.054       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | 0.200         | 0.12        | 0.00030 | 0.0001 | 0.0001 | 0.0001 |
| OCDD                | ND            | 0.15        | 0.00030 | 0.0001 | 0.0001 | 0.0001 |

**0.19 ng/Kg      0.20 ng/Kg      0.22 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Blank Analysis Results**

|                        |                           |             |                  |
|------------------------|---------------------------|-------------|------------------|
| Lab Sample ID          | BLANK-39465               | Matrix      | Solid            |
| Filename               | P140301A_05               | Dilution    | NA               |
| Total Amount Extracted | 10.4 g                    | Extracted   | 02/26/2014 21:30 |
| ICAL ID                | P130624                   | Analyzed    | 03/01/2014 10:03 |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Injected By | BAL              |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | -----      | 0.14       | 0.140 IJ | 2,3,7,8-TCDF-13C         | 2.00       | 75               |
| Total TCDF          | ND         | -----      | 0.140    | 2,3,7,8-TCDD-13C         | 2.00       | 82               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 75               |
| 2,3,7,8-TCDD        | ND         | -----      | 0.150    | 2,3,4,7,8-PeCDF-13C      | 2.00       | 70               |
| Total TCDD          | ND         | -----      | 0.150    | 1,2,3,7,8-PeCDD-13C      | 2.00       | 73               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 83               |
| 1,2,3,7,8-PeCDF     | ND         | -----      | 0.160    | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 87               |
| 2,3,4,7,8-PeCDF     | ND         | -----      | 0.110    | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 83               |
| Total PeCDF         | ND         | -----      | 0.130    | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 83               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 79               |
| 1,2,3,7,8-PeCDD     | ND         | -----      | 0.160    | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 72               |
| Total PeCDD         | ND         | -----      | 0.160    | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 77               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 74               |
| 1,2,3,4,7,8-HxCDF   | ND         | -----      | 0.110    | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 76               |
| 1,2,3,6,7,8-HxCDF   | ND         | -----      | 0.084    | OCDD-13C                 | 4.00       | 81 Y             |
| 2,3,4,6,7,8-HxCDF   | ND         | -----      | 0.091    |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | ND         | -----      | 0.110    | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | ND         | -----      | 0.099    | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | ND         | -----      | 0.130    | 2,3,7,8-TCDD-37Cl4       | 0.20       | 90               |
| 1,2,3,6,7,8-HxCDD   | ND         | -----      | 0.130    |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | ND         | -----      | 0.130    |                          |            |                  |
| Total HxCDD         | ND         | -----      | 0.130    |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | -----      | 0.20       | 0.120 IJ | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | ND         | -----      | 0.170    | Equivalence: 0.018 ng/Kg |            |                  |
| Total HpCDF         | ND         | -----      | 0.140    | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 0.17       | -----      | 0.140 J  |                          |            |                  |
| Total HpCDD         | 0.44       | -----      | 0.140 J  |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | ND         | -----      | 0.260    |                          |            |                  |
| OCDD                | 0.95       | -----      | 0.330 J  |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

J = Estimated value

I = Interference present

Y = Calculated using average of daily RFs

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |  |
|------------------------|---------------------------|-----------|------------------|--|--|
| Client's Sample ID     |                           |           |                  |  |  |
| Lab Sample ID          | BLANK-39465               |           |                  |  |  |
| Filename               | P140301A_05               |           |                  |  |  |
| Injected By            | BAL                       |           |                  |  |  |
| Total Amount Extracted | 10.4 g                    | Matrix    | Solid            |  |  |
| % Moisture             | 0.0                       | Dilution  | NA               |  |  |
| Dry Weight Extracted   | 10.4 g                    | Collected | 02/25/2014 20:04 |  |  |
| ICAL ID                | P130624                   | Received  | 02/25/2014 20:04 |  |  |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Extracted | 02/26/2014 21:30 |  |  |
| Method Blank ID        |                           | Analyzed  | 03/01/2014 10:03 |  |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | ND            | 0.14        | 0.10000 | 0.0143 | 0.0143 | 0.0143 |
| Total TCDF          | ND            | 0.14        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | ND            | 0.15        | 1.00000 | 0.0000 | 0.0752 | 0.1504 |
| Total TCDD          | ND            | 0.15        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | ND            | 0.16        | 0.03000 | 0.0000 | 0.0023 | 0.0047 |
| 2,3,4,7,8-PeCDF     | ND            | 0.11        | 0.30000 | 0.0000 | 0.0169 | 0.0339 |
| Total PeCDF         | ND            | 0.13        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | ND            | 0.16        | 1.00000 | 0.0000 | 0.0775 | 0.1550 |
| Total PeCDD         | ND            | 0.16        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | ND            | 0.11        | 0.10000 | 0.0000 | 0.0053 | 0.0107 |
| 1,2,3,6,7,8-HxCDF   | ND            | 0.084       | 0.10000 | 0.0000 | 0.0042 | 0.0084 |
| 2,3,4,6,7,8-HxCDF   | ND            | 0.091       | 0.10000 | 0.0000 | 0.0045 | 0.0091 |
| 1,2,3,7,8,9-HxCDF   | ND            | 0.11        | 0.10000 | 0.0000 | 0.0057 | 0.0114 |
| Total HxCDF         | ND            | 0.099       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | ND            | 0.13        | 0.10000 | 0.0000 | 0.0067 | 0.0134 |
| 1,2,3,6,7,8-HxCDD   | ND            | 0.13        | 0.10000 | 0.0000 | 0.0064 | 0.0129 |
| 1,2,3,7,8,9-HxCDD   | ND            | 0.13        | 0.10000 | 0.0000 | 0.0065 | 0.0130 |
| Total HxCDD         | ND            | 0.13        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | ND            | 0.12        | 0.01000 | 0.0020 | 0.0020 | 0.0020 |
| 1,2,3,4,7,8,9-HpCDF | ND            | 0.17        | 0.01000 | 0.0000 | 0.0008 | 0.0017 |
| Total HpCDF         | ND            | 0.14        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | 0.17          | 0.14        | 0.01000 | 0.0017 | 0.0017 | 0.0017 |
| Total HpCDD         | 0.44          | 0.14        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | ND            | 0.26        | 0.00030 | 0.0000 | 0.0000 | 0.0001 |
| OCDD                | 0.95          | 0.33        | 0.00030 | 0.0003 | 0.0003 | 0.0003 |

**0.018 ng/Kg      0.23 ng/Kg      0.44 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Laboratory Control Spike Results**

|                        |                           |             |                  |
|------------------------|---------------------------|-------------|------------------|
| Lab Sample ID          | LCS-39447                 | Matrix      | Solid            |
| Filename               | F140226B_15               | Dilution    | NA               |
| Total Amount Extracted | 10.0 g                    | Extracted   | 02/24/2014 21:00 |
| ICAL ID                | F131125                   | Analyzed    | 02/27/2014 06:48 |
| CCal Filename(s)       | F140226B_01 & F140226B_17 | Injected By | BAL              |
| Method Blank ID        | BLANK-39446               |             |                  |

| Native Isomers      | Qs (ng) | Qm (ng) | % Rec. | Internal Standards      | ng's Added | Percent Recovery |
|---------------------|---------|---------|--------|-------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 0.20    | 0.23    | 115    | 2,3,7,8-TCDF-13C        | 2.0        | 75               |
| Total TCDF          |         |         |        | 2,3,7,8-TCDD-13C        | 2.0        | 85               |
|                     |         |         |        | 1,2,3,7,8-PeCDF-13C     | 2.0        | 85               |
| 2,3,7,8-TCDD        | 0.20    | 0.18    | 89     | 2,3,4,7,8-PeCDF-13C     | 2.0        | 85               |
| Total TCDD          |         |         |        | 1,2,3,7,8-PeCDD-13C     | 2.0        | 97               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDF-13C   | 2.0        | 81               |
| 1,2,3,7,8-PeCDF     | 1.0     | 1.2     | 117    | 1,2,3,6,7,8-HxCDF-13C   | 2.0        | 89               |
| 2,3,4,7,8-PeCDF     | 1.0     | 1.1     | 109    | 2,3,4,6,7,8-HxCDF-13C   | 2.0        | 87               |
| Total PeCDF         |         |         |        | 1,2,3,7,8,9-HxCDF-13C   | 2.0        | 82               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDD-13C   | 2.0        | 77               |
| 1,2,3,7,8-PeCDD     | 1.0     | 1.0     | 100    | 1,2,3,6,7,8-HxCDD-13C   | 2.0        | 78               |
| Total PeCDD         |         |         |        | 1,2,3,4,6,7,8-HpCDF-13C | 2.0        | 78               |
|                     |         |         |        | 1,2,3,4,7,8,9-HpCDF-13C | 2.0        | 75               |
| 1,2,3,4,7,8-HxCDF   | 1.0     | 1.1     | 110    | 1,2,3,4,6,7,8-HpCDD-13C | 2.0        | 83               |
| 1,2,3,6,7,8-HxCDF   | 1.0     | 1.1     | 109    | OCDD-13C                | 4.0        | 69               |
| 2,3,4,6,7,8-HxCDF   | 1.0     | 1.0     | 100    |                         |            |                  |
| 1,2,3,7,8,9-HxCDF   | 1.0     | 1.1     | 107    | 1,2,3,4-TCDD-13C        | 2.0        | NA               |
| Total HxCDF         |         |         |        | 1,2,3,7,8,9-HxCDD-13C   | 2.0        | NA               |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,7,8-HxCDD   | 1.0     | 1.2     | 115    | 2,3,7,8-TCDD-37Cl4      | 0.20       | 87               |
| 1,2,3,6,7,8-HxCDD   | 1.0     | 1.3     | 128    |                         |            |                  |
| 1,2,3,7,8,9-HxCDD   | 1.0     | 1.2     | 119    |                         |            |                  |
| Total HxCDD         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 1.0     | 1.1     | 106    |                         |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 1.0     | 0.94    | 94     |                         |            |                  |
| Total HpCDF         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 1.0     | 1.0     | 102    |                         |            |                  |
| Total HpCDD         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| OCDF                | 2.0     | 2.0     | 99     |                         |            |                  |
| OCDD                | 2.0     | 2.2     | 110    |                         |            |                  |

Qs = Quantity Spiked  
Qm = Quantity Measured  
Rec. = Recovery (Expressed as Percent)  
R = Recovery outside of target range

Y = RF averaging used in calculations  
Nn = Value obtained from additional analysis  
NA = Not Applicable  
\* = See Discussion

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**Method 8290 Laboratory Control Spike Results**

|                        |                           |             |                  |
|------------------------|---------------------------|-------------|------------------|
| Lab Sample ID          | LCS-39466                 | Matrix      | Solid            |
| Filename               | P140301A_01               | Dilution    | NA               |
| Total Amount Extracted | 10.6 g                    | Extracted   | 02/26/2014 21:30 |
| ICAL ID                | P130624                   | Analyzed    | 03/01/2014 07:12 |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Injected By | BAL              |
| Method Blank ID        | BLANK-39465               |             |                  |

| Native Isomers      | Qs (ng) | Qm (ng) | % Rec. | Internal Standards      | ng's Added | Percent Recovery |
|---------------------|---------|---------|--------|-------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 0.20    | 0.22    | 112    | 2,3,7,8-TCDF-13C        | 2.0        | 79               |
| Total TCDF          |         |         |        | 2,3,7,8-TCDD-13C        | 2.0        | 85               |
|                     |         |         |        | 1,2,3,7,8-PeCDF-13C     | 2.0        | 75               |
| 2,3,7,8-TCDD        | 0.20    | 0.18    | 89     | 2,3,4,7,8-PeCDF-13C     | 2.0        | 71               |
| Total TCDD          |         |         |        | 1,2,3,7,8-PeCDD-13C     | 2.0        | 75               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDF-13C   | 2.0        | 88               |
| 1,2,3,7,8-PeCDF     | 1.0     | 1.2     | 117    | 1,2,3,6,7,8-HxCDF-13C   | 2.0        | 90               |
| 2,3,4,7,8-PeCDF     | 1.0     | 1.1     | 111    | 2,3,4,6,7,8-HxCDF-13C   | 2.0        | 86               |
| Total PeCDF         |         |         |        | 1,2,3,7,8,9-HxCDF-13C   | 2.0        | 88               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDD-13C   | 2.0        | 80               |
| 1,2,3,7,8-PeCDD     | 1.0     | 0.97    | 97     | 1,2,3,6,7,8-HxCDD-13C   | 2.0        | 75               |
| Total PeCDD         |         |         |        | 1,2,3,4,6,7,8-HpCDF-13C | 2.0        | 82               |
|                     |         |         |        | 1,2,3,4,7,8,9-HpCDF-13C | 2.0        | 77               |
| 1,2,3,4,7,8-HxCDF   | 1.0     | 1.1     | 111    | 1,2,3,4,6,7,8-HpCDD-13C | 2.0        | 80               |
| 1,2,3,6,7,8-HxCDF   | 1.0     | 1.1     | 112    | OCDD-13C                | 4.0        | 100 Y            |
| 2,3,4,6,7,8-HxCDF   | 1.0     | 1.1     | 112    |                         |            |                  |
| 1,2,3,7,8,9-HxCDF   | 1.0     | 1.1     | 106    | 1,2,3,4-TCDD-13C        | 2.0        | NA               |
| Total HxCDF         |         |         |        | 1,2,3,7,8,9-HxCDD-13C   | 2.0        | NA               |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,7,8-HxCDD   | 1.0     | 1.2     | 115    | 2,3,7,8-TCDD-37Cl4      | 0.20       | 85               |
| 1,2,3,6,7,8-HxCDD   | 1.0     | 1.2     | 122    |                         |            |                  |
| 1,2,3,7,8,9-HxCDD   | 1.0     | 1.2     | 118    |                         |            |                  |
| Total HxCDD         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 1.0     | 1.1     | 114    |                         |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 1.0     | 1.0     | 103    |                         |            |                  |
| Total HpCDF         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 1.0     | 1.1     | 106    |                         |            |                  |
| Total HpCDD         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| OCDF                | 2.0     | 2.2     | 108    |                         |            |                  |
| OCDD                | 2.0     | 2.2     | 110    |                         |            |                  |

Qs = Quantity Spiked  
Qm = Quantity Measured  
Rec. = Recovery (Expressed as Percent)  
R = Recovery outside of target range

Y = RF averaging used in calculations  
Nn = Value obtained from additional analysis  
NA = Not Applicable  
\* = See Discussion

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**Method 8290 Laboratory Control Spike Results**

|                        |                           |             |                  |
|------------------------|---------------------------|-------------|------------------|
| Lab Sample ID          | LCSD-39448                | Matrix      | Solid            |
| Filename               | F140226B_16               | Dilution    | NA               |
| Total Amount Extracted | 10.5 g                    | Extracted   | 02/24/2014 21:00 |
| ICAL ID                | F131125                   | Analyzed    | 02/27/2014 07:34 |
| CCal Filename(s)       | F140226B_01 & F140226B_17 | Injected By | BAL              |
| Method Blank ID        | BLANK-39446               |             |                  |

| Native Isomers      | Qs (ng) | Qm (ng) | % Rec. | Internal Standards      | ng's Added | Percent Recovery |
|---------------------|---------|---------|--------|-------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 0.20    | 0.24    | 119    | 2,3,7,8-TCDF-13C        | 2.0        | 64               |
| Total TCDF          |         |         |        | 2,3,7,8-TCDD-13C        | 2.0        | 74               |
|                     |         |         |        | 1,2,3,7,8-PeCDF-13C     | 2.0        | 78               |
| 2,3,7,8-TCDD        | 0.20    | 0.18    | 91     | 2,3,4,7,8-PeCDF-13C     | 2.0        | 81               |
| Total TCDD          |         |         |        | 1,2,3,7,8-PeCDD-13C     | 2.0        | 90               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDF-13C   | 2.0        | 78               |
| 1,2,3,7,8-PeCDF     | 1.0     | 1.2     | 119    | 1,2,3,6,7,8-HxCDF-13C   | 2.0        | 90               |
| 2,3,4,7,8-PeCDF     | 1.0     | 1.1     | 109    | 2,3,4,6,7,8-HxCDF-13C   | 2.0        | 87               |
| Total PeCDF         |         |         |        | 1,2,3,7,8,9-HxCDF-13C   | 2.0        | 78               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDD-13C   | 2.0        | 73               |
| 1,2,3,7,8-PeCDD     | 1.0     | 1.0     | 103    | 1,2,3,6,7,8-HxCDD-13C   | 2.0        | 81               |
| Total PeCDD         |         |         |        | 1,2,3,4,6,7,8-HpCDF-13C | 2.0        | 78               |
|                     |         |         |        | 1,2,3,4,7,8,9-HpCDF-13C | 2.0        | 74               |
| 1,2,3,4,7,8-HxCDF   | 1.0     | 1.2     | 115    | 1,2,3,4,6,7,8-HpCDD-13C | 2.0        | 84               |
| 1,2,3,6,7,8-HxCDF   | 1.0     | 1.1     | 107    | OCDD-13C                | 4.0        | 73               |
| 2,3,4,6,7,8-HxCDF   | 1.0     | 1.0     | 102    |                         |            |                  |
| 1,2,3,7,8,9-HxCDF   | 1.0     | 1.1     | 107    | 1,2,3,4-TCDD-13C        | 2.0        | NA               |
| Total HxCDF         |         |         |        | 1,2,3,7,8,9-HxCDD-13C   | 2.0        | NA               |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,7,8-HxCDD   | 1.0     | 1.2     | 124    | 2,3,7,8-TCDD-37Cl4      | 0.20       | 75               |
| 1,2,3,6,7,8-HxCDD   | 1.0     | 1.2     | 123    |                         |            |                  |
| 1,2,3,7,8,9-HxCDD   | 1.0     | 1.2     | 121    |                         |            |                  |
| Total HxCDD         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 1.0     | 1.1     | 108    |                         |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 1.0     | 0.94    | 94     |                         |            |                  |
| Total HpCDF         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 1.0     | 1.0     | 102    |                         |            |                  |
| Total HpCDD         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| OCDF                | 2.0     | 2.0     | 101    |                         |            |                  |
| OCDD                | 2.0     | 2.2     | 110    |                         |            |                  |

Qs = Quantity Spiked  
Qm = Quantity Measured  
Rec. = Recovery (Expressed as Percent)  
R = Recovery outside of target range

Y = RF averaging used in calculations  
Nn = Value obtained from additional analysis  
NA = Not Applicable  
\* = See Discussion

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**Method 8290**

**Spike Recovery Relative Percent Difference (RPD) Results**

Client ESN Pacific

Spike 1 ID LCS-39447  
Spike 1 Filename F140226B\_15

Spike 2 ID LCSD-39448  
Spike 2 Filename F140226B\_16

| Compound            | Spike 1<br>%REC | Spike 2<br>%REC | %RPD |
|---------------------|-----------------|-----------------|------|
| 2,3,7,8-TCDF        | 115             | 119             | 3.4  |
| 2,3,7,8-TCDD        | 89              | 91              | 2.2  |
| 1,2,3,7,8-PeCDF     | 117             | 119             | 1.7  |
| 2,3,4,7,8-PeCDF     | 109             | 109             | 0.0  |
| 1,2,3,7,8-PeCDD     | 100             | 103             | 3.0  |
| 1,2,3,4,7,8-HxCDF   | 110             | 115             | 4.4  |
| 1,2,3,6,7,8-HxCDF   | 109             | 107             | 1.9  |
| 2,3,4,6,7,8-HxCDF   | 100             | 102             | 2.0  |
| 1,2,3,7,8,9-HxCDF   | 107             | 107             | 0.0  |
| 1,2,3,4,7,8-HxCDD   | 115             | 124             | 7.5  |
| 1,2,3,6,7,8-HxCDD   | 128             | 123             | 4.0  |
| 1,2,3,7,8,9-HxCDD   | 119             | 121             | 1.7  |
| 1,2,3,4,6,7,8-HpCDF | 106             | 108             | 1.9  |
| 1,2,3,4,7,8,9-HpCDF | 94              | 94              | 0.0  |
| 1,2,3,4,6,7,8-HpCDD | 102             | 102             | 0.0  |
| OCDF                | 99              | 101             | 2.0  |
| OCDD                | 110             | 110             | 0.0  |

%REC = Percent Recovered

RPD = The difference between the two values divided by the mean value

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### Method 8290 Spiked Sample Report

Client - ESN Pacific

|                        |                           |             |                  |
|------------------------|---------------------------|-------------|------------------|
| Client's Sample ID     | RHS-DU-1.2-MS             | Matrix      | Soil             |
| Lab Sample ID          | 10257346002-S-MS          | Dilution    | NA               |
| Filename               | P140301A_02               | Extracted   | 02/26/2014 21:30 |
| Total Amount Extracted | 10.0 g                    | Analyzed    | 03/01/2014 07:55 |
| ICAL ID                | P130624                   | Injected By | BAL              |
| CCal Filename(s)       | P140228B_19 & P140301A_15 |             |                  |
| Method Blank ID        | BLANK-39465               |             |                  |

| Native Isomers      | Qs (ng) | Qm (ng) | % Rec. | Internal Standards      | ng's Added | Percent Recovery |
|---------------------|---------|---------|--------|-------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 0.20    | 5.59    | 2796 E | 2,3,7,8-TCDF-13C        | 2.00       | 77               |
|                     |         |         |        | 2,3,7,8-TCDD-13C        | 2.00       | 82               |
|                     |         |         |        | 1,2,3,7,8-PeCDF-13C     | 2.00       | 73               |
| 2,3,7,8-TCDD        | 0.20    | 0.54    | 268    | 2,3,4,7,8-PeCDF-13C     | 2.00       | 70               |
|                     |         |         |        | 1,2,3,7,8-PeCDD-13C     | 2.00       | 72               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDF-13C   | 2.00       | 92               |
| 1,2,3,7,8-PeCDF     | 1.00    | 4.91    | 491    | 1,2,3,6,7,8-HxCDF-13C   | 2.00       | 90               |
| 2,3,4,7,8-PeCDF     | 1.00    | 8.91    | 891    | 2,3,4,6,7,8-HxCDF-13C   | 2.00       | 86               |
|                     |         |         |        | 1,2,3,7,8,9-HxCDF-13C   | 2.00       | 85               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDD-13C   | 2.00       | 80               |
| 1,2,3,7,8-PeCDD     | 1.00    | 1.92    | 192    | 1,2,3,6,7,8-HxCDD-13C   | 2.00       | 75               |
|                     |         |         |        | 1,2,3,4,6,7,8-HpCDF-13C | 2.00       | 75               |
|                     |         |         |        | 1,2,3,4,7,8,9-HpCDF-13C | 2.00       | 70               |
| 1,2,3,4,7,8-HxCDF   | 1.00    | 6.83    | 683    | 1,2,3,4,6,7,8-HpCDD-13C | 2.00       | 74               |
| 1,2,3,6,7,8-HxCDF   | 1.00    | 6.06    | 606    | OCDD-13C                | 4.00       | 79 Y             |
| 2,3,4,6,7,8-HxCDF   | 1.00    | 8.25    | 825    |                         |            |                  |
| 1,2,3,7,8,9-HxCDF   | 1.00    | 1.89    | 189    | 1,2,3,4-TCDD-13C        | 2.00       | NA               |
|                     |         |         |        | 1,2,3,7,8,9-HxCDD-13C   | 2.00       | NA               |
| 1,2,3,4,7,8-HxCDD   | 1.00    | 1.88    | 188    | 2,3,7,8-TCDD-37Cl4      | 0.20       | 84               |
| 1,2,3,6,7,8-HxCDD   | 1.00    | 2.54    | 254    |                         |            |                  |
| 1,2,3,7,8,9-HxCDD   | 1.00    | 2.25    | 225    |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 1.00    | 24.43   | 2443 E |                         |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 1.00    | 1.78    | 178    |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 1.00    | 8.00    | 800    |                         |            |                  |
| OCDF                | 2.00    | 7.05    | 352    |                         |            |                  |
| OCDD                | 2.00    | 19.55   | 977    |                         |            |                  |

Qs = Quantity Spiked                      Qm = Quantity Measured                      Rec. = Recovery (Expressed as Percent)

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

E = Exceeds calibration range

Y = Calculated using average of daily RFs

### REPORT OF LABORATORY ANALYSIS

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### Method 8290 Spiked Sample Report

Client - ESN Pacific

|                        |                           |             |                  |
|------------------------|---------------------------|-------------|------------------|
| Client's Sample ID     | RHS-DU-1.2-MSD            | Matrix      | Soil             |
| Lab Sample ID          | 10257346002-S-MSD         | Dilution    | NA               |
| Filename               | P140301A_03               | Extracted   | 02/26/2014 21:30 |
| Total Amount Extracted | 10.1 g                    | Analyzed    | 03/01/2014 08:38 |
| ICAL ID                | P130624                   | Injected By | BAL              |
| CCal Filename(s)       | P140228B_19 & P140301A_15 |             |                  |
| Method Blank ID        | BLANK-39465               |             |                  |

| Native Isomers      | Qs (ng) | Qm (ng) | % Rec. | Internal Standards      | ng's Added | Percent Recovery |
|---------------------|---------|---------|--------|-------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 0.20    | 5.58    | 2791 E | 2,3,7,8-TCDF-13C        | 2.00       | 80               |
|                     |         |         |        | 2,3,7,8-TCDD-13C        | 2.00       | 87               |
|                     |         |         |        | 1,2,3,7,8-PeCDF-13C     | 2.00       | 79               |
| 2,3,7,8-TCDD        | 0.20    | 0.53    | 266    | 2,3,4,7,8-PeCDF-13C     | 2.00       | 77               |
|                     |         |         |        | 1,2,3,7,8-PeCDD-13C     | 2.00       | 81               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDF-13C   | 2.00       | 90               |
| 1,2,3,7,8-PeCDF     | 1.00    | 5.14    | 514    | 1,2,3,6,7,8-HxCDF-13C   | 2.00       | 90               |
| 2,3,4,7,8-PeCDF     | 1.00    | 8.80    | 880    | 2,3,4,6,7,8-HxCDF-13C   | 2.00       | 86               |
|                     |         |         |        | 1,2,3,7,8,9-HxCDF-13C   | 2.00       | 83               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDD-13C   | 2.00       | 82               |
| 1,2,3,7,8-PeCDD     | 1.00    | 1.85    | 185    | 1,2,3,6,7,8-HxCDD-13C   | 2.00       | 74               |
|                     |         |         |        | 1,2,3,4,6,7,8-HpCDF-13C | 2.00       | 77               |
|                     |         |         |        | 1,2,3,4,7,8,9-HpCDF-13C | 2.00       | 77               |
| 1,2,3,4,7,8-HxCDF   | 1.00    | 6.92    | 692    | 1,2,3,4,6,7,8-HpCDD-13C | 2.00       | 81               |
| 1,2,3,6,7,8-HxCDF   | 1.00    | 6.11    | 611    | OCDD-13C                | 4.00       | 88 Y             |
| 2,3,4,6,7,8-HxCDF   | 1.00    | 8.12    | 812    |                         |            |                  |
| 1,2,3,7,8,9-HxCDF   | 1.00    | 1.94    | 194    | 1,2,3,4-TCDD-13C        | 2.00       | NA               |
|                     |         |         |        | 1,2,3,7,8,9-HxCDD-13C   | 2.00       | NA               |
| 1,2,3,4,7,8-HxCDD   | 1.00    | 1.98    | 198    | 2,3,7,8-TCDD-37Cl4      | 0.20       | 87               |
| 1,2,3,6,7,8-HxCDD   | 1.00    | 2.58    | 258    |                         |            |                  |
| 1,2,3,7,8,9-HxCDD   | 1.00    | 2.40    | 240    |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 1.00    | 24.04   | 2404 E |                         |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 1.00    | 1.82    | 182    |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 1.00    | 7.85    | 785    |                         |            |                  |
| OCDF                | 2.00    | 7.18    | 359    |                         |            |                  |
| OCDD                | 2.00    | 19.02   | 951    |                         |            |                  |

Qs = Quantity Spiked                      Qm = Quantity Measured                      Rec. = Recovery (Expressed as Percent)

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

E = Exceeds calibration range

Y = Calculated using average of daily RFs

### REPORT OF LABORATORY ANALYSIS

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### Method 8290 Spike Sample Results

Client - ESN Pacific

|                  |                   |                 |             |                    |        |
|------------------|-------------------|-----------------|-------------|--------------------|--------|
| Client Sample ID | RHS-DU-1.2        | Sample Filename | P140301A_06 | <u>Dry Weights</u> |        |
| Lab Sample ID    | 10257346002-S     | MS Filename     | P140301A_02 | Sample Amount      | 9.33 g |
| MS ID            | 10257346002-S-MS  | MSD Filename    | P140301A_03 | MS Amount          | 9.2 g  |
| MSD ID           | 10257346002-S-MSD |                 |             | MSD Amount         | 9.2 g  |

| Analyte             | Sample Conc.<br>ng/Kg | MS/MSD Qs<br>(ng) | MS Qm<br>(ng) | MSD Qm<br>(ng) | RPD | Background Subtracted |            |       |
|---------------------|-----------------------|-------------------|---------------|----------------|-----|-----------------------|------------|-------|
|                     |                       |                   |               |                |     | MS % Rec.             | MSD % Rec. | RPD   |
| 2,3,7,8-TCDF        | 630.003               | 0.20              | 5.59          | 5.58           | 0.2 | 0                     | 0          | 0.0   |
| 2,3,7,8-TCDD        | 38.298                | 0.20              | 0.54          | 0.53           | 0.6 | 93                    | 89         | 3.6   |
| 1,2,3,7,8-PeCDF     | 471.535               | 1.00              | 4.91          | 5.14           | 4.7 | 59                    | 79         | 27.8  |
| 2,3,4,7,8-PeCDF     | 969.248               | 1.00              | 8.91          | 8.80           | 1.2 | 4                     | 0          | 200.0 |
| 1,2,3,7,8-PeCDD     | 117.881               | 1.00              | 1.92          | 1.85           | 3.8 | 84                    | 76         | 10.3  |
| 1,2,3,4,7,8-HxCDF   | 872.673               | 1.00              | 6.83          | 6.92           | 1.3 | 0                     | 0          | 0.0   |
| 1,2,3,6,7,8-HxCDF   | 740.624               | 1.00              | 6.06          | 6.11           | 0.9 | 0                     | 0          | 0.0   |
| 2,3,4,6,7,8-HxCDF   | 1019.369              | 1.00              | 8.25          | 8.12           | 1.6 | 0                     | 0          | 0.0   |
| 1,2,3,7,8,9-HxCDF   | 121.000               | 1.00              | 1.89          | 1.94           | 2.5 | 79                    | 82         | 4.5   |
| 1,2,3,4,7,8-HxCDD   | 106.051               | 1.00              | 1.88          | 1.98           | 5.1 | 91                    | 100        | 9.3   |
| 1,2,3,6,7,8-HxCDD   | 199.110               | 1.00              | 2.54          | 2.58           | 1.5 | 72                    | 74         | 2.6   |
| 1,2,3,7,8,9-HxCDD   | 154.994               | 1.00              | 2.25          | 2.40           | 6.3 | 83                    | 97         | 14.7  |
| 1,2,3,4,6,7,8-HpCDF | 3233.560              | 1.00              | 24.43         | 24.04          | 1.6 | 0                     | 0          | 0.0   |
| 1,2,3,4,7,8,9-HpCDF | 104.864               | 1.00              | 1.78          | 1.82           | 2.2 | 82                    | 85         | 3.6   |
| 1,2,3,4,6,7,8-HpCDD | 849.687               | 1.00              | 8.00          | 7.85           | 2.0 | 23                    | 0          | 200.0 |
| OCDF                | 562.760               | 2.00              | 7.05          | 7.18           | 1.9 | 95                    | 99         | 4.2   |
| OCDD                | 1906.520              | 2.00              | 19.55         | 19.02          | 2.7 | 105                   | 70         | 39.9  |

**Definitions**

|                                   |                                    |
|-----------------------------------|------------------------------------|
| MS = Matrix Spike                 | CDD = Chlorinated dibenzo-p-dioxin |
| MSD = Matrix Spike Duplicate      | CDF = Chlorinated dibenzo-p-furan  |
| Qm = Quantity Measured            | T = Tetra                          |
| Qs = Quantity Spiked              | Pe = Penta                         |
| % Rec. = Percent Recovery         | Hx = Hexa                          |
| RPD = Relative Percent Difference | Hp = Hepta                         |
| NA = Not Applicable               | O = Octa                           |
| NC = Not Calculated               |                                    |

**Report Prepared for:**

Karen Carvallo  
ESN Pacific  
2020-B Kahai Street  
Honolulu HI 96819

**REPORT OF  
LABORATORY  
ANALYSIS FOR  
PCDD/PCDF**

**Report Information:**

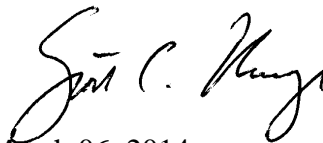
**Pace Project #: 10257348**  
**Sample Receipt Date: 02/11/2014**  
**Client Project #: D1402050037**  
**Client Sub PO #: N/A**  
**State Cert #: SLD**

**Invoicing & Reporting Options:**

The report provided has been invoiced as a Level 2 PCDD/PCDF Report. If an upgrade of this report package is requested, an additional charge may be applied.

Please review the attached invoice for accuracy and forward any questions to Scott Unze, your Pace Project Manager.

**This report has been reviewed by:**



March 06, 2014

Scott Unze, Project Manager  
(612) 607-6383  
(612) 607-6444 (fax)  
scott.unze@pacelabs.com

**Report Prepared Date:**

March 6, 2014



**Report of Laboratory Analysis**

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The results relate only to the samples included in this report.



## **DISCUSSION**

This report presents the results from the analyses performed on seventeen samples submitted by a representative of ESN Pacific. The samples were analyzed for the presence or absence of polychlorodibenzo-p-dioxins (PCDDs) and polychlorodibenzofurans (PCDFs) using a modified version of USEPA Method 8290. The reporting limits were based on signal-to-noise measurements.

The results provided in this report include lower-bound, mid-bound, and upper-bound toxic equivalence (TEQ) results. Lower bound TEQ results include only contributions from positive values in the sample. Mid-bound TEQ results include one-half of the reporting limit in cases where an analyte was not detected, along with contributions from positives in the sample. Upper-bound TEQ results include the full reporting limit in cases where an analyte was not detected, along with contributions from positives in the sample.

Second column confirmation analyses of 2,3,7,8-TCDF values obtained from the primary (DB5-MS) column are performed only when specifically requested for a project and only when the values are above the concentration of the lowest calibration standard. Typical resolution for this isomer using the DB5-MS column ranges from 25-30%.

The recoveries of the isotopically-labeled PCDD/PCDF internal standards in the sample extracts ranged from 47-134%. All of the labeled standard recoveries obtained for this project were within the 40-135% target range specified in Method 8290. Also, since the quantification of the native 2,3,7,8-substituted congeners was based on isotope dilution, the data were automatically corrected for variation in recovery and accurate values were obtained.

In some cases, interfering substances impacted the determinations of PCDD or PCDF congeners; the affected values were flagged "I" where incorrect isotope ratios were obtained or "P" where polychlorinated diphenyl ethers were present. Concentrations below the calibration range were flagged "J" and should be regarded as estimates. Concentrations above the calibration range were flagged "E" and should also be regarded as estimates.

A laboratory method blank was prepared and analyzed with each sample batch as part of our routine quality control procedures. The results show the blanks to contain trace levels of selected congeners. These levels were below the calibration range of the method. The levels reported for the affected congeners in the field samples were higher than the corresponding blank levels by one or more orders of magnitude. These results indicate that the sample processing steps did not contribute significantly to the levels reported for the field samples.

Laboratory and matrix spike samples were also prepared with the sample batches using clean sand or sample matrix that had been fortified with native standard materials. The results show that the spiked native compounds were generally recovered at 89-129% with relative percent differences of 0.0-8.0%. The background-subtracted recovery values obtained for TCDF and OCDD in the matrix spike duplicate were above the 70-130% target range, possibly due to sample inhomogeneity. Matrix spikes were prepared with the 02/26/2014 sample batch using sample material from a separate project; results from these analyses will be provided upon request. Matrix spikes were not prepared with the 02/24/2014 sample batch.

The responses obtained for selected native and labeled congeners in calibration standard analyses U140226B\_20, P140301A\_15, and U140304B\_20 were outside the target ranges. As specified in the method, the averages of the daily response factors for these compounds were used in the calculations for

## **REPORT OF LABORATORY ANALYSIS**

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**Pace Analytical Services, Inc.**  
1700 Elm Street  
Minneapolis, MN 55414  
Phone: 612.607.1700  
Fax: 612.607.6444

**DISCUSSION**

the samples from these runshifts. The affected values were flagged "Y" on the results tables.

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**REPORT OF LABORATORY ANALYSIS**

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## Minnesota Laboratory Certifications

| Authority       | Certificate # | Authority       | Certificate # |
|-----------------|---------------|-----------------|---------------|
| A2LA            | 2926.01       | Minnesota       | 027-053-137   |
| Alabama         | 40770         | Mississippi     | MN00064       |
| Alaska          | MN00064       | Montana         | 92            |
| Arizona         | AZ0014        | Nebraska        |               |
| Arkansas        | 88-0680       | Nevada          | MN_00064_200  |
| California      | 01155CA       | New Jersey (NE) | MN002         |
| Colorado        | MN00064       | New York (NEL)  | 11647         |
| Connecticut     | PH-0256       | North Carolina  | 27700         |
| EPA Region 8    | 8TMS-Q        | North Dakota    | R-036         |
| Florida (NELAP) | E87605        | Ohio            | 4150          |
| Georgia (DNR)   | 959           | Oklahoma        | D9922         |
| Guam            | 959           | Oregon (ELAP)   | MN200001-005  |
| Hawaii          | SLD           | Oregon (OREL)   | MN300001-001  |
| Idaho           | MN00064       | Pennsylvania    | 68-00563      |
| Illinois        | 200012        | Puerto Rico     | MN00064       |
| Indiana         | C-MN-01       | Saipan          | MP0003        |
| Indiana         | C-MN-01       | South Carolina  | 74003001      |
| Iowa            | 368           | Texas           | T104704192-08 |
| Kansas          | E-10167       | Utah (NELAP)    | MN00064       |
| Kentucky        | 90062         | Virginia        | 00251         |
| Louisiana       | 03086         | Washington      | C755          |
| Maine           | 2007029       | West Virginia   | 9952C         |
| Maryland        | 322           | Wisconsin       | 999407970     |
| Michigan        | 9909          | Wyoming         | 8TMS-Q        |

## REPORT OF LABORATORY ANALYSIS

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Report No.....10257348

## Appendix A

### Sample Management

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# ESN PACIFIC'S CHAIN-OF-CUSTODY RECORD

10257348

CLIENT: ESN Pacific  
 ADDRESS: 2020-B Kahai St Honolulu, HI 96819  
 PHONE: 8088470067 FAX: 8088470917  
 EMAIL: esn@esnpacific.com  
 CLIENT PROJECT #: \_\_\_\_\_ Project Manager: K. Carroll  
 TAT: 10-day  
 DATE: 2-7-14 PAGE 2 OF 2  
 ESN PROJECT # D1402050037  
 LOCATION/PROJECT NAME: BV - Radford High School  
 COLLECTOR: \_\_\_\_\_ DATE COLLECTED: 2/4-2/5

| Sample ID#     | Date | Time | Sample Type | Container Type | X | Comments | # of Containers |
|----------------|------|------|-------------|----------------|---|----------|-----------------|
| 1 RHS-DU-17    | 2/4  | 1436 | soil        |                | X |          | 1               |
| 2 RHS-DU-17.2  | 2/4  | 1449 |             |                | X |          | 1               |
| 3 RHS-DU-17.3  | 2/4  | 1502 |             |                | X |          | 1               |
| 4 RHS-DU-18    | 2/4  | 1300 |             |                | X |          | 1               |
| 5 RHS-DU-19    | 2/4  | 1422 |             |                | X |          | 1               |
| 6 RHS-DU-20    | 2/4  | 1146 |             |                | X |          | 1               |
| 7 RHS-DU-21    | 2/4  | 1226 |             |                | X |          | 1               |
| 8 RHS-DU-22    | 2/4  | 1015 |             |                | X |          | 1               |
| 9 RHS-DU-23    | 2/4  | 1023 |             |                | X |          | 1               |
| 10 RHS-DU-24   | 2/4  | 1050 |             |                | X |          | 1               |
| 11 RHS-DU-25   | 2/4  | 1115 |             |                | X |          | 1               |
| 12 RHS-DU-26   | 2/4  | 1020 |             |                | X |          | 1               |
| 13 RHS-DU-26.2 | 2/4  | 1059 |             |                | X |          | 1               |
| 14 RHS-DU-26.3 | 2/4  | 1145 |             |                | X |          | 1               |
| 15 RHS-DU-27   | 2/4  | 1000 |             |                | X |          | 1               |
| 16 RHS-DU-28   | 2/4  | 1007 |             |                | X |          | 1               |
| 17 RHS-DU-29   | 2/4  | 945  |             |                | X |          | 1               |

Dioxins/Furans 8290

Last Entry

|  |                     |  |                      |   |
|--|---------------------|--|----------------------|---|
| RELINQUISHED BY: (Signature)<br><i>K. Carroll</i>      | DATE/TIME<br>2/7/14 | RECEIVED BY: (Signature)<br><i>Carroll</i> | DATE/TIME<br>2/14/14 | SAMPLE RECEIPT:<br>TOTAL # OF CONTAINERS <u>17</u> (of 35)<br>COC SEALS <u>Y / N / NA</u><br>SEALS INTACT <u>Y / N / NA</u> |
| RELINQUISHED BY: (Signature)                           |                     |  |                      | RECEIVED TEMP: _____  |
| LABORATORY NOTES:<br>Samples MI prepped at ESN Pacific |                     |  |                      |   |



Sample Condition Upon Receipt

Client Name: ESN Pacific

Project #:

WO#: 10257348



Courier: Fed Ex, UPS, USPS, Client, Commercial, Pace, Other

Tracking Number: 1Z V23 U3F B 9904 0403

Custody Seal on Cooler/Box Present? Yes No, Seals Intact? Yes No

Optional: Proj. Due Date: Proj. Name:

Packing Material: Bubble Wrap, Bubble Bags, None, Other

Temp Blank? Yes No

Thermom. Used: 80512447, 72337080

B88A912167504, B88A9132501491

Type of Ice: Wet, Blue, None, Samples on ice, cooling process has begun

Cooler Temp Read (°C): 24, Temp should be above freezing to 6°C

Cooler Temp Corrected (°C): 24, Correction Factor: TRUE

Biological Tissue Frozen? Yes No N/A

Date and Initials of Person Examining Contents: CMB 2/11/14

Comments:

Table with 15 rows of inspection items and checkboxes. Includes items like 'Chain of Custody Present?', 'Samples Arrived within Hold Time?', 'Short Hold Time Analysis (<72 hr)?', 'Sufficient Volume?', 'Containers Intact?', 'Sample Labels Match COC?', 'Filtered Volume Received for Dissolved Tests?', 'Sample #', 'Initial when completed:', 'Lot # of added preservative:'.

CLIENT NOTIFICATION/RESOLUTION

Field Data Required? Yes No

Person Contacted: Date/Time:

Comments/Resolution:

Project Manager Review:

Date: 02/11/14

Note: Whenever there is a discrepancy affecting North Carolina compliance samples, a copy of this form will be sent to the North Carolina DEHNR Certification Office (be out of no d. incorrect preservative, out of temp, incorrect containers)

## Reporting Flags

- A = Reporting Limit based on signal to noise
- B = Less than 10x higher than method blank level
- C = Result obtained from confirmation analysis
- D = Result obtained from analysis of diluted sample
- E = Exceeds calibration range
- I = Interference present
- J = Estimated value
- Nn = Value obtained from additional analysis
- P = PCDE Interference
- R = Recovery outside target range
- S = Peak saturated
- U = Analyte not detected
- V = Result verified by confirmation analysis
- X = %D Exceeds limits
- Y = Calculated using average of daily RFs
- \* = See Discussion

### REPORT OF LABORATORY ANALYSIS

Report No.....10257348

## Appendix B

### Sample Analysis Summary

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-17                 |           |                  |  |
| Lab Sample ID          | 10257348001-S             |           |                  |  |
| Filename               | U140226B_16               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.2 g                    | Matrix    | Soil             |  |
| % Moisture             | 11.0                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.08 g                    | Collected | 02/04/2014 14:36 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140226B_06 & U140226B_20 | Extracted | 02/24/2014 21:00 |  |
| Method Blank ID        | BLANK-39446               | Analyzed  | 02/27/2014 05:14 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 190.0      | ----       | 0.26     | 2,3,7,8-TCDF-13C         | 2.00       | 86 Y             |
| Total TCDF          | 5000.0     | ----       | 0.26 E   | 2,3,7,8-TCDD-13C         | 2.00       | 91               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 96 Y             |
| 2,3,7,8-TCDD        | 9.7        | ----       | 0.17 Y   | 2,3,4,7,8-PeCDF-13C      | 2.00       | 92 Y             |
| Total TCDD          | 290.0      | ----       | 0.17 Y   | 1,2,3,7,8-PeCDD-13C      | 2.00       | 91 Y             |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 102              |
| 1,2,3,7,8-PeCDF     | 130.0      | ----       | 0.51     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 104              |
| 2,3,4,7,8-PeCDF     | 250.0      | ----       | 0.55     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 101              |
| Total PeCDF         | 3100.0     | ----       | 0.53     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 95               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 80               |
| 1,2,3,7,8-PeCDD     | 34.0       | ----       | 0.29     | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 73               |
| Total PeCDD         | 430.0      | ----       | 0.29     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 68               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 66               |
| 1,2,3,4,7,8-HxCDF   | 170.0      | ----       | 0.34     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 66               |
| 1,2,3,6,7,8-HxCDF   | -----      | 150        | 0.34 P   | OCDD-13C                 | 4.00       | 85               |
| 2,3,4,6,7,8-HxCDF   | 190.0      | ----       | 0.29     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 33.0       | ----       | 0.55     | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 1700.0     | ----       | 0.38     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 30.0       | ----       | 0.40     | 2,3,7,8-TCDD-37Cl4       | 0.20       | 102              |
| 1,2,3,6,7,8-HxCDD   | 71.0       | ----       | 0.61     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 62.0       | ----       | 0.35     |                          |            |                  |
| Total HxCDD         | 850.0      | ----       | 0.46     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 480.0      | ----       | 0.28     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 35.0       | ----       | 0.44     | Equivalence: 230 ng/Kg   |            |                  |
| Total HpCDF         | 880.0      | ----       | 0.36     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 1000.0     | ----       | 0.77     |                          |            |                  |
| Total HpCDD         | 2000.0     | ----       | 0.77     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 650.0      | ----       | 0.20 Y   |                          |            |                  |
| OCDD                | 6900.0     | ----       | 0.29 E   |                          |            |                  |

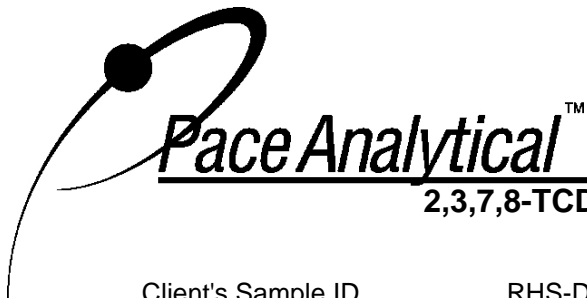
Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
P = PCDE Interference  
E = Exceeds calibration range  
Y = Calculated using average of daily RFs

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-17                 |           |                  |  |
| Lab Sample ID          | 10257348001-S             |           |                  |  |
| Filename               | U140226B_16               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.2 g                    | Matrix    | Soil             |  |
| % Moisture             | 11.0                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.08 g                    | Collected | 02/04/2014 14:36 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140226B_06 & U140226B_20 | Extracted | 02/24/2014 21:00 |  |
| Method Blank ID        | BLANK-39446               | Analyzed  | 02/27/2014 05:14 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB      | MB      | UB      |
|---------------------|---------------|-------------|---------|---------|---------|---------|
| 2,3,7,8-TCDF        | 190.0         | 0.26        | 0.10000 | 19.4520 | 19.4520 | 19.4520 |
| Total TCDF          | 5000.0        | 0.26        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 2,3,7,8-TCDD        | 9.7           | 0.17        | 1.00000 | 9.6688  | 9.6688  | 9.6688  |
| Total TCDD          | 290.0         | 0.17        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDF     | 130.0         | 0.51        | 0.03000 | 3.8938  | 3.8938  | 3.8938  |
| 2,3,4,7,8-PeCDF     | 250.0         | 0.55        | 0.30000 | 75.6271 | 75.6271 | 75.6271 |
| Total PeCDF         | 3100.0        | 0.53        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDD     | 34.0          | 0.29        | 1.00000 | 34.1050 | 34.1050 | 34.1050 |
| Total PeCDD         | 430.0         | 0.29        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDF   | 170.0         | 0.34        | 0.10000 | 16.7288 | 16.7288 | 16.7288 |
| 1,2,3,6,7,8-HxCDF   | ND            | 0.34        | 0.10000 | 15.0064 | 15.0064 | 15.0064 |
| 2,3,4,6,7,8-HxCDF   | 190.0         | 0.29        | 0.10000 | 19.3926 | 19.3926 | 19.3926 |
| 1,2,3,7,8,9-HxCDF   | 33.0          | 0.55        | 0.10000 | 3.3375  | 3.3375  | 3.3375  |
| Total HxCDF         | 1700.0        | 0.38        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDD   | 30.0          | 0.40        | 0.10000 | 3.0218  | 3.0218  | 3.0218  |
| 1,2,3,6,7,8-HxCDD   | 71.0          | 0.61        | 0.10000 | 7.1371  | 7.1371  | 7.1371  |
| 1,2,3,7,8,9-HxCDD   | 62.0          | 0.35        | 0.10000 | 6.1690  | 6.1690  | 6.1690  |
| Total HxCDD         | 850.0         | 0.46        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDF | 480.0         | 0.28        | 0.01000 | 4.7697  | 4.7697  | 4.7697  |
| 1,2,3,4,7,8,9-HpCDF | 35.0          | 0.44        | 0.01000 | 0.3527  | 0.3527  | 0.3527  |
| Total HpCDF         | 880.0         | 0.36        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDD | 1000.0        | 0.77        | 0.01000 | 10.4521 | 10.4521 | 10.4521 |
| Total HpCDD         | 2000.0        | 0.77        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| OCDF                | 650.0         | 0.20        | 0.00030 | 0.1955  | 0.1955  | 0.1955  |
| OCDD                | 6900.0        | 0.29        | 0.00030 | 2.0742  | 2.0742  | 2.0742  |

**230 ng/Kg      230 ng/Kg      230 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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### Method 8290 Sample Analysis Results

Client - ESN Pacific

|                        |                           |           |                  |  |  |
|------------------------|---------------------------|-----------|------------------|--|--|
| Client's Sample ID     | RHS-DU-17.2               |           |                  |  |  |
| Lab Sample ID          | 10257348002-S             |           |                  |  |  |
| Filename               | P140301A_08               |           |                  |  |  |
| Injected By            | BAL                       |           |                  |  |  |
| Total Amount Extracted | 10.3 g                    | Matrix    | Soil             |  |  |
| % Moisture             | 7.8                       | Dilution  | NA               |  |  |
| Dry Weight Extracted   | 9.50 g                    | Collected | 02/04/2014 14:49 |  |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |  |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Extracted | 02/26/2014 21:30 |  |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 12:12 |  |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 270        | ----       | 0.98     | 2,3,7,8-TCDF-13C         | 2.00       | 73               |
| Total TCDF          | 6500       | ----       | 0.98 E   | 2,3,7,8-TCDD-13C         | 2.00       | 79               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 66               |
| 2,3,7,8-TCDD        | 15         | ----       | 1.00     | 2,3,4,7,8-PeCDF-13C      | 2.00       | 64               |
| Total TCDD          | 400        | ----       | 1.00     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 65               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 89               |
| 1,2,3,7,8-PeCDF     | 200        | ----       | 2.80     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 85               |
| 2,3,4,7,8-PeCDF     | 350        | ----       | 3.50     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 82               |
| Total PeCDF         | 3900       | ----       | 3.20     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 74               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 81               |
| 1,2,3,7,8-PeCDD     | 38         | ----       | 1.20     | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 70               |
| Total PeCDD         | 480        | ----       | 1.20     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 67               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 65               |
| 1,2,3,4,7,8-HxCDF   | 220        | ----       | 1.50     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 71               |
| 1,2,3,6,7,8-HxCDF   | 130        | ----       | 2.10     | OCDD-13C                 | 4.00       | 70 Y             |
| 2,3,4,6,7,8-HxCDF   | 270        | ----       | 2.00     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 42         | ----       | 2.90     | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 2300       | ----       | 2.10     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 33         | ----       | 1.90     | 2,3,7,8-TCDD-37Cl4       | 0.20       | 115              |
| 1,2,3,6,7,8-HxCDD   | 70         | ----       | 1.80     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 55         | ----       | 2.50     |                          |            |                  |
| Total HxCDD         | 790        | ----       | 2.10     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 810        | ----       | 1.70     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 40         | ----       | 3.40     | Equivalence: 290 ng/Kg   |            |                  |
| Total HpCDF         | 1200       | ----       | 2.50     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 970        | ----       | 3.50     |                          |            |                  |
| Total HpCDD         | 1900       | ----       | 3.50     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 570        | ----       | 2.20     |                          |            |                  |
| OCDD                | 6200       | ----       | 2.70 E   |                          |            |                  |

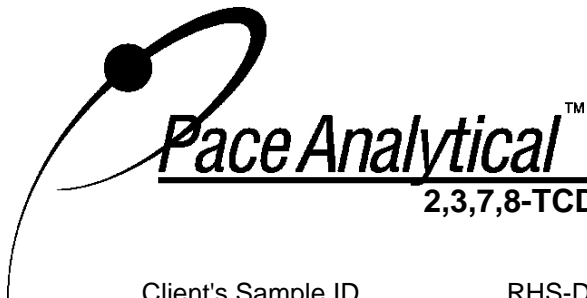
Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
E = Exceeds calibration range  
Y = Calculated using average of daily RfFs

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-17.2               |           |                  |  |
| Lab Sample ID          | 10257348002-S             |           |                  |  |
| Filename               | P140301A_08               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.3 g                    | Matrix    | Soil             |  |
| % Moisture             | 7.8                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.50 g                    | Collected | 02/04/2014 14:49 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 12:12 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB       | MB       | UB       |
|---------------------|---------------|-------------|---------|----------|----------|----------|
| 2,3,7,8-TCDF        | 270           | 0.98        | 0.10000 | 27.3743  | 27.3743  | 27.3743  |
| Total TCDF          | 6500          | 0.98        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 2,3,7,8-TCDD        | 15            | 1.00        | 1.00000 | 14.5551  | 14.5551  | 14.5551  |
| Total TCDD          | 400           | 1.00        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,7,8-PeCDF     | 200           | 2.8         | 0.03000 | 6.0424   | 6.0424   | 6.0424   |
| 2,3,4,7,8-PeCDF     | 350           | 3.5         | 0.30000 | 104.3071 | 104.3071 | 104.3071 |
| Total PeCDF         | 3900          | 3.2         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,7,8-PeCDD     | 38            | 1.2         | 1.00000 | 37.5398  | 37.5398  | 37.5398  |
| Total PeCDD         | 480           | 1.2         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,7,8-HxCDF   | 220           | 1.5         | 0.10000 | 21.7027  | 21.7027  | 21.7027  |
| 1,2,3,6,7,8-HxCDF   | 130           | 2.1         | 0.10000 | 12.7962  | 12.7962  | 12.7962  |
| 2,3,4,6,7,8-HxCDF   | 270           | 2.0         | 0.10000 | 26.9467  | 26.9467  | 26.9467  |
| 1,2,3,7,8,9-HxCDF   | 42            | 2.9         | 0.10000 | 4.1947   | 4.1947   | 4.1947   |
| Total HxCDF         | 2300          | 2.1         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,7,8-HxCDD   | 33            | 1.9         | 0.10000 | 3.2895   | 3.2895   | 3.2895   |
| 1,2,3,6,7,8-HxCDD   | 70            | 1.8         | 0.10000 | 7.0085   | 7.0085   | 7.0085   |
| 1,2,3,7,8,9-HxCDD   | 55            | 2.5         | 0.10000 | 5.4564   | 5.4564   | 5.4564   |
| Total HxCDD         | 790           | 2.1         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,6,7,8-HpCDF | 810           | 1.7         | 0.01000 | 8.0837   | 8.0837   | 8.0837   |
| 1,2,3,4,7,8,9-HpCDF | 40            | 3.4         | 0.01000 | 0.3984   | 0.3984   | 0.3984   |
| Total HpCDF         | 1200          | 2.5         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,6,7,8-HpCDD | 970           | 3.5         | 0.01000 | 9.6899   | 9.6899   | 9.6899   |
| Total HpCDD         | 1900          | 3.5         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| OCDF                | 570           | 2.2         | 0.00030 | 0.1702   | 0.1702   | 0.1702   |
| OCDD                | 6200          | 2.7         | 0.00030 | 1.8702   | 1.8702   | 1.8702   |

**290 ng/Kg      290 ng/Kg      290 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-17.3               |           |                  |  |
| Lab Sample ID          | 10257348003-S             |           |                  |  |
| Filename               | P140301A_09               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.2 g                    | Matrix    | Soil             |  |
| % Moisture             | 9.0                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.28 g                    | Collected | 02/04/2014 15:02 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 12:54 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg |   | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|---|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 260        | ----       | 0.33     |   | 2,3,7,8-TCDF-13C         | 2.00       | 73               |
| Total TCDF          | 5100       | ----       | 0.33     | E | 2,3,7,8-TCDD-13C         | 2.00       | 68               |
|                     |            |            |          |   | 1,2,3,7,8-PeCDF-13C      | 2.00       | 61               |
| 2,3,7,8-TCDD        | -----      | 12         | 0.28     | I | 2,3,4,7,8-PeCDF-13C      | 2.00       | 71               |
| Total TCDD          | 430        | ----       | 0.28     |   | 1,2,3,7,8-PeCDD-13C      | 2.00       | 77               |
|                     |            |            |          |   | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 115              |
| 1,2,3,7,8-PeCDF     | 180        | ----       | 1.40     |   | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 107              |
| 2,3,4,7,8-PeCDF     | 320        | ----       | 0.60     |   | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 100              |
| Total PeCDF         | 3800       | ----       | 1.00     |   | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 55               |
|                     |            |            |          |   | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 98               |
| 1,2,3,7,8-PeCDD     | 34         | ----       | 0.38     |   | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 85               |
| Total PeCDD         | 410        | ----       | 0.38     |   | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 70               |
|                     |            |            |          |   | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 57               |
| 1,2,3,4,7,8-HxCDF   | 190        | ----       | 0.71     |   | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 71               |
| 1,2,3,6,7,8-HxCDF   | -----      | 200        | 0.53     | P | OCDD-13C                 | 4.00       | 49 Y             |
| 2,3,4,6,7,8-HxCDF   | 270        | ----       | 0.60     |   |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 39         | ----       | 0.68     |   | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 2300       | ----       | 0.63     |   | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |   |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 34         | ----       | 1.10     |   | 2,3,7,8-TCDD-37Cl4       | 0.20       | 133              |
| 1,2,3,6,7,8-HxCDD   | 80         | ----       | 0.95     |   |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 53         | ----       | 0.87     |   |                          |            |                  |
| Total HxCDD         | 440        | ----       | 0.98     |   |                          |            |                  |
|                     |            |            |          |   |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 840        | ----       | 0.63     |   | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 41         | ----       | 0.81     |   | Equivalence: 280 ng/Kg   |            |                  |
| Total HpCDF         | 1300       | ----       | 0.72     |   | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |   |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 1300       | ----       | 0.80     |   |                          |            |                  |
| Total HpCDD         | 2600       | ----       | 0.80     |   |                          |            |                  |
|                     |            |            |          |   |                          |            |                  |
| OCDF                | 730        | ----       | 1.40     |   |                          |            |                  |
| OCDD                | 9700       | ----       | 1.20     | E |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

P = PCDE Interference

E = Exceeds calibration range

I = Interference present

Y = Calculated using average of daily RFs

**REPORT OF LABORATORY ANALYSIS**

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-17.3               |           |                  |  |
| Lab Sample ID          | 10257348003-S             |           |                  |  |
| Filename               | P140301A_09               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.2 g                    | Matrix    | Soil             |  |
| % Moisture             | 9.0                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.28 g                    | Collected | 02/04/2014 15:02 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 12:54 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB      | MB      | UB      |
|---------------------|---------------|-------------|---------|---------|---------|---------|
| 2,3,7,8-TCDF        | 260           | 0.33        | 0.10000 | 25.6769 | 25.6769 | 25.6769 |
| Total TCDF          | 5100          | 0.33        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 2,3,7,8-TCDD        | ND            | 0.28        | 1.00000 | 11.8140 | 11.8140 | 11.8140 |
| Total TCDD          | 430           | 0.28        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDF     | 180           | 1.4         | 0.03000 | 5.2986  | 5.2986  | 5.2986  |
| 2,3,4,7,8-PeCDF     | 320           | 0.60        | 0.30000 | 95.6536 | 95.6536 | 95.6536 |
| Total PeCDF         | 3800          | 1.0         | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,7,8-PeCDD     | 34            | 0.38        | 1.00000 | 34.3129 | 34.3129 | 34.3129 |
| Total PeCDD         | 410           | 0.38        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDF   | 190           | 0.71        | 0.10000 | 18.6103 | 18.6103 | 18.6103 |
| 1,2,3,6,7,8-HxCDF   | ND            | 0.53        | 0.10000 | 20.4754 | 20.4754 | 20.4754 |
| 2,3,4,6,7,8-HxCDF   | 270           | 0.60        | 0.10000 | 27.4135 | 27.4135 | 27.4135 |
| 1,2,3,7,8,9-HxCDF   | 39            | 0.68        | 0.10000 | 3.9382  | 3.9382  | 3.9382  |
| Total HxCDF         | 2300          | 0.63        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,7,8-HxCDD   | 34            | 1.1         | 0.10000 | 3.4311  | 3.4311  | 3.4311  |
| 1,2,3,6,7,8-HxCDD   | 80            | 0.95        | 0.10000 | 7.9997  | 7.9997  | 7.9997  |
| 1,2,3,7,8,9-HxCDD   | 53            | 0.87        | 0.10000 | 5.2647  | 5.2647  | 5.2647  |
| Total HxCDD         | 440           | 0.98        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDF | 840           | 0.63        | 0.01000 | 8.3920  | 8.3920  | 8.3920  |
| 1,2,3,4,7,8,9-HpCDF | 41            | 0.81        | 0.01000 | 0.4076  | 0.4076  | 0.4076  |
| Total HpCDF         | 1300          | 0.72        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| 1,2,3,4,6,7,8-HpCDD | 1300          | 0.80        | 0.01000 | 12.5752 | 12.5752 | 12.5752 |
| Total HpCDD         | 2600          | 0.80        | 0.00000 | 0.0000  | 0.0000  | 0.0000  |
| OCDF                | 730           | 1.4         | 0.00030 | 0.2191  | 0.2191  | 0.2191  |
| OCDD                | 9700          | 1.2         | 0.00030 | 2.9066  | 2.9066  | 2.9066  |

**280 ng/Kg      280 ng/Kg      280 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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### Method 8290 Sample Analysis Results

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-18                 |           |                  |  |
| Lab Sample ID          | 10257348004-S             |           |                  |  |
| Filename               | P140301A_10               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.3 g                    | Matrix    | Soil             |  |
| % Moisture             | 9.5                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.32 g                    | Collected | 02/04/2014 13:00 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 13:37 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg |   | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|---|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 570        | ----       | 0.45     | E | 2,3,7,8-TCDF-13C         | 2.00       | 80               |
| Total TCDF          | 11000      | ----       | 0.45     | E | 2,3,7,8-TCDD-13C         | 2.00       | 87               |
|                     |            |            |          |   | 1,2,3,7,8-PeCDF-13C      | 2.00       | 78               |
| 2,3,7,8-TCDD        | 28         | ----       | 0.49     |   | 2,3,4,7,8-PeCDF-13C      | 2.00       | 75               |
| Total TCDD          | 1200       | ----       | 0.49     |   | 1,2,3,7,8-PeCDD-13C      | 2.00       | 77               |
|                     |            |            |          |   | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 99               |
| 1,2,3,7,8-PeCDF     | 410        | ----       | 0.77     |   | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 96               |
| 2,3,4,7,8-PeCDF     | 800        | ----       | 1.80     |   | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 91               |
| Total PeCDF         | 8500       | ----       | 1.30     |   | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 72               |
|                     |            |            |          |   | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 87               |
| 1,2,3,7,8-PeCDD     | 90         | ----       | 0.44     |   | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 77               |
| Total PeCDD         | 1400       | ----       | 0.44     |   | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 75               |
|                     |            |            |          |   | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 68               |
| 1,2,3,4,7,8-HxCDF   | 570        | ----       | 0.90     |   | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 77               |
| 1,2,3,6,7,8-HxCDF   | 420        | ----       | 1.50     |   | OCDD-13C                 | 4.00       | 69 Y             |
| 2,3,4,6,7,8-HxCDF   | 730        | ----       | 1.20     |   |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 100        | ----       | 1.00     |   | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 6000       | ----       | 1.20     |   | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |   |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 84         | ----       | 1.40     |   | 2,3,7,8-TCDD-37Cl4       | 0.20       | 104              |
| 1,2,3,6,7,8-HxCDD   | 160        | ----       | 1.40     |   |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 130        | ----       | 1.10     |   |                          |            |                  |
| Total HxCDD         | 2100       | ----       | 1.30     |   |                          |            |                  |
|                     |            |            |          |   |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 2400       | ----       | 0.84     | E | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 88         | ----       | 1.30     |   | Equivalence: 690 ng/Kg   |            |                  |
| Total HpCDF         | 3100       | ----       | 1.10     | E | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |   |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 1800       | ----       | 2.00     |   |                          |            |                  |
| Total HpCDD         | 3600       | ----       | 2.00     |   |                          |            |                  |
|                     |            |            |          |   |                          |            |                  |
| OCDF                | 850        | ----       | 1.10     |   |                          |            |                  |
| OCDD                | 11000      | ----       | 1.80     | E |                          |            |                  |

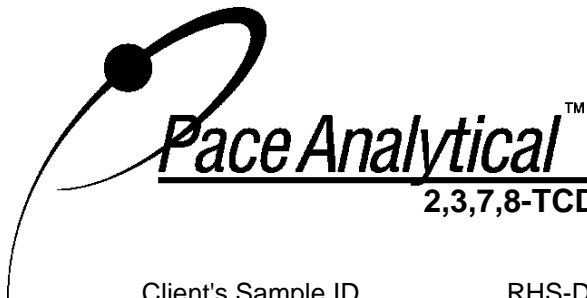
Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
E = Exceeds calibration range  
Y = Calculated using average of daily RfFs

## REPORT OF LABORATORY ANALYSIS

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-18                 |           |                  |  |
| Lab Sample ID          | 10257348004-S             |           |                  |  |
| Filename               | P140301A_10               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.3 g                    | Matrix    | Soil             |  |
| % Moisture             | 9.5                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.32 g                    | Collected | 02/04/2014 13:00 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 13:37 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB       | MB       | UB       |
|---------------------|---------------|-------------|---------|----------|----------|----------|
| 2,3,7,8-TCDF        | 570           | 0.45        | 0.10000 | 56.6262  | 56.6262  | 56.6262  |
| Total TCDF          | 11000         | 0.45        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 2,3,7,8-TCDD        | 28            | 0.49        | 1.00000 | 28.3706  | 28.3706  | 28.3706  |
| Total TCDD          | 1200          | 0.49        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,7,8-PeCDF     | 410           | 0.77        | 0.03000 | 12.2387  | 12.2387  | 12.2387  |
| 2,3,4,7,8-PeCDF     | 800           | 1.8         | 0.30000 | 239.9010 | 239.9010 | 239.9010 |
| Total PeCDF         | 8500          | 1.3         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,7,8-PeCDD     | 90            | 0.44        | 1.00000 | 89.8140  | 89.8140  | 89.8140  |
| Total PeCDD         | 1400          | 0.44        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,7,8-HxCDF   | 570           | 0.90        | 0.10000 | 56.8898  | 56.8898  | 56.8898  |
| 1,2,3,6,7,8-HxCDF   | 420           | 1.5         | 0.10000 | 41.5281  | 41.5281  | 41.5281  |
| 2,3,4,6,7,8-HxCDF   | 730           | 1.2         | 0.10000 | 73.2362  | 73.2362  | 73.2362  |
| 1,2,3,7,8,9-HxCDF   | 100           | 1.0         | 0.10000 | 9.9793   | 9.9793   | 9.9793   |
| Total HxCDF         | 6000          | 1.2         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,7,8-HxCDD   | 84            | 1.4         | 0.10000 | 8.4178   | 8.4178   | 8.4178   |
| 1,2,3,6,7,8-HxCDD   | 160           | 1.4         | 0.10000 | 16.3243  | 16.3243  | 16.3243  |
| 1,2,3,7,8,9-HxCDD   | 130           | 1.1         | 0.10000 | 13.0199  | 13.0199  | 13.0199  |
| Total HxCDD         | 2100          | 1.3         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,6,7,8-HpCDF | 2400          | 0.84        | 0.01000 | 23.8558  | 23.8558  | 23.8558  |
| 1,2,3,4,7,8,9-HpCDF | 88            | 1.3         | 0.01000 | 0.8779   | 0.8779   | 0.8779   |
| Total HpCDF         | 3100          | 1.1         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,6,7,8-HpCDD | 1800          | 2.0         | 0.01000 | 17.5083  | 17.5083  | 17.5083  |
| Total HpCDD         | 3600          | 2.0         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| OCDF                | 850           | 1.1         | 0.00030 | 0.2556   | 0.2556   | 0.2556   |
| OCDD                | 11000         | 1.8         | 0.00030 | 3.1712   | 3.1712   | 3.1712   |

**690 ng/Kg      690 ng/Kg      690 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-19                 |           |                  |  |
| Lab Sample ID          | 10257348005-S             |           |                  |  |
| Filename               | P140301A_11               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.3 g                    | Matrix    | Soil             |  |
| % Moisture             | 7.7                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.51 g                    | Collected | 02/04/2014 14:22 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 14:20 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | -----      | 370        | 0.18 P   | 2,3,7,8-TCDF-13C         | 2.00       | 73               |
| Total TCDF          | 6800       | -----      | 0.18 E   | 2,3,7,8-TCDD-13C         | 2.00       | 79               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 73               |
| 2,3,7,8-TCDD        | 23         | -----      | 0.18     | 2,3,4,7,8-PeCDF-13C      | 2.00       | 70               |
| Total TCDD          | 750        | -----      | 0.18     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 73               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 92               |
| 1,2,3,7,8-PeCDF     | 300        | -----      | 0.44     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 88               |
| 2,3,4,7,8-PeCDF     | 610        | -----      | 0.63     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 82               |
| Total PeCDF         | 6500       | -----      | 0.53     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 52               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 83               |
| 1,2,3,7,8-PeCDD     | 72         | -----      | 0.23     | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 71               |
| Total PeCDD         | 970        | -----      | 0.23     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 66               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 56               |
| 1,2,3,4,7,8-HxCDF   | 410        | -----      | 0.67     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 63               |
| 1,2,3,6,7,8-HxCDF   | 390        | -----      | 0.98     | OCDD-13C                 | 4.00       | 47 Y             |
| 2,3,4,6,7,8-HxCDF   | 600        | -----      | 1.20     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 83         | -----      | 0.80     | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 4500       | -----      | 0.90     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 68         | -----      | 1.60     | 2,3,7,8-TCDD-37Cl4       | 0.20       | 88               |
| 1,2,3,6,7,8-HxCDD   | 120        | -----      | 0.77     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 99         | -----      | 0.64     |                          |            |                  |
| Total HxCDD         | 1500       | -----      | 1.00     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 2100       | -----      | 0.61     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 87         | -----      | 0.96     | Equivalence: 540 ng/Kg   |            |                  |
| Total HpCDF         | 2700       | -----      | 0.78     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 1200       | -----      | 1.40     |                          |            |                  |
| Total HpCDD         | 2400       | -----      | 1.40     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 690        | -----      | 1.00     |                          |            |                  |
| OCDD                | 5600       | -----      | 1.10     |                          |            |                  |

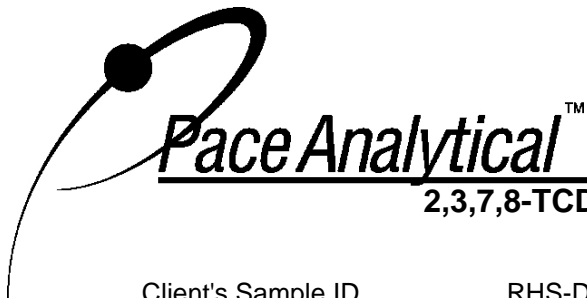
Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
P = PCDE Interference  
E = Exceeds calibration range  
Y = Calculated using average of daily RFs

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-19                 |           |                  |  |
| Lab Sample ID          | 10257348005-S             |           |                  |  |
| Filename               | P140301A_11               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.3 g                    | Matrix    | Soil             |  |
| % Moisture             | 7.7                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.51 g                    | Collected | 02/04/2014 14:22 |  |
| ICAL ID                | P130624                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Extracted | 02/26/2014 21:30 |  |
| Method Blank ID        | BLANK-39465               | Analyzed  | 03/01/2014 14:20 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB       | MB       | UB       |
|---------------------|---------------|-------------|---------|----------|----------|----------|
| 2,3,7,8-TCDF        | ND            | 0.18        | 0.10000 | 36.7462  | 36.7462  | 36.7462  |
| Total TCDF          | 6800          | 0.18        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 2,3,7,8-TCDD        | 23            | 0.18        | 1.00000 | 23.1303  | 23.1303  | 23.1303  |
| Total TCDD          | 750           | 0.18        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,7,8-PeCDF     | 300           | 0.44        | 0.03000 | 9.1296   | 9.1296   | 9.1296   |
| 2,3,4,7,8-PeCDF     | 610           | 0.63        | 0.30000 | 182.6852 | 182.6852 | 182.6852 |
| Total PeCDF         | 6500          | 0.53        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,7,8-PeCDD     | 72            | 0.23        | 1.00000 | 72.4801  | 72.4801  | 72.4801  |
| Total PeCDD         | 970           | 0.23        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,7,8-HxCDF   | 410           | 0.67        | 0.10000 | 41.3009  | 41.3009  | 41.3009  |
| 1,2,3,6,7,8-HxCDF   | 390           | 0.98        | 0.10000 | 38.6998  | 38.6998  | 38.6998  |
| 2,3,4,6,7,8-HxCDF   | 600           | 1.2         | 0.10000 | 60.3981  | 60.3981  | 60.3981  |
| 1,2,3,7,8,9-HxCDF   | 83            | 0.80        | 0.10000 | 8.2573   | 8.2573   | 8.2573   |
| Total HxCDF         | 4500          | 0.90        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,7,8-HxCDD   | 68            | 1.6         | 0.10000 | 6.7685   | 6.7685   | 6.7685   |
| 1,2,3,6,7,8-HxCDD   | 120           | 0.77        | 0.10000 | 12.2455  | 12.2455  | 12.2455  |
| 1,2,3,7,8,9-HxCDD   | 99            | 0.64        | 0.10000 | 9.9064   | 9.9064   | 9.9064   |
| Total HxCDD         | 1500          | 1.00        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,6,7,8-HpCDF | 2100          | 0.61        | 0.01000 | 20.7621  | 20.7621  | 20.7621  |
| 1,2,3,4,7,8,9-HpCDF | 87            | 0.96        | 0.01000 | 0.8747   | 0.8747   | 0.8747   |
| Total HpCDF         | 2700          | 0.78        | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| 1,2,3,4,6,7,8-HpCDD | 1200          | 1.4         | 0.01000 | 11.6360  | 11.6360  | 11.6360  |
| Total HpCDD         | 2400          | 1.4         | 0.00000 | 0.0000   | 0.0000   | 0.0000   |
| OCDF                | 690           | 1.0         | 0.00030 | 0.2064   | 0.2064   | 0.2064   |
| OCDD                | 5600          | 1.1         | 0.00030 | 1.6893   | 1.6893   | 1.6893   |

**540 ng/Kg      540 ng/Kg      540 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |
|------------------------|---------------------------|-----------|------------------|
| Client's Sample ID     | RHS-DU-20                 |           |                  |
| Lab Sample ID          | 10257348006-S             |           |                  |
| Filename               | U140304B_07               |           |                  |
| Injected By            | SMT                       |           |                  |
| Total Amount Extracted | 10.0 g                    | Matrix    | Soil             |
| % Moisture             | 15.1                      | Dilution  | NA               |
| Dry Weight Extracted   | 8.49 g                    | Collected | 02/04/2014 11:46 |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/04/2014 21:10 |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 9.10       | ----       | 0.16     | 2,3,7,8-TCDF-13C         | 2.00       | 96               |
| Total TCDF          | 240.00     | ----       | 0.16     | 2,3,7,8-TCDD-13C         | 2.00       | 100              |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 107              |
| 2,3,7,8-TCDD        | 0.40       | ----       | 0.16 J   | 2,3,4,7,8-PeCDF-13C      | 2.00       | 83 Y             |
| Total TCDD          | 8.50       | ----       | 0.16     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 113              |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 89               |
| 1,2,3,7,8-PeCDF     | 9.60       | ----       | 0.25     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 97               |
| 2,3,4,7,8-PeCDF     | 21.00      | ----       | 0.27     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 103              |
| Total PeCDF         | 270.00     | ----       | 0.26     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 94               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 81               |
| 1,2,3,7,8-PeCDD     | -----      | 2.3        | 0.15 IJ  | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 80               |
| Total PeCDD         | 18.00      | -----      | 0.15     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 80               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 81               |
| 1,2,3,4,7,8-HxCDF   | 16.00      | ----       | 0.38     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 89               |
| 1,2,3,6,7,8-HxCDF   | 18.00      | ----       | 0.26     | OCDD-13C                 | 4.00       | 77               |
| 2,3,4,6,7,8-HxCDF   | 20.00      | ----       | 0.28     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 3.30       | ----       | 0.29 J   | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 200.00     | ----       | 0.30     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 2.50       | ----       | 0.37 J   | 2,3,7,8-TCDD-37Cl4       | 0.20       | 100              |
| 1,2,3,6,7,8-HxCDD   | 7.20       | ----       | 0.52     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 4.80       | ----       | 0.28 J   |                          |            |                  |
| Total HxCDD         | 64.00      | ----       | 0.39     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 72.00      | ----       | 0.16     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 3.90       | ----       | 0.33 J   | Equivalence: 19 ng/Kg    |            |                  |
| Total HpCDF         | 120.00     | ----       | 0.24     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 100.00     | ----       | 0.44     |                          |            |                  |
| Total HpCDD         | 190.00     | ----       | 0.44     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 62.00      | ----       | 0.30     |                          |            |                  |
| OCDD                | 670.00     | ----       | 0.39     |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

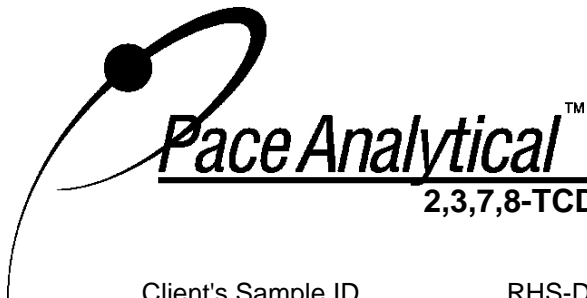
J = Estimated value

I = Interference present

Y = Calculated using average of daily RFs

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |
|------------------------|---------------------------|-----------|------------------|
| Client's Sample ID     | RHS-DU-20                 |           |                  |
| Lab Sample ID          | 10257348006-S             |           |                  |
| Filename               | U140304B_07               |           |                  |
| Injected By            | SMT                       |           |                  |
| Total Amount Extracted | 10.0 g                    | Matrix    | Soil             |
| % Moisture             | 15.1                      | Dilution  | NA               |
| Dry Weight Extracted   | 8.49 g                    | Collected | 02/04/2014 11:46 |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/04/2014 21:10 |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | 9.10          | 0.16        | 0.10000 | 0.9118 | 0.9118 | 0.9118 |
| Total TCDF          | 240.00        | 0.16        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | 0.40          | 0.16        | 1.00000 | 0.4016 | 0.4016 | 0.4016 |
| Total TCDD          | 8.50          | 0.16        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | 9.60          | 0.25        | 0.03000 | 0.2878 | 0.2878 | 0.2878 |
| 2,3,4,7,8-PeCDF     | 21.00         | 0.27        | 0.30000 | 6.3737 | 6.3737 | 6.3737 |
| Total PeCDF         | 270.00        | 0.26        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | ND            | 0.15        | 1.00000 | 2.3376 | 2.3376 | 2.3376 |
| Total PeCDD         | 18.00         | 0.15        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | 16.00         | 0.38        | 0.10000 | 1.6158 | 1.6158 | 1.6158 |
| 1,2,3,6,7,8-HxCDF   | 18.00         | 0.26        | 0.10000 | 1.7658 | 1.7658 | 1.7658 |
| 2,3,4,6,7,8-HxCDF   | 20.00         | 0.28        | 0.10000 | 1.9601 | 1.9601 | 1.9601 |
| 1,2,3,7,8,9-HxCDF   | 3.30          | 0.29        | 0.10000 | 0.3292 | 0.3292 | 0.3292 |
| Total HxCDF         | 200.00        | 0.30        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | 2.50          | 0.37        | 0.10000 | 0.2476 | 0.2476 | 0.2476 |
| 1,2,3,6,7,8-HxCDD   | 7.20          | 0.52        | 0.10000 | 0.7186 | 0.7186 | 0.7186 |
| 1,2,3,7,8,9-HxCDD   | 4.80          | 0.28        | 0.10000 | 0.4806 | 0.4806 | 0.4806 |
| Total HxCDD         | 64.00         | 0.39        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | 72.00         | 0.16        | 0.01000 | 0.7244 | 0.7244 | 0.7244 |
| 1,2,3,4,7,8,9-HpCDF | 3.90          | 0.33        | 0.01000 | 0.0394 | 0.0394 | 0.0394 |
| Total HpCDF         | 120.00        | 0.24        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | 100.00        | 0.44        | 0.01000 | 1.0448 | 1.0448 | 1.0448 |
| Total HpCDD         | 190.00        | 0.44        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | 62.00         | 0.30        | 0.00030 | 0.0186 | 0.0186 | 0.0186 |
| OCDD                | 670.00        | 0.39        | 0.00030 | 0.2002 | 0.2002 | 0.2002 |

**19 ng/Kg      19 ng/Kg      19 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-21                 |           |                  |  |
| Lab Sample ID          | 10257348007-S             |           |                  |  |
| Filename               | U140304B_08               |           |                  |  |
| Injected By            | SMT                       |           |                  |  |
| Total Amount Extracted | 10.5 g                    | Matrix    | Soil             |  |
| % Moisture             | 11.2                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.32 g                    | Collected | 02/04/2014 12:26 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |  |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/04/2014 21:54 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg |    | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|----|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 8.5        | ----       | 0.32     |    | 2,3,7,8-TCDF-13C         | 2.00       | 97               |
| Total TCDF          | 220.0      | ----       | 0.32     |    | 2,3,7,8-TCDD-13C         | 2.00       | 101              |
|                     |            |            |          |    | 1,2,3,7,8-PeCDF-13C      | 2.00       | 110              |
| 2,3,7,8-TCDD        | -----      | 0.35       | 0.26     | IJ | 2,3,4,7,8-PeCDF-13C      | 2.00       | 88 Y             |
| Total TCDD          | 8.8        | ----       | 0.26     |    | 1,2,3,7,8-PeCDD-13C      | 2.00       | 114              |
|                     |            |            |          |    | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 94               |
| 1,2,3,7,8-PeCDF     | 8.6        | ----       | 0.22     |    | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 98               |
| 2,3,4,7,8-PeCDF     | 20.0       | ----       | 0.22     |    | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 105              |
| Total PeCDF         | 230.0      | ----       | 0.22     |    | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 93               |
|                     |            |            |          |    | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 83               |
| 1,2,3,7,8-PeCDD     | 2.2        | ----       | 0.16     | J  | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 80               |
| Total PeCDD         | 19.0       | ----       | 0.16     |    | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 80               |
|                     |            |            |          |    | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 82               |
| 1,2,3,4,7,8-HxCDF   | 14.0       | ----       | 0.45     |    | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 90               |
| 1,2,3,6,7,8-HxCDF   | 15.0       | ----       | 0.39     |    | OCDD-13C                 | 4.00       | 76               |
| 2,3,4,6,7,8-HxCDF   | 19.0       | ----       | 0.31     |    |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | -----      | 2.60       | 0.39     | IJ | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 180.0      | ----       | 0.38     |    | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 2.1        | ----       | 0.40     | J  | 2,3,7,8-TCDD-37Cl4       | 0.20       | 99               |
| 1,2,3,6,7,8-HxCDD   | 6.5        | ----       | 0.37     |    |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 5.1        | ----       | 0.39     | J  |                          |            |                  |
| Total HxCDD         | 64.0       | ----       | 0.39     |    |                          |            |                  |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 68.0       | ----       | 0.28     |    | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 3.4        | ----       | 0.37     | J  | Equivalence: 18 ng/Kg    |            |                  |
| Total HpCDF         | 110.0      | ----       | 0.32     |    | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 94.0       | ----       | 0.40     |    |                          |            |                  |
| Total HpCDD         | 170.0      | ----       | 0.40     |    |                          |            |                  |
|                     |            |            |          |    |                          |            |                  |
| OCDF                | 58.0       | ----       | 0.44     |    |                          |            |                  |
| OCDD                | 620.0      | ----       | 0.58     |    |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

J = Estimated value

I = Interference present

Y = Calculated using average of daily RFs

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |
|------------------------|---------------------------|-----------|------------------|
| Client's Sample ID     | RHS-DU-21                 |           |                  |
| Lab Sample ID          | 10257348007-S             |           |                  |
| Filename               | U140304B_08               |           |                  |
| Injected By            | SMT                       |           |                  |
| Total Amount Extracted | 10.5 g                    | Matrix    | Soil             |
| % Moisture             | 11.2                      | Dilution  | NA               |
| Dry Weight Extracted   | 9.32 g                    | Collected | 02/04/2014 12:26 |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/04/2014 21:54 |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | 8.5           | 0.32        | 0.10000 | 0.8454 | 0.8454 | 0.8454 |
| Total TCDF          | 220.0         | 0.32        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | ND            | 0.26        | 1.00000 | 0.3533 | 0.3533 | 0.3533 |
| Total TCDD          | 8.8           | 0.26        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | 8.6           | 0.22        | 0.03000 | 0.2590 | 0.2590 | 0.2590 |
| 2,3,4,7,8-PeCDF     | 20.0          | 0.22        | 0.30000 | 5.9384 | 5.9384 | 5.9384 |
| Total PeCDF         | 230.0         | 0.22        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | 2.2           | 0.16        | 1.00000 | 2.2467 | 2.2467 | 2.2467 |
| Total PeCDD         | 19.0          | 0.16        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | 14.0          | 0.45        | 0.10000 | 1.4476 | 1.4476 | 1.4476 |
| 1,2,3,6,7,8-HxCDF   | 15.0          | 0.39        | 0.10000 | 1.5036 | 1.5036 | 1.5036 |
| 2,3,4,6,7,8-HxCDF   | 19.0          | 0.31        | 0.10000 | 1.8663 | 1.8663 | 1.8663 |
| 1,2,3,7,8,9-HxCDF   | ND            | 0.39        | 0.10000 | 0.2636 | 0.2636 | 0.2636 |
| Total HxCDF         | 180.0         | 0.38        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | 2.1           | 0.40        | 0.10000 | 0.2106 | 0.2106 | 0.2106 |
| 1,2,3,6,7,8-HxCDD   | 6.5           | 0.37        | 0.10000 | 0.6453 | 0.6453 | 0.6453 |
| 1,2,3,7,8,9-HxCDD   | 5.1           | 0.39        | 0.10000 | 0.5101 | 0.5101 | 0.5101 |
| Total HxCDD         | 64.0          | 0.39        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | 68.0          | 0.28        | 0.01000 | 0.6767 | 0.6767 | 0.6767 |
| 1,2,3,4,7,8,9-HpCDF | 3.4           | 0.37        | 0.01000 | 0.0339 | 0.0339 | 0.0339 |
| Total HpCDF         | 110.0         | 0.32        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | 94.0          | 0.40        | 0.01000 | 0.9413 | 0.9413 | 0.9413 |
| Total HpCDD         | 170.0         | 0.40        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | 58.0          | 0.44        | 0.00030 | 0.0173 | 0.0173 | 0.0173 |
| OCDD                | 620.0         | 0.58        | 0.00030 | 0.1846 | 0.1846 | 0.1846 |

**18 ng/Kg      18 ng/Kg      18 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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### Method 8290 Sample Analysis Results

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-22                 |           |                  |  |
| Lab Sample ID          | 10257348008-S             |           |                  |  |
| Filename               | U140226B_17               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.5 g                    | Matrix    | Soil             |  |
| % Moisture             | 9.9                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.46 g                    | Collected | 02/04/2014 10:15 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140226B_06 & U140226B_20 | Extracted | 02/24/2014 21:00 |  |
| Method Blank ID        | BLANK-39446               | Analyzed  | 02/27/2014 06:01 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 15.0       | ----       | 0.27     | 2,3,7,8-TCDF-13C         | 2.00       | 87 Y             |
| Total TCDF          | 440.0      | ----       | 0.27     | 2,3,7,8-TCDD-13C         | 2.00       | 91               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 94 Y             |
| 2,3,7,8-TCDD        | 1.1        | ----       | 0.14 Y   | 2,3,4,7,8-PeCDF-13C      | 2.00       | 89 Y             |
| Total TCDD          | 45.0       | ----       | 0.14 Y   | 1,2,3,7,8-PeCDD-13C      | 2.00       | 94 Y             |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 102              |
| 1,2,3,7,8-PeCDF     | 15.0       | ----       | 0.23     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 100              |
| 2,3,4,7,8-PeCDF     | 36.0       | ----       | 0.17     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 99               |
| Total PeCDF         | 450.0      | ----       | 0.20     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 91               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 83               |
| 1,2,3,7,8-PeCDD     | 4.8        | ----       | 0.30 J   | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 75               |
| Total PeCDD         | 89.0       | ----       | 0.30     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 64               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 59               |
| 1,2,3,4,7,8-HxCDF   | 31.0       | ----       | 0.22     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 60               |
| 1,2,3,6,7,8-HxCDF   | ----       | 31         | 0.20 P   | OCDD-13C                 | 4.00       | 63               |
| 2,3,4,6,7,8-HxCDF   | 39.0       | ----       | 0.15     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 6.8        | ----       | 0.18     | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 380.0      | ----       | 0.19     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 5.7        | ----       | 0.37     | 2,3,7,8-TCDD-37Cl4       | 0.20       | 101              |
| 1,2,3,6,7,8-HxCDD   | 19.0       | ----       | 0.32     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 12.0       | ----       | 0.36     |                          |            |                  |
| Total HxCDD         | 220.0      | ----       | 0.35     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 150.0      | ----       | 0.24     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 9.7        | ----       | 0.20     | Equivalence: 39 ng/Kg    |            |                  |
| Total HpCDF         | 290.0      | ----       | 0.22     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 300.0      | ----       | 0.66     |                          |            |                  |
| Total HpCDD         | 560.0      | ----       | 0.66     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 170.0      | ----       | 0.48 Y   |                          |            |                  |
| OCDD                | 3000.0     | ----       | 0.48     |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
J = Estimated value  
P = PCDE Interference  
Y = Calculated using average of daily RFs

## REPORT OF LABORATORY ANALYSIS

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-22                 |           |                  |  |
| Lab Sample ID          | 10257348008-S             |           |                  |  |
| Filename               | U140226B_17               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.5 g                    | Matrix    | Soil             |  |
| % Moisture             | 9.9                       | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.46 g                    | Collected | 02/04/2014 10:15 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140226B_06 & U140226B_20 | Extracted | 02/24/2014 21:00 |  |
| Method Blank ID        | BLANK-39446               | Analyzed  | 02/27/2014 06:01 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB              | MB              | UB              |
|---------------------|---------------|-------------|---------|-----------------|-----------------|-----------------|
| 2,3,7,8-TCDF        | 15.0          | 0.27        | 0.10000 | 1.5439          | 1.5439          | 1.5439          |
| Total TCDF          | 440.0         | 0.27        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 2,3,7,8-TCDD        | 1.1           | 0.14        | 1.00000 | 1.0975          | 1.0975          | 1.0975          |
| Total TCDD          | 45.0          | 0.14        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,7,8-PeCDF     | 15.0          | 0.23        | 0.03000 | 0.4583          | 0.4583          | 0.4583          |
| 2,3,4,7,8-PeCDF     | 36.0          | 0.17        | 0.30000 | 10.9110         | 10.9110         | 10.9110         |
| Total PeCDF         | 450.0         | 0.20        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,7,8-PeCDD     | 4.8           | 0.30        | 1.00000 | 4.7781          | 4.7781          | 4.7781          |
| Total PeCDD         | 89.0          | 0.30        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,4,7,8-HxCDF   | 31.0          | 0.22        | 0.10000 | 3.0775          | 3.0775          | 3.0775          |
| 1,2,3,6,7,8-HxCDF   | ND            | 0.20        | 0.10000 | 3.1058          | 3.1058          | 3.1058          |
| 2,3,4,6,7,8-HxCDF   | 39.0          | 0.15        | 0.10000 | 3.8650          | 3.8650          | 3.8650          |
| 1,2,3,7,8,9-HxCDF   | 6.8           | 0.18        | 0.10000 | 0.6755          | 0.6755          | 0.6755          |
| Total HxCDF         | 380.0         | 0.19        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,4,7,8-HxCDD   | 5.7           | 0.37        | 0.10000 | 0.5663          | 0.5663          | 0.5663          |
| 1,2,3,6,7,8-HxCDD   | 19.0          | 0.32        | 0.10000 | 1.9266          | 1.9266          | 1.9266          |
| 1,2,3,7,8,9-HxCDD   | 12.0          | 0.36        | 0.10000 | 1.2018          | 1.2018          | 1.2018          |
| Total HxCDD         | 220.0         | 0.35        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,4,6,7,8-HpCDF | 150.0         | 0.24        | 0.01000 | 1.4927          | 1.4927          | 1.4927          |
| 1,2,3,4,7,8,9-HpCDF | 9.7           | 0.20        | 0.01000 | 0.0974          | 0.0974          | 0.0974          |
| Total HpCDF         | 290.0         | 0.22        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,4,6,7,8-HpCDD | 300.0         | 0.66        | 0.01000 | 2.9534          | 2.9534          | 2.9534          |
| Total HpCDD         | 560.0         | 0.66        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| OCDF                | 170.0         | 0.48        | 0.00030 | 0.0519          | 0.0519          | 0.0519          |
| OCDD                | 3000.0        | 0.48        | 0.00030 | 0.8973          | 0.8973          | 0.8973          |
|                     |               |             |         | <b>39 ng/Kg</b> | <b>39 ng/Kg</b> | <b>39 ng/Kg</b> |

Final values are valid to only 2 significant figures  
 TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
 LB = Lower Bound, Where "ND", TEQ Conc = 0  
 MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
 UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
 RL = Reporting Limit

**REPORT OF LABORATORY ANALYSIS**

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-23                 |           |                  |  |
| Lab Sample ID          | 10257348009-S             |           |                  |  |
| Filename               | U140304B_09               |           |                  |  |
| Injected By            | SMT                       |           |                  |  |
| Total Amount Extracted | 10.4 g                    | Matrix    | Soil             |  |
| % Moisture             | 11.6                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.19 g                    | Collected | 02/04/2014 10:23 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |  |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/04/2014 22:39 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg |    | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|----|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 8.6        | ----       | 0.36     |    | 2,3,7,8-TCDF-13C         | 2.00       | 96               |
| Total TCDF          | 220.0      | ----       | 0.36     |    | 2,3,7,8-TCDD-13C         | 2.00       | 99               |
|                     |            |            |          |    | 1,2,3,7,8-PeCDF-13C      | 2.00       | 112              |
| 2,3,7,8-TCDD        | -----      | 0.43       | 0.40     | IJ | 2,3,4,7,8-PeCDF-13C      | 2.00       | 86 Y             |
| Total TCDD          | 11.0       | ----       | 0.40     |    | 1,2,3,7,8-PeCDD-13C      | 2.00       | 113              |
|                     |            |            |          |    | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 93               |
| 1,2,3,7,8-PeCDF     | 9.0        | ----       | 0.15     |    | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 101              |
| 2,3,4,7,8-PeCDF     | 19.0       | ----       | 0.20     |    | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 105              |
| Total PeCDF         | 240.0      | ----       | 0.18     |    | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 95               |
|                     |            |            |          |    | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 87               |
| 1,2,3,7,8-PeCDD     | 2.0        | ----       | 0.15     | J  | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 79               |
| Total PeCDD         | 26.0       | ----       | 0.15     |    | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 80               |
|                     |            |            |          |    | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 82               |
| 1,2,3,4,7,8-HxCDF   | 16.0       | ----       | 0.48     |    | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 91               |
| 1,2,3,6,7,8-HxCDF   | 16.0       | ----       | 0.24     |    | OCDD-13C                 | 4.00       | 77               |
| 2,3,4,6,7,8-HxCDF   | 18.0       | ----       | 0.21     |    |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 3.6        | ----       | 0.31     | J  | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 190.0      | ----       | 0.31     |    | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 2.0        | ----       | 0.35     | J  | 2,3,7,8-TCDD-37Cl4       | 0.20       | 97               |
| 1,2,3,6,7,8-HxCDD   | 7.3        | ----       | 0.37     |    |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 4.8        | ----       | 0.32     | J  |                          |            |                  |
| Total HxCDD         | 66.0       | ----       | 0.35     |    |                          |            |                  |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 69.0       | ----       | 0.30     |    | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 3.5        | ----       | 0.45     | J  | Equivalence: 18 ng/Kg    |            |                  |
| Total HpCDF         | 110.0      | ----       | 0.37     |    | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 97.0       | ----       | 0.50     |    |                          |            |                  |
| Total HpCDD         | 180.0      | ----       | 0.50     |    |                          |            |                  |
|                     |            |            |          |    |                          |            |                  |
| OCDF                | 57.0       | ----       | 0.96     |    |                          |            |                  |
| OCDD                | 610.0      | ----       | 0.59     |    |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

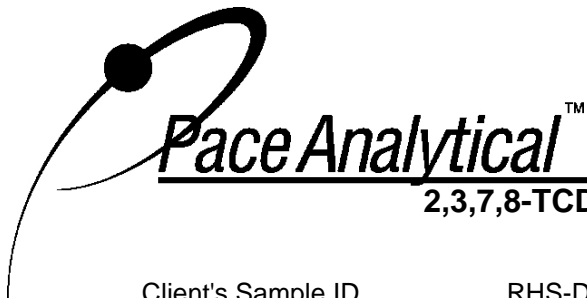
J = Estimated value

I = Interference present

Y = Calculated using average of daily RFs

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |  |
|------------------------|---------------------------|-----------|------------------|--|--|
| Client's Sample ID     | RHS-DU-23                 |           |                  |  |  |
| Lab Sample ID          | 10257348009-S             |           |                  |  |  |
| Filename               | U140304B_09               |           |                  |  |  |
| Injected By            | SMT                       |           |                  |  |  |
| Total Amount Extracted | 10.4 g                    | Matrix    | Soil             |  |  |
| % Moisture             | 11.6                      | Dilution  | NA               |  |  |
| Dry Weight Extracted   | 9.19 g                    | Collected | 02/04/2014 10:23 |  |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |  |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |  |  |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/04/2014 22:39 |  |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | 8.6           | 0.36        | 0.10000 | 0.8634 | 0.8634 | 0.8634 |
| Total TCDF          | 220.0         | 0.36        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | ND            | 0.40        | 1.00000 | 0.4348 | 0.4348 | 0.4348 |
| Total TCDD          | 11.0          | 0.40        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | 9.0           | 0.15        | 0.03000 | 0.2689 | 0.2689 | 0.2689 |
| 2,3,4,7,8-PeCDF     | 19.0          | 0.20        | 0.30000 | 5.7835 | 5.7835 | 5.7835 |
| Total PeCDF         | 240.0         | 0.18        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | 2.0           | 0.15        | 1.00000 | 2.0346 | 2.0346 | 2.0346 |
| Total PeCDD         | 26.0          | 0.15        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | 16.0          | 0.48        | 0.10000 | 1.6048 | 1.6048 | 1.6048 |
| 1,2,3,6,7,8-HxCDF   | 16.0          | 0.24        | 0.10000 | 1.5691 | 1.5691 | 1.5691 |
| 2,3,4,6,7,8-HxCDF   | 18.0          | 0.21        | 0.10000 | 1.8447 | 1.8447 | 1.8447 |
| 1,2,3,7,8,9-HxCDF   | 3.6           | 0.31        | 0.10000 | 0.3598 | 0.3598 | 0.3598 |
| Total HxCDF         | 190.0         | 0.31        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | 2.0           | 0.35        | 0.10000 | 0.2037 | 0.2037 | 0.2037 |
| 1,2,3,6,7,8-HxCDD   | 7.3           | 0.37        | 0.10000 | 0.7264 | 0.7264 | 0.7264 |
| 1,2,3,7,8,9-HxCDD   | 4.8           | 0.32        | 0.10000 | 0.4823 | 0.4823 | 0.4823 |
| Total HxCDD         | 66.0          | 0.35        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | 69.0          | 0.30        | 0.01000 | 0.6929 | 0.6929 | 0.6929 |
| 1,2,3,4,7,8,9-HpCDF | 3.5           | 0.45        | 0.01000 | 0.0351 | 0.0351 | 0.0351 |
| Total HpCDF         | 110.0         | 0.37        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | 97.0          | 0.50        | 0.01000 | 0.9730 | 0.9730 | 0.9730 |
| Total HpCDD         | 180.0         | 0.50        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | 57.0          | 0.96        | 0.00030 | 0.0171 | 0.0171 | 0.0171 |
| OCDD                | 610.0         | 0.59        | 0.00030 | 0.1825 | 0.1825 | 0.1825 |

**18 ng/Kg      18 ng/Kg      18 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-24                 |           |                  |  |
| Lab Sample ID          | 10257348010-S             |           |                  |  |
| Filename               | U140304B_10               |           |                  |  |
| Injected By            | SMT                       |           |                  |  |
| Total Amount Extracted | 11.4 g                    | Matrix    | Soil             |  |
| % Moisture             | 14.5                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.75 g                    | Collected | 02/04/2014 10:50 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |  |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/04/2014 23:24 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 8.2        | ----       | 0.47     | 2,3,7,8-TCDF-13C         | 2.00       | 97               |
| Total TCDF          | 210.0      | ----       | 0.47     | 2,3,7,8-TCDD-13C         | 2.00       | 100              |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 111              |
| 2,3,7,8-TCDD        | ND         | ----       | 0.39     | 2,3,4,7,8-PeCDF-13C      | 2.00       | 85 Y             |
| Total TCDD          | 9.0        | ----       | 0.39     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 114              |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 98               |
| 1,2,3,7,8-PeCDF     | 8.3        | ----       | 0.26     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 97               |
| 2,3,4,7,8-PeCDF     | 20.0       | ----       | 0.20     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 107              |
| Total PeCDF         | 230.0      | ----       | 0.23     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 96               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 86               |
| 1,2,3,7,8-PeCDD     | 2.1        | ----       | 0.16 J   | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 77               |
| Total PeCDD         | 22.0       | ----       | 0.16     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 80               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 82               |
| 1,2,3,4,7,8-HxCDF   | 13.0       | ----       | 0.45     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 90               |
| 1,2,3,6,7,8-HxCDF   | 15.0       | ----       | 0.38     | OCDD-13C                 | 4.00       | 76               |
| 2,3,4,6,7,8-HxCDF   | 18.0       | ----       | 0.38     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | ----       | 2.5        | 0.19 IJ  | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 170.0      | ----       | 0.35     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | ----       | 2.0        | 0.13 IJ  | 2,3,7,8-TCDD-37Cl4       | 0.20       | 104              |
| 1,2,3,6,7,8-HxCDD   | 6.9        | ----       | 0.35     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 4.8        | ----       | 0.45 J   |                          |            |                  |
| Total HxCDD         | 61.0       | ----       | 0.31     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 66.0       | ----       | 0.30     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 3.5        | ----       | 0.34 J   | Equivalence: 17 ng/Kg    |            |                  |
| Total HpCDF         | 110.0      | ----       | 0.32     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 92.0       | ----       | 0.50     |                          |            |                  |
| Total HpCDD         | 170.0      | ----       | 0.50     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 54.0       | ----       | 0.68     |                          |            |                  |
| OCDD                | 580.0      | ----       | 0.58     |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

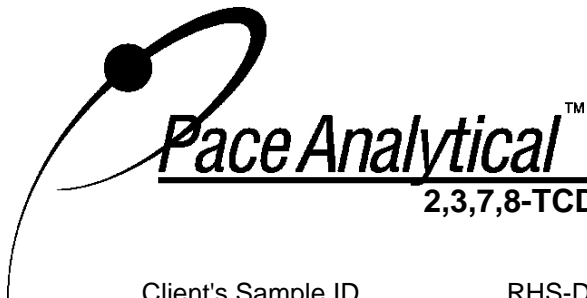
J = Estimated value

I = Interference present

Y = Calculated using average of daily RFs

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-24                 |           |                  |  |
| Lab Sample ID          | 10257348010-S             |           |                  |  |
| Filename               | U140304B_10               |           |                  |  |
| Injected By            | SMT                       |           |                  |  |
| Total Amount Extracted | 11.4 g                    | Matrix    | Soil             |  |
| % Moisture             | 14.5                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 9.75 g                    | Collected | 02/04/2014 10:50 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |  |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/04/2014 23:24 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | 8.2           | 0.47        | 0.10000 | 0.8214 | 0.8214 | 0.8214 |
| Total TCDF          | 210.0         | 0.47        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | ND            | 0.39        | 1.00000 | 0.0000 | 0.1959 | 0.3918 |
| Total TCDD          | 9.0           | 0.39        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | 8.3           | 0.26        | 0.03000 | 0.2504 | 0.2504 | 0.2504 |
| 2,3,4,7,8-PeCDF     | 20.0          | 0.20        | 0.30000 | 5.9093 | 5.9093 | 5.9093 |
| Total PeCDF         | 230.0         | 0.23        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | 2.1           | 0.16        | 1.00000 | 2.1120 | 2.1120 | 2.1120 |
| Total PeCDD         | 22.0          | 0.16        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | 13.0          | 0.45        | 0.10000 | 1.2935 | 1.2935 | 1.2935 |
| 1,2,3,6,7,8-HxCDF   | 15.0          | 0.38        | 0.10000 | 1.5347 | 1.5347 | 1.5347 |
| 2,3,4,6,7,8-HxCDF   | 18.0          | 0.38        | 0.10000 | 1.8307 | 1.8307 | 1.8307 |
| 1,2,3,7,8,9-HxCDF   | ND            | 0.19        | 0.10000 | 0.2511 | 0.2511 | 0.2511 |
| Total HxCDF         | 170.0         | 0.35        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | ND            | 0.13        | 0.10000 | 0.1992 | 0.1992 | 0.1992 |
| 1,2,3,6,7,8-HxCDD   | 6.9           | 0.35        | 0.10000 | 0.6950 | 0.6950 | 0.6950 |
| 1,2,3,7,8,9-HxCDD   | 4.8           | 0.45        | 0.10000 | 0.4797 | 0.4797 | 0.4797 |
| Total HxCDD         | 61.0          | 0.31        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | 66.0          | 0.30        | 0.01000 | 0.6597 | 0.6597 | 0.6597 |
| 1,2,3,4,7,8,9-HpCDF | 3.5           | 0.34        | 0.01000 | 0.0355 | 0.0355 | 0.0355 |
| Total HpCDF         | 110.0         | 0.32        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | 92.0          | 0.50        | 0.01000 | 0.9189 | 0.9189 | 0.9189 |
| Total HpCDD         | 170.0         | 0.50        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | 54.0          | 0.68        | 0.00030 | 0.0161 | 0.0161 | 0.0161 |
| OCDD                | 580.0         | 0.58        | 0.00030 | 0.1740 | 0.1740 | 0.1740 |

**17 ng/Kg      17 ng/Kg      18 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-25                 |           |                  |  |
| Lab Sample ID          | 10257348011-S             |           |                  |  |
| Filename               | U140304B_11               |           |                  |  |
| Injected By            | SMT                       |           |                  |  |
| Total Amount Extracted | 10.0 g                    | Matrix    | Soil             |  |
| % Moisture             | 14.9                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 8.51 g                    | Collected | 02/04/2014 11:15 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |  |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/05/2014 00:09 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 9.5        | ----       | 0.48     | 2,3,7,8-TCDF-13C         | 2.00       | 97               |
| Total TCDF          | 240.0      | ----       | 0.48     | 2,3,7,8-TCDD-13C         | 2.00       | 102              |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 111              |
| 2,3,7,8-TCDD        | ND         | ----       | 0.42     | 2,3,4,7,8-PeCDF-13C      | 2.00       | 86 Y             |
| Total TCDD          | 10.0       | ----       | 0.42     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 116              |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 95               |
| 1,2,3,7,8-PeCDF     | 9.5        | ----       | 0.36     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 93               |
| 2,3,4,7,8-PeCDF     | 22.0       | ----       | 0.43     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 104              |
| Total PeCDF         | 260.0      | ----       | 0.40     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 96               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 86               |
| 1,2,3,7,8-PeCDD     | 2.5        | ----       | 0.25 J   | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 79               |
| Total PeCDD         | 26.0       | ----       | 0.25     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 81               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 83               |
| 1,2,3,4,7,8-HxCDF   | 15.0       | ----       | 0.64     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 91               |
| 1,2,3,6,7,8-HxCDF   | 18.0       | ----       | 0.61     | OCDD-13C                 | 4.00       | 75               |
| 2,3,4,6,7,8-HxCDF   | 21.0       | ----       | 0.53     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 3.5        | ----       | 0.59 J   | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 210.0      | ----       | 0.59     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 2.7        | ----       | 1.00 J   | 2,3,7,8-TCDD-37Cl4       | 0.20       | 102              |
| 1,2,3,6,7,8-HxCDD   | 7.6        | ----       | 0.44     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | ----       | 4.6        | 0.69 IJ  |                          |            |                  |
| Total HxCDD         | 65.0       | ----       | 0.72     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 74.0       | ----       | 0.19     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 3.5        | ----       | 0.40 J   | Equivalence: 20 ng/Kg    |            |                  |
| Total HpCDF         | 110.0      | ----       | 0.30     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 110.0      | ----       | 0.50     |                          |            |                  |
| Total HpCDD         | 190.0      | ----       | 0.50     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 61.0       | ----       | 0.86     |                          |            |                  |
| OCDD                | 660.0      | ----       | 0.74     |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

J = Estimated value  
I = Interference present  
Y = Calculated using average of daily RFs

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |
|------------------------|---------------------------|-----------|------------------|
| Client's Sample ID     | RHS-DU-25                 |           |                  |
| Lab Sample ID          | 10257348011-S             |           |                  |
| Filename               | U140304B_11               |           |                  |
| Injected By            | SMT                       |           |                  |
| Total Amount Extracted | 10.0 g                    | Matrix    | Soil             |
| % Moisture             | 14.9                      | Dilution  | NA               |
| Dry Weight Extracted   | 8.51 g                    | Collected | 02/04/2014 11:15 |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/05/2014 00:09 |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | 9.5           | 0.48        | 0.10000 | 0.9458 | 0.9458 | 0.9458 |
| Total TCDF          | 240.0         | 0.48        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | ND            | 0.42        | 1.00000 | 0.0000 | 0.2122 | 0.4243 |
| Total TCDD          | 10.0          | 0.42        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | 9.5           | 0.36        | 0.03000 | 0.2842 | 0.2842 | 0.2842 |
| 2,3,4,7,8-PeCDF     | 22.0          | 0.43        | 0.30000 | 6.7302 | 6.7302 | 6.7302 |
| Total PeCDF         | 260.0         | 0.40        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | 2.5           | 0.25        | 1.00000 | 2.4768 | 2.4768 | 2.4768 |
| Total PeCDD         | 26.0          | 0.25        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | 15.0          | 0.64        | 0.10000 | 1.5121 | 1.5121 | 1.5121 |
| 1,2,3,6,7,8-HxCDF   | 18.0          | 0.61        | 0.10000 | 1.8005 | 1.8005 | 1.8005 |
| 2,3,4,6,7,8-HxCDF   | 21.0          | 0.53        | 0.10000 | 2.0779 | 2.0779 | 2.0779 |
| 1,2,3,7,8,9-HxCDF   | 3.5           | 0.59        | 0.10000 | 0.3544 | 0.3544 | 0.3544 |
| Total HxCDF         | 210.0         | 0.59        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | 2.7           | 1.0         | 0.10000 | 0.2728 | 0.2728 | 0.2728 |
| 1,2,3,6,7,8-HxCDD   | 7.6           | 0.44        | 0.10000 | 0.7609 | 0.7609 | 0.7609 |
| 1,2,3,7,8,9-HxCDD   | ND            | 0.69        | 0.10000 | 0.4614 | 0.4614 | 0.4614 |
| Total HxCDD         | 65.0          | 0.72        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | 74.0          | 0.19        | 0.01000 | 0.7394 | 0.7394 | 0.7394 |
| 1,2,3,4,7,8,9-HpCDF | 3.5           | 0.40        | 0.01000 | 0.0346 | 0.0346 | 0.0346 |
| Total HpCDF         | 110.0         | 0.30        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | 110.0         | 0.50        | 0.01000 | 1.0506 | 1.0506 | 1.0506 |
| Total HpCDD         | 190.0         | 0.50        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | 61.0          | 0.86        | 0.00030 | 0.0182 | 0.0182 | 0.0182 |
| OCDD                | 660.0         | 0.74        | 0.00030 | 0.1976 | 0.1976 | 0.1976 |

**20 ng/Kg      20 ng/Kg      20 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-26                 |           |                  |  |
| Lab Sample ID          | 10257348012-S             |           |                  |  |
| Filename               | U140226B_18               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.2 g                    | Matrix    | Soil             |  |
| % Moisture             | 25.1                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 7.64 g                    | Collected | 02/04/2014 10:20 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140226B_06 & U140226B_20 | Extracted | 02/24/2014 21:00 |  |
| Method Blank ID        | BLANK-39446               | Analyzed  | 02/27/2014 06:49 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg |     | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|-----|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 12.0       | ----       | 0.25     |     | 2,3,7,8-TCDF-13C         | 2.00       | 82 Y             |
| Total TCDF          | 290.0      | ----       | 0.25     |     | 2,3,7,8-TCDD-13C         | 2.00       | 90               |
|                     |            |            |          |     | 1,2,3,7,8-PeCDF-13C      | 2.00       | 98 Y             |
| 2,3,7,8-TCDD        | -----      | 0.74       | 0.23     | IJY | 2,3,4,7,8-PeCDF-13C      | 2.00       | 89 Y             |
| Total TCDD          | 16.0       | ----       | 0.23     | Y   | 1,2,3,7,8-PeCDD-13C      | 2.00       | 95 Y             |
|                     |            |            |          |     | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 96               |
| 1,2,3,7,8-PeCDF     | 12.0       | ----       | 0.24     |     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 102              |
| 2,3,4,7,8-PeCDF     | 27.0       | ----       | 0.42     |     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 95               |
| Total PeCDF         | 270.0      | ----       | 0.33     |     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 92               |
|                     |            |            |          |     | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 80               |
| 1,2,3,7,8-PeCDD     | -----      | 3.80       | 0.20     | IJ  | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 72               |
| Total PeCDD         | 36.0       | ----       | 0.20     |     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 67               |
|                     |            |            |          |     | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 64               |
| 1,2,3,4,7,8-HxCDF   | 20.0       | ----       | 0.34     |     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 65               |
| 1,2,3,6,7,8-HxCDF   | -----      | 21.00      | 0.27     | P   | OCDD-13C                 | 4.00       | 74               |
| 2,3,4,6,7,8-HxCDF   | 30.0       | ----       | 0.32     |     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 3.6        | ----       | 0.35     | J   | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 290.0      | ----       | 0.32     |     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |     |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 5.6        | ----       | 0.35     | J   | 2,3,7,8-TCDD-37Cl4       | 0.20       | 106              |
| 1,2,3,6,7,8-HxCDD   | 22.0       | ----       | 0.59     |     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 12.0       | ----       | 0.38     |     |                          |            |                  |
| Total HxCDD         | 160.0      | ----       | 0.44     |     |                          |            |                  |
|                     |            |            |          |     |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 140.0      | ----       | 0.30     |     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 7.3        | ----       | 0.46     |     | Equivalence: 33 ng/Kg    |            |                  |
| Total HpCDF         | 330.0      | ----       | 0.38     |     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |     |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 440.0      | ----       | 0.56     |     |                          |            |                  |
| Total HpCDD         | 760.0      | ----       | 0.56     |     |                          |            |                  |
|                     |            |            |          |     |                          |            |                  |
| OCDF                | 250.0      | ----       | 0.36     | Y   |                          |            |                  |
| OCDD                | 5400.0     | ----       | 0.51     |     |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

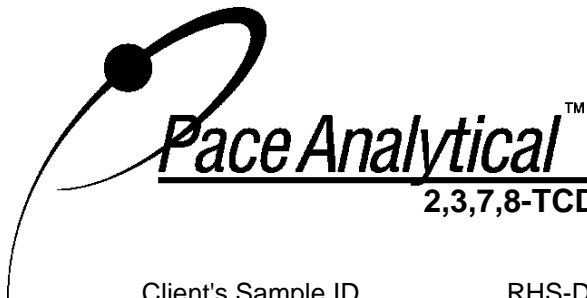
ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

J = Estimated value  
P = PCDE Interference  
I = Interference present  
Y = Calculated using average of daily RFs

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-26                 |           |                  |  |
| Lab Sample ID          | 10257348012-S             |           |                  |  |
| Filename               | U140226B_18               |           |                  |  |
| Injected By            | BAL                       |           |                  |  |
| Total Amount Extracted | 10.2 g                    | Matrix    | Soil             |  |
| % Moisture             | 25.1                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 7.64 g                    | Collected | 02/04/2014 10:20 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140226B_06 & U140226B_20 | Extracted | 02/24/2014 21:00 |  |
| Method Blank ID        | BLANK-39446               | Analyzed  | 02/27/2014 06:49 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | 12.0          | 0.25        | 0.10000 | 1.1923 | 1.1923 | 1.1923 |
| Total TCDF          | 290.0         | 0.25        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | ND            | 0.23        | 1.00000 | 0.7429 | 0.7429 | 0.7429 |
| Total TCDD          | 16.0          | 0.23        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | 12.0          | 0.24        | 0.03000 | 0.3468 | 0.3468 | 0.3468 |
| 2,3,4,7,8-PeCDF     | 27.0          | 0.42        | 0.30000 | 8.1524 | 8.1524 | 8.1524 |
| Total PeCDF         | 270.0         | 0.33        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | ND            | 0.20        | 1.00000 | 3.8311 | 3.8311 | 3.8311 |
| Total PeCDD         | 36.0          | 0.20        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | 20.0          | 0.34        | 0.10000 | 2.0023 | 2.0023 | 2.0023 |
| 1,2,3,6,7,8-HxCDF   | ND            | 0.27        | 0.10000 | 2.1271 | 2.1271 | 2.1271 |
| 2,3,4,6,7,8-HxCDF   | 30.0          | 0.32        | 0.10000 | 2.9534 | 2.9534 | 2.9534 |
| 1,2,3,7,8,9-HxCDF   | 3.6           | 0.35        | 0.10000 | 0.3650 | 0.3650 | 0.3650 |
| Total HxCDF         | 290.0         | 0.32        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | 5.6           | 0.35        | 0.10000 | 0.5628 | 0.5628 | 0.5628 |
| 1,2,3,6,7,8-HxCDD   | 22.0          | 0.59        | 0.10000 | 2.2083 | 2.2083 | 2.2083 |
| 1,2,3,7,8,9-HxCDD   | 12.0          | 0.38        | 0.10000 | 1.1817 | 1.1817 | 1.1817 |
| Total HxCDD         | 160.0         | 0.44        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | 140.0         | 0.30        | 0.01000 | 1.3929 | 1.3929 | 1.3929 |
| 1,2,3,4,7,8,9-HpCDF | 7.3           | 0.46        | 0.01000 | 0.0729 | 0.0729 | 0.0729 |
| Total HpCDF         | 330.0         | 0.38        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | 440.0         | 0.56        | 0.01000 | 4.4479 | 4.4479 | 4.4479 |
| Total HpCDD         | 760.0         | 0.56        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | 250.0         | 0.36        | 0.00030 | 0.0756 | 0.0756 | 0.0756 |
| OCDD                | 5400.0        | 0.51        | 0.00030 | 1.6071 | 1.6071 | 1.6071 |

**33 ng/Kg      33 ng/Kg      33 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-26.2               |           |                  |  |
| Lab Sample ID          | 10257348013-S             |           |                  |  |
| Filename               | U140304B_12               |           |                  |  |
| Injected By            | SMT                       |           |                  |  |
| Total Amount Extracted | 10.9 g                    | Matrix    | Soil             |  |
| % Moisture             | 18.2                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 8.92 g                    | Collected | 02/04/2014 10:59 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |  |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/05/2014 00:53 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 8.60       | ----       | 0.39     | 2,3,7,8-TCDF-13C         | 2.00       | 98               |
| Total TCDF          | 230.00     | ----       | 0.39     | 2,3,7,8-TCDD-13C         | 2.00       | 102              |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 109              |
| 2,3,7,8-TCDD        | 0.41       | ----       | 0.32 J   | 2,3,4,7,8-PeCDF-13C      | 2.00       | 85 Y             |
| Total TCDD          | 13.00      | ----       | 0.32     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 114              |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 93               |
| 1,2,3,7,8-PeCDF     | 9.40       | ----       | 0.24     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 94               |
| 2,3,4,7,8-PeCDF     | 21.00      | ----       | 0.21     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 101              |
| Total PeCDF         | 260.00     | ----       | 0.23     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 94               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 83               |
| 1,2,3,7,8-PeCDD     | 2.50       | ----       | 0.36 J   | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 79               |
| Total PeCDD         | 24.00      | ----       | 0.36     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 78               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 82               |
| 1,2,3,4,7,8-HxCDF   | 15.00      | ----       | 0.40     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 90               |
| 1,2,3,6,7,8-HxCDF   | 16.00      | ----       | 0.55     | OCDD-13C                 | 4.00       | 72               |
| 2,3,4,6,7,8-HxCDF   | 19.00      | ----       | 0.41     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 3.20       | ----       | 0.51 J   | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 190.00     | ----       | 0.47     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 2.70       | ----       | 0.38 J   | 2,3,7,8-TCDD-37Cl4       | 0.20       | 100              |
| 1,2,3,6,7,8-HxCDD   | 6.70       | ----       | 0.41     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 5.50       | ----       | 0.43 J   |                          |            |                  |
| Total HxCDD         | 69.00      | ----       | 0.40     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 70.00      | ----       | 0.31     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 3.30       | ----       | 0.34 J   | Equivalence: 19 ng/Kg    |            |                  |
| Total HpCDF         | 110.00     | ----       | 0.33     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 98.00      | ----       | 0.37     |                          |            |                  |
| Total HpCDD         | 180.00     | ----       | 0.37     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 59.00      | ----       | 0.82     |                          |            |                  |
| OCDD                | 640.00     | ----       | 0.76     |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

J = Estimated value

Y = Calculated using average of daily RFs

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |
|------------------------|---------------------------|-----------|------------------|
| Client's Sample ID     | RHS-DU-26.2               |           |                  |
| Lab Sample ID          | 10257348013-S             |           |                  |
| Filename               | U140304B_12               |           |                  |
| Injected By            | SMT                       |           |                  |
| Total Amount Extracted | 10.9 g                    | Matrix    | Soil             |
| % Moisture             | 18.2                      | Dilution  | NA               |
| Dry Weight Extracted   | 8.92 g                    | Collected | 02/04/2014 10:59 |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/05/2014 00:53 |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | 8.60          | 0.39        | 0.10000 | 0.8649 | 0.8649 | 0.8649 |
| Total TCDF          | 230.00        | 0.39        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | 0.41          | 0.32        | 1.00000 | 0.4072 | 0.4072 | 0.4072 |
| Total TCDD          | 13.00         | 0.32        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | 9.40          | 0.24        | 0.03000 | 0.2813 | 0.2813 | 0.2813 |
| 2,3,4,7,8-PeCDF     | 21.00         | 0.21        | 0.30000 | 6.2647 | 6.2647 | 6.2647 |
| Total PeCDF         | 260.00        | 0.23        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | 2.50          | 0.36        | 1.00000 | 2.5449 | 2.5449 | 2.5449 |
| Total PeCDD         | 24.00         | 0.36        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | 15.00         | 0.40        | 0.10000 | 1.5027 | 1.5027 | 1.5027 |
| 1,2,3,6,7,8-HxCDF   | 16.00         | 0.55        | 0.10000 | 1.6224 | 1.6224 | 1.6224 |
| 2,3,4,6,7,8-HxCDF   | 19.00         | 0.41        | 0.10000 | 1.9275 | 1.9275 | 1.9275 |
| 1,2,3,7,8,9-HxCDF   | 3.20          | 0.51        | 0.10000 | 0.3179 | 0.3179 | 0.3179 |
| Total HxCDF         | 190.00        | 0.47        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | 2.70          | 0.38        | 0.10000 | 0.2750 | 0.2750 | 0.2750 |
| 1,2,3,6,7,8-HxCDD   | 6.70          | 0.41        | 0.10000 | 0.6667 | 0.6667 | 0.6667 |
| 1,2,3,7,8,9-HxCDD   | 5.50          | 0.43        | 0.10000 | 0.5466 | 0.5466 | 0.5466 |
| Total HxCDD         | 69.00         | 0.40        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | 70.00         | 0.31        | 0.01000 | 0.6981 | 0.6981 | 0.6981 |
| 1,2,3,4,7,8,9-HpCDF | 3.30          | 0.34        | 0.01000 | 0.0326 | 0.0326 | 0.0326 |
| Total HpCDF         | 110.00        | 0.33        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | 98.00         | 0.37        | 0.01000 | 0.9794 | 0.9794 | 0.9794 |
| Total HpCDD         | 180.00        | 0.37        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | 59.00         | 0.82        | 0.00030 | 0.0176 | 0.0176 | 0.0176 |
| OCDD                | 640.00        | 0.76        | 0.00030 | 0.1926 | 0.1926 | 0.1926 |

**19 ng/Kg      19 ng/Kg      19 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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### Method 8290 Sample Analysis Results

Client - ESN Pacific

|                        |                           |           |                  |  |  |
|------------------------|---------------------------|-----------|------------------|--|--|
| Client's Sample ID     | RHS-DU-26.3               |           |                  |  |  |
| Lab Sample ID          | 10257348014-S             |           |                  |  |  |
| Filename               | U140304B_13               |           |                  |  |  |
| Injected By            | SMT                       |           |                  |  |  |
| Total Amount Extracted | 10.5 g                    | Matrix    | Soil             |  |  |
| % Moisture             | 18.7                      | Dilution  | NA               |  |  |
| Dry Weight Extracted   | 8.54 g                    | Collected | 02/04/2014 11:45 |  |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |  |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |  |  |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/05/2014 01:38 |  |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg |    | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|----|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 9.1        | ----       | 0.38     |    | 2,3,7,8-TCDF-13C         | 2.00       | 94               |
| Total TCDF          | 230.0      | ----       | 0.38     |    | 2,3,7,8-TCDD-13C         | 2.00       | 99               |
|                     |            |            |          |    | 1,2,3,7,8-PeCDF-13C      | 2.00       | 108              |
| 2,3,7,8-TCDD        | -----      | 0.43       | 0.33     | IJ | 2,3,4,7,8-PeCDF-13C      | 2.00       | 82 Y             |
| Total TCDD          | 11.0       | ----       | 0.33     |    | 1,2,3,7,8-PeCDD-13C      | 2.00       | 110              |
|                     |            |            |          |    | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 95               |
| 1,2,3,7,8-PeCDF     | 9.7        | ----       | 0.31     |    | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 96               |
| 2,3,4,7,8-PeCDF     | 22.0       | ----       | 0.29     |    | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 104              |
| Total PeCDF         | 270.0      | ----       | 0.30     |    | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 95               |
|                     |            |            |          |    | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 85               |
| 1,2,3,7,8-PeCDD     | -----      | 2.20       | 0.24     | IJ | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 79               |
| Total PeCDD         | 22.0       | ----       | 0.24     |    | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 79               |
|                     |            |            |          |    | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 81               |
| 1,2,3,4,7,8-HxCDF   | 15.0       | ----       | 0.53     |    | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 92               |
| 1,2,3,6,7,8-HxCDF   | 17.0       | ----       | 0.45     |    | OCDD-13C                 | 4.00       | 74               |
| 2,3,4,6,7,8-HxCDF   | 21.0       | ----       | 0.33     |    |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 3.5        | ----       | 0.46     | J  | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 210.0      | ----       | 0.44     |    | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 3.0        | ----       | 0.85     | J  | 2,3,7,8-TCDD-37Cl4       | 0.20       | 98               |
| 1,2,3,6,7,8-HxCDD   | 7.0        | ----       | 0.28     |    |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 5.4        | ----       | 0.79     | J  |                          |            |                  |
| Total HxCDD         | 70.0       | ----       | 0.64     |    |                          |            |                  |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 73.0       | ----       | 0.30     |    | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 3.4        | ----       | 0.49     | J  | Equivalence: 20 ng/Kg    |            |                  |
| Total HpCDF         | 120.0      | ----       | 0.40     |    | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 100.0      | ----       | 0.27     |    |                          |            |                  |
| Total HpCDD         | 190.0      | ----       | 0.27     |    |                          |            |                  |
|                     |            |            |          |    |                          |            |                  |
| OCDF                | 59.0       | ----       | 0.75     |    |                          |            |                  |
| OCDD                | 670.0      | ----       | 0.69     |    |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

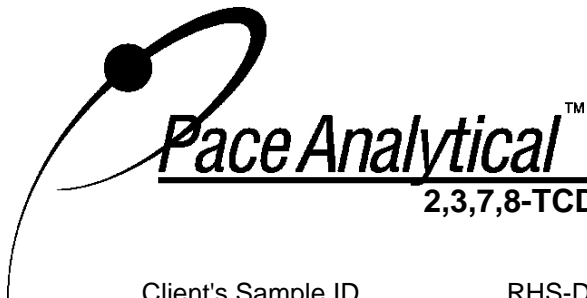
ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

J = Estimated value  
I = Interference present  
Y = Calculated using average of daily RFs

## REPORT OF LABORATORY ANALYSIS

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |
|------------------------|---------------------------|-----------|------------------|
| Client's Sample ID     | RHS-DU-26.3               |           |                  |
| Lab Sample ID          | 10257348014-S             |           |                  |
| Filename               | U140304B_13               |           |                  |
| Injected By            | SMT                       |           |                  |
| Total Amount Extracted | 10.5 g                    | Matrix    | Soil             |
| % Moisture             | 18.7                      | Dilution  | NA               |
| Dry Weight Extracted   | 8.54 g                    | Collected | 02/04/2014 11:45 |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/05/2014 01:38 |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | 9.1           | 0.38        | 0.10000 | 0.9149 | 0.9149 | 0.9149 |
| Total TCDF          | 230.0         | 0.38        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | ND            | 0.33        | 1.00000 | 0.4277 | 0.4277 | 0.4277 |
| Total TCDD          | 11.0          | 0.33        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | 9.7           | 0.31        | 0.03000 | 0.2911 | 0.2911 | 0.2911 |
| 2,3,4,7,8-PeCDF     | 22.0          | 0.29        | 0.30000 | 6.5593 | 6.5593 | 6.5593 |
| Total PeCDF         | 270.0         | 0.30        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | ND            | 0.24        | 1.00000 | 2.2206 | 2.2206 | 2.2206 |
| Total PeCDD         | 22.0          | 0.24        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | 15.0          | 0.53        | 0.10000 | 1.4629 | 1.4629 | 1.4629 |
| 1,2,3,6,7,8-HxCDF   | 17.0          | 0.45        | 0.10000 | 1.7089 | 1.7089 | 1.7089 |
| 2,3,4,6,7,8-HxCDF   | 21.0          | 0.33        | 0.10000 | 2.0703 | 2.0703 | 2.0703 |
| 1,2,3,7,8,9-HxCDF   | 3.5           | 0.46        | 0.10000 | 0.3540 | 0.3540 | 0.3540 |
| Total HxCDF         | 210.0         | 0.44        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | 3.0           | 0.85        | 0.10000 | 0.2999 | 0.2999 | 0.2999 |
| 1,2,3,6,7,8-HxCDD   | 7.0           | 0.28        | 0.10000 | 0.7035 | 0.7035 | 0.7035 |
| 1,2,3,7,8,9-HxCDD   | 5.4           | 0.79        | 0.10000 | 0.5370 | 0.5370 | 0.5370 |
| Total HxCDD         | 70.0          | 0.64        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | 73.0          | 0.30        | 0.01000 | 0.7261 | 0.7261 | 0.7261 |
| 1,2,3,4,7,8,9-HpCDF | 3.4           | 0.49        | 0.01000 | 0.0337 | 0.0337 | 0.0337 |
| Total HpCDF         | 120.0         | 0.40        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | 100.0         | 0.27        | 0.01000 | 1.0233 | 1.0233 | 1.0233 |
| Total HpCDD         | 190.0         | 0.27        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | 59.0          | 0.75        | 0.00030 | 0.0177 | 0.0177 | 0.0177 |
| OCDD                | 670.0         | 0.69        | 0.00030 | 0.1996 | 0.1996 | 0.1996 |

**20 ng/Kg      20 ng/Kg      20 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

**REPORT OF LABORATORY ANALYSIS**

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |
|------------------------|---------------------------|-----------|------------------|
| Client's Sample ID     | RHS-DU-27                 |           |                  |
| Lab Sample ID          | 10257348015-S             |           |                  |
| Filename               | U140304B_14               |           |                  |
| Injected By            | SMT                       |           |                  |
| Total Amount Extracted | 10.0 g                    | Matrix    | Soil             |
| % Moisture             | 21.7                      | Dilution  | NA               |
| Dry Weight Extracted   | 7.83 g                    | Collected | 02/04/2014 10:00 |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/05/2014 02:23 |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 11.00      | ----       | 0.46     | 2,3,7,8-TCDF-13C         | 2.00       | 95               |
| Total TCDF          | 270.00     | ----       | 0.46     | 2,3,7,8-TCDD-13C         | 2.00       | 100              |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 110              |
| 2,3,7,8-TCDD        | 0.67       | ----       | 0.39 J   | 2,3,4,7,8-PeCDF-13C      | 2.00       | 86 Y             |
| Total TCDD          | 14.00      | ----       | 0.39     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 116              |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 93               |
| 1,2,3,7,8-PeCDF     | 11.00      | ----       | 0.24     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 93               |
| 2,3,4,7,8-PeCDF     | 24.00      | ----       | 0.39     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 104              |
| Total PeCDF         | 280.00     | ----       | 0.32     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 94               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 86               |
| 1,2,3,7,8-PeCDD     | -----      | 2.4        | 0.21 IJ  | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 76               |
| Total PeCDD         | 22.00      | -----      | 0.21     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 80               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 81               |
| 1,2,3,4,7,8-HxCDF   | 17.00      | ----       | 0.57     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 91               |
| 1,2,3,6,7,8-HxCDF   | 20.00      | ----       | 0.61     | OCDD-13C                 | 4.00       | 74               |
| 2,3,4,6,7,8-HxCDF   | 21.00      | ----       | 0.34     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 4.00       | ----       | 0.33 J   | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 220.00     | ----       | 0.46     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 2.80       | ----       | 0.36 J   | 2,3,7,8-TCDD-37Cl4       | 0.20       | 100              |
| 1,2,3,6,7,8-HxCDD   | -----      | 7.5        | 0.42 I   |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 5.80       | ----       | 0.31 J   |                          |            |                  |
| Total HxCDD         | 70.00      | ----       | 0.37     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 80.00      | ----       | 0.22     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 4.10       | ----       | 0.50 J   | Equivalence: 22 ng/Kg    |            |                  |
| Total HpCDF         | 120.00     | ----       | 0.36     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 110.00     | ----       | 0.56     |                          |            |                  |
| Total HpCDD         | 200.00     | ----       | 0.56     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 67.00      | ----       | 0.70     |                          |            |                  |
| OCDD                | 720.00     | ----       | 0.77     |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.  
J = Estimated value  
I = Interference present  
Y = Calculated using average of daily RFs

**REPORT OF LABORATORY ANALYSIS**

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |
|------------------------|---------------------------|-----------|------------------|
| Client's Sample ID     | RHS-DU-27                 |           |                  |
| Lab Sample ID          | 10257348015-S             |           |                  |
| Filename               | U140304B_14               |           |                  |
| Injected By            | SMT                       |           |                  |
| Total Amount Extracted | 10.0 g                    | Matrix    | Soil             |
| % Moisture             | 21.7                      | Dilution  | NA               |
| Dry Weight Extracted   | 7.83 g                    | Collected | 02/04/2014 10:00 |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/05/2014 02:23 |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | 11.00         | 0.46        | 0.10000 | 1.0538 | 1.0538 | 1.0538 |
| Total TCDF          | 270.00        | 0.46        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | 0.67          | 0.39        | 1.00000 | 0.6667 | 0.6667 | 0.6667 |
| Total TCDD          | 14.00         | 0.39        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | 11.00         | 0.24        | 0.03000 | 0.3194 | 0.3194 | 0.3194 |
| 2,3,4,7,8-PeCDF     | 24.00         | 0.39        | 0.30000 | 7.1498 | 7.1498 | 7.1498 |
| Total PeCDF         | 280.00        | 0.32        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | ND            | 0.21        | 1.00000 | 2.4427 | 2.4427 | 2.4427 |
| Total PeCDD         | 22.00         | 0.21        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | 17.00         | 0.57        | 0.10000 | 1.7217 | 1.7217 | 1.7217 |
| 1,2,3,6,7,8-HxCDF   | 20.00         | 0.61        | 0.10000 | 1.9854 | 1.9854 | 1.9854 |
| 2,3,4,6,7,8-HxCDF   | 21.00         | 0.34        | 0.10000 | 2.1190 | 2.1190 | 2.1190 |
| 1,2,3,7,8,9-HxCDF   | 4.00          | 0.33        | 0.10000 | 0.4026 | 0.4026 | 0.4026 |
| Total HxCDF         | 220.00        | 0.46        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | 2.80          | 0.36        | 0.10000 | 0.2841 | 0.2841 | 0.2841 |
| 1,2,3,6,7,8-HxCDD   | ND            | 0.42        | 0.10000 | 0.7527 | 0.7527 | 0.7527 |
| 1,2,3,7,8,9-HxCDD   | 5.80          | 0.31        | 0.10000 | 0.5810 | 0.5810 | 0.5810 |
| Total HxCDD         | 70.00         | 0.37        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | 80.00         | 0.22        | 0.01000 | 0.7958 | 0.7958 | 0.7958 |
| 1,2,3,4,7,8,9-HpCDF | 4.10          | 0.50        | 0.01000 | 0.0406 | 0.0406 | 0.0406 |
| Total HpCDF         | 120.00        | 0.36        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | 110.00        | 0.56        | 0.01000 | 1.1320 | 1.1320 | 1.1320 |
| Total HpCDD         | 200.00        | 0.56        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | 67.00         | 0.70        | 0.00030 | 0.0200 | 0.0200 | 0.0200 |
| OCDD                | 720.00        | 0.77        | 0.00030 | 0.2153 | 0.2153 | 0.2153 |

**22 ng/Kg      22 ng/Kg      22 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

**REPORT OF LABORATORY ANALYSIS**

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-28                 |           |                  |  |
| Lab Sample ID          | 10257348016-S             |           |                  |  |
| Filename               | U140304B_15               |           |                  |  |
| Injected By            | SMT                       |           |                  |  |
| Total Amount Extracted | 6.40 g                    | Matrix    | Soil             |  |
| % Moisture             | 18.9                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 5.19 g                    | Collected | 02/04/2014 10:07 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |  |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/05/2014 03:08 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 3.8        | ----       | 0.65     | 2,3,7,8-TCDF-13C         | 2.00       | 92               |
| Total TCDF          | 79.0       | ----       | 0.65     | 2,3,7,8-TCDD-13C         | 2.00       | 96               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 108              |
| 2,3,7,8-TCDD        | ND         | ----       | 0.52     | 2,3,4,7,8-PeCDF-13C      | 2.00       | 84 Y             |
| Total TCDD          | 3.3        | ----       | 0.52     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 113              |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 92               |
| 1,2,3,7,8-PeCDF     | 3.3        | ----       | 0.45 J   | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 101              |
| 2,3,4,7,8-PeCDF     | 8.0        | ----       | 0.40 J   | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 103              |
| Total PeCDF         | 93.0       | ----       | 0.43     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 95               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 86               |
| 1,2,3,7,8-PeCDD     | ----       | 1.2        | 0.27 IJ  | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 78               |
| Total PeCDD         | 2.9        | ----       | 0.27 J   | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 86               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 88               |
| 1,2,3,4,7,8-HxCDF   | 6.2        | ----       | 0.57 J   | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 97               |
| 1,2,3,6,7,8-HxCDF   | 8.7        | ----       | 0.75 J   | OCDD-13C                 | 4.00       | 91               |
| 2,3,4,6,7,8-HxCDF   | 10.0       | ----       | 0.74     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 1.4        | ----       | 0.69 J   | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 150.0      | ----       | 0.69     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 3.2        | ----       | 0.57 J   | 2,3,7,8-TCDD-37Cl4       | 0.20       | 93               |
| 1,2,3,6,7,8-HxCDD   | 12.0       | ----       | 0.76     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 6.7        | ----       | 0.69 J   |                          |            |                  |
| Total HxCDD         | 74.0       | ----       | 0.67     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 84.0       | ----       | 0.46     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 3.4        | ----       | 1.00 J   | Equivalence: 14 ng/Kg    |            |                  |
| Total HpCDF         | 220.0      | ----       | 0.73     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 300.0      | ----       | 0.59     |                          |            |                  |
| Total HpCDD         | 520.0      | ----       | 0.59     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 200.0      | ----       | 0.91     |                          |            |                  |
| OCDD                | 3600.0     | ----       | 0.81     |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

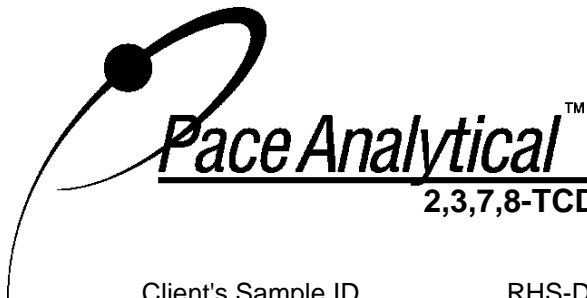
J = Estimated value

I = Interference present

Y = Calculated using average of daily RFs

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |  |
|------------------------|---------------------------|-----------|------------------|--|--|
| Client's Sample ID     | RHS-DU-28                 |           |                  |  |  |
| Lab Sample ID          | 10257348016-S             |           |                  |  |  |
| Filename               | U140304B_15               |           |                  |  |  |
| Injected By            | SMT                       |           |                  |  |  |
| Total Amount Extracted | 6.40 g                    | Matrix    | Soil             |  |  |
| % Moisture             | 18.9                      | Dilution  | NA               |  |  |
| Dry Weight Extracted   | 5.19 g                    | Collected | 02/04/2014 10:07 |  |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |  |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |  |  |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/05/2014 03:08 |  |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | 3.8           | 0.65        | 0.10000 | 0.3763 | 0.3763 | 0.3763 |
| Total TCDF          | 79.0          | 0.65        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | ND            | 0.52        | 1.00000 | 0.0000 | 0.2592 | 0.5184 |
| Total TCDD          | 3.3           | 0.52        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | 3.3           | 0.45        | 0.03000 | 0.0992 | 0.0992 | 0.0992 |
| 2,3,4,7,8-PeCDF     | 8.0           | 0.40        | 0.30000 | 2.4084 | 2.4084 | 2.4084 |
| Total PeCDF         | 93.0          | 0.43        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | ND            | 0.27        | 1.00000 | 1.2273 | 1.2273 | 1.2273 |
| Total PeCDD         | 2.9           | 0.27        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | 6.2           | 0.57        | 0.10000 | 0.6152 | 0.6152 | 0.6152 |
| 1,2,3,6,7,8-HxCDF   | 8.7           | 0.75        | 0.10000 | 0.8650 | 0.8650 | 0.8650 |
| 2,3,4,6,7,8-HxCDF   | 10.0          | 0.74        | 0.10000 | 1.0415 | 1.0415 | 1.0415 |
| 1,2,3,7,8,9-HxCDF   | 1.4           | 0.69        | 0.10000 | 0.1411 | 0.1411 | 0.1411 |
| Total HxCDF         | 150.0         | 0.69        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | 3.2           | 0.57        | 0.10000 | 0.3218 | 0.3218 | 0.3218 |
| 1,2,3,6,7,8-HxCDD   | 12.0          | 0.76        | 0.10000 | 1.2422 | 1.2422 | 1.2422 |
| 1,2,3,7,8,9-HxCDD   | 6.7           | 0.69        | 0.10000 | 0.6679 | 0.6679 | 0.6679 |
| Total HxCDD         | 74.0          | 0.67        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | 84.0          | 0.46        | 0.01000 | 0.8394 | 0.8394 | 0.8394 |
| 1,2,3,4,7,8,9-HpCDF | 3.4           | 1.00        | 0.01000 | 0.0343 | 0.0343 | 0.0343 |
| Total HpCDF         | 220.0         | 0.73        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | 300.0         | 0.59        | 0.01000 | 3.0432 | 3.0432 | 3.0432 |
| Total HpCDD         | 520.0         | 0.59        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | 200.0         | 0.91        | 0.00030 | 0.0611 | 0.0611 | 0.0611 |
| OCDD                | 3600.0        | 0.81        | 0.00030 | 1.0855 | 1.0855 | 1.0855 |

**14 ng/Kg      14 ng/Kg      15 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Sample Analysis Results**

Client - ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-29                 |           |                  |  |
| Lab Sample ID          | 10257348017-S             |           |                  |  |
| Filename               | U140304B_16               |           |                  |  |
| Injected By            | SMT                       |           |                  |  |
| Total Amount Extracted | 10.7 g                    | Matrix    | Soil             |  |
| % Moisture             | 16.1                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 8.98 g                    | Collected | 02/04/2014 09:45 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |  |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/05/2014 03:52 |  |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 4.50       | ----       | 0.32     | 2,3,7,8-TCDF-13C         | 2.00       | 104              |
| Total TCDF          | 100.00     | ----       | 0.32     | 2,3,7,8-TCDD-13C         | 2.00       | 109              |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 127              |
| 2,3,7,8-TCDD        | 0.48       | ----       | 0.25 J   | 2,3,4,7,8-PeCDF-13C      | 2.00       | 96 Y             |
| Total TCDD          | 6.70       | ----       | 0.25     | 1,2,3,7,8-PeCDD-13C      | 2.00       | 134              |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 100              |
| 1,2,3,7,8-PeCDF     | 5.80       | ----       | 0.21     | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 107              |
| 2,3,4,7,8-PeCDF     | 13.00      | ----       | 0.16     | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 113              |
| Total PeCDF         | 140.00     | ----       | 0.18     | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 104              |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 91               |
| 1,2,3,7,8-PeCDD     | 2.50       | ----       | 0.12 J   | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 87               |
| Total PeCDD         | 22.00      | ----       | 0.12     | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 95               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 99               |
| 1,2,3,4,7,8-HxCDF   | 12.00      | ----       | 0.53     | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 108              |
| 1,2,3,6,7,8-HxCDF   | 14.00      | ----       | 0.23     | OCDD-13C                 | 4.00       | 108              |
| 2,3,4,6,7,8-HxCDF   | 18.00      | ----       | 0.21     |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 2.70       | ----       | 0.26 J   | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 200.00     | ----       | 0.31     | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 4.40       | ----       | 0.28 J   | 2,3,7,8-TCDD-37Cl4       | 0.20       | 108              |
| 1,2,3,6,7,8-HxCDD   | 17.00      | ----       | 0.30     |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | 8.50       | ----       | 0.21     |                          |            |                  |
| Total HxCDD         | 110.00     | ----       | 0.26     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 120.00     | ----       | 0.16     | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 5.20       | ----       | 0.31 J   | Equivalence: 21 ng/Kg    |            |                  |
| Total HpCDF         | 290.00     | ----       | 0.23     | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 360.00     | ----       | 0.40     |                          |            |                  |
| Total HpCDD         | 610.00     | ----       | 0.40     |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | 240.00     | ----       | 0.31     |                          |            |                  |
| OCDD                | 4200.00    | ----       | 0.34     |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit.

ND = Not Detected  
NA = Not Applicable  
NC = Not Calculated

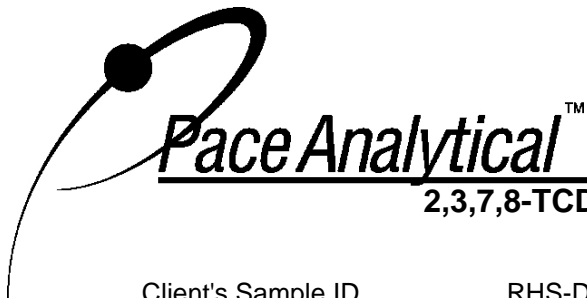
Results reported on a dry weight basis and are valid to no more than 2 significant figures.

J = Estimated value

Y = Calculated using average of daily RfFs

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |
|------------------------|---------------------------|-----------|------------------|--|
| Client's Sample ID     | RHS-DU-29                 |           |                  |  |
| Lab Sample ID          | 10257348017-S             |           |                  |  |
| Filename               | U140304B_16               |           |                  |  |
| Injected By            | SMT                       |           |                  |  |
| Total Amount Extracted | 10.7 g                    | Matrix    | Soil             |  |
| % Moisture             | 16.1                      | Dilution  | NA               |  |
| Dry Weight Extracted   | 8.98 g                    | Collected | 02/04/2014 09:45 |  |
| ICAL ID                | U140224                   | Received  | 02/11/2014 09:07 |  |
| CCal Filename(s)       | U140304B_02 & U140304B_20 | Extracted | 02/27/2014 19:30 |  |
| Method Blank ID        | BLANK-39477               | Analyzed  | 03/05/2014 03:52 |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB              | MB              | UB              |
|---------------------|---------------|-------------|---------|-----------------|-----------------|-----------------|
| 2,3,7,8-TCDF        | 4.50          | 0.32        | 0.10000 | 0.4526          | 0.4526          | 0.4526          |
| Total TCDF          | 100.00        | 0.32        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 2,3,7,8-TCDD        | 0.48          | 0.25        | 1.00000 | 0.4751          | 0.4751          | 0.4751          |
| Total TCDD          | 6.70          | 0.25        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,7,8-PeCDF     | 5.80          | 0.21        | 0.03000 | 0.1752          | 0.1752          | 0.1752          |
| 2,3,4,7,8-PeCDF     | 13.00         | 0.16        | 0.30000 | 4.0438          | 4.0438          | 4.0438          |
| Total PeCDF         | 140.00        | 0.18        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,7,8-PeCDD     | 2.50          | 0.12        | 1.00000 | 2.4535          | 2.4535          | 2.4535          |
| Total PeCDD         | 22.00         | 0.12        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,4,7,8-HxCDF   | 12.00         | 0.53        | 0.10000 | 1.1988          | 1.1988          | 1.1988          |
| 1,2,3,6,7,8-HxCDF   | 14.00         | 0.23        | 0.10000 | 1.4133          | 1.4133          | 1.4133          |
| 2,3,4,6,7,8-HxCDF   | 18.00         | 0.21        | 0.10000 | 1.7704          | 1.7704          | 1.7704          |
| 1,2,3,7,8,9-HxCDF   | 2.70          | 0.26        | 0.10000 | 0.2678          | 0.2678          | 0.2678          |
| Total HxCDF         | 200.00        | 0.31        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,4,7,8-HxCDD   | 4.40          | 0.28        | 0.10000 | 0.4447          | 0.4447          | 0.4447          |
| 1,2,3,6,7,8-HxCDD   | 17.00         | 0.30        | 0.10000 | 1.6783          | 1.6783          | 1.6783          |
| 1,2,3,7,8,9-HxCDD   | 8.50          | 0.21        | 0.10000 | 0.8481          | 0.8481          | 0.8481          |
| Total HxCDD         | 110.00        | 0.26        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,4,6,7,8-HpCDF | 120.00        | 0.16        | 0.01000 | 1.1781          | 1.1781          | 1.1781          |
| 1,2,3,4,7,8,9-HpCDF | 5.20          | 0.31        | 0.01000 | 0.0522          | 0.0522          | 0.0522          |
| Total HpCDF         | 290.00        | 0.23        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| 1,2,3,4,6,7,8-HpCDD | 360.00        | 0.40        | 0.01000 | 3.6495          | 3.6495          | 3.6495          |
| Total HpCDD         | 610.00        | 0.40        | 0.00000 | 0.0000          | 0.0000          | 0.0000          |
| OCDF                | 240.00        | 0.31        | 0.00030 | 0.0724          | 0.0724          | 0.0724          |
| OCDD                | 4200.00       | 0.34        | 0.00030 | 1.2711          | 1.2711          | 1.2711          |
|                     |               |             |         | <b>21 ng/Kg</b> | <b>21 ng/Kg</b> | <b>21 ng/Kg</b> |

Final values are valid to only 2 significant figures  
 TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
 LB = Lower Bound, Where "ND", TEQ Conc = 0  
 MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
 UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
 RL = Reporting Limit

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**Method 8290 Blank Analysis Results**

|                        |                           |             |                  |
|------------------------|---------------------------|-------------|------------------|
| Lab Sample ID          | BLANK-39446               | Matrix      | Solid            |
| Filename               | F140226B_09               | Dilution    | NA               |
| Total Amount Extracted | 10.5 g                    | Extracted   | 02/24/2014 21:00 |
| ICAL ID                | F131125                   | Analyzed    | 02/27/2014 02:08 |
| CCal Filename(s)       | F140226B_01 & F140226B_17 | Injected By | BAL              |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg |    | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|----|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | -----      | 0.140      | 0.035    | IJ | 2,3,7,8-TCDF-13C         | 2.00       | 74               |
| Total TCDF          | 0.430      | -----      | 0.035    | J  | 2,3,7,8-TCDD-13C         | 2.00       | 81               |
|                     |            |            |          |    | 1,2,3,7,8-PeCDF-13C      | 2.00       | 81               |
| 2,3,7,8-TCDD        | -----      | 0.060      | 0.057    | IJ | 2,3,4,7,8-PeCDF-13C      | 2.00       | 82               |
| Total TCDD          | ND         | -----      | 0.057    |    | 1,2,3,7,8-PeCDD-13C      | 2.00       | 88               |
|                     |            |            |          |    | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 78               |
| 1,2,3,7,8-PeCDF     | -----      | 0.150      | 0.062    | IJ | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 91               |
| 2,3,4,7,8-PeCDF     | ND         | -----      | 0.037    |    | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 88               |
| Total PeCDF         | ND         | -----      | 0.049    |    | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 82               |
|                     |            |            |          |    | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 73               |
| 1,2,3,7,8-PeCDD     | -----      | 0.086      | 0.039    | IJ | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 77               |
| Total PeCDD         | ND         | -----      | 0.039    |    | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 74               |
|                     |            |            |          |    | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 73               |
| 1,2,3,4,7,8-HxCDF   | ND         | -----      | 0.048    |    | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 79               |
| 1,2,3,6,7,8-HxCDF   | 0.094      | -----      | 0.035    | J  | OCDD-13C                 | 4.00       | 70               |
| 2,3,4,6,7,8-HxCDF   | -----      | 0.028      | 0.028    | IJ |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | ND         | -----      | 0.042    |    | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 0.094      | -----      | 0.038    | J  | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | ND         | -----      | 0.050    |    | 2,3,7,8-TCDD-37Cl4       | 0.20       | 87               |
| 1,2,3,6,7,8-HxCDD   | -----      | 0.081      | 0.052    | IJ |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | ND         | -----      | 0.048    |    |                          |            |                  |
| Total HxCDD         | ND         | -----      | 0.050    |    |                          |            |                  |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | -----      | 0.096      | 0.056    | IJ | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | ND         | -----      | 0.047    |    | Equivalence: 0.19 ng/Kg  |            |                  |
| Total HpCDF         | ND         | -----      | 0.051    |    | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | -----      | 0.068      | 0.054    | IJ |                          |            |                  |
| Total HpCDD         | 0.100      | -----      | 0.054    | J  |                          |            |                  |
|                     |            |            |          |    |                          |            |                  |
| OCDF                | 0.200      | -----      | 0.120    | J  |                          |            |                  |
| OCDD                | -----      | 0.370      | 0.150    | IJ |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

J = Estimated value

I = Interference present

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |  |
|------------------------|---------------------------|-----------|------------------|--|--|
| Client's Sample ID     |                           |           |                  |  |  |
| Lab Sample ID          | BLANK-39446               |           |                  |  |  |
| Filename               | F140226B_09               |           |                  |  |  |
| Injected By            | BAL                       |           |                  |  |  |
| Total Amount Extracted | 10.5 g                    | Matrix    | Solid            |  |  |
| % Moisture             | 0.0                       | Dilution  | NA               |  |  |
| Dry Weight Extracted   | 10.5 g                    | Collected | 02/24/2014 14:08 |  |  |
| ICAL ID                | F131125                   | Received  | 02/24/2014 14:08 |  |  |
| CCal Filename(s)       | F140226B_01 & F140226B_17 | Extracted | 02/24/2014 21:00 |  |  |
| Method Blank ID        |                           | Analyzed  | 02/27/2014 02:08 |  |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | ND            | 0.035       | 0.10000 | 0.0136 | 0.0136 | 0.0136 |
| Total TCDF          | 0.430         | 0.035       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | ND            | 0.057       | 1.00000 | 0.0601 | 0.0601 | 0.0601 |
| Total TCDD          | ND            | 0.057       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | ND            | 0.062       | 0.03000 | 0.0044 | 0.0044 | 0.0044 |
| 2,3,4,7,8-PeCDF     | ND            | 0.037       | 0.30000 | 0.0000 | 0.0056 | 0.0112 |
| Total PeCDF         | ND            | 0.049       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | ND            | 0.039       | 1.00000 | 0.0857 | 0.0857 | 0.0857 |
| Total PeCDD         | ND            | 0.039       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | ND            | 0.048       | 0.10000 | 0.0000 | 0.0024 | 0.0048 |
| 1,2,3,6,7,8-HxCDF   | 0.094         | 0.035       | 0.10000 | 0.0094 | 0.0094 | 0.0094 |
| 2,3,4,6,7,8-HxCDF   | ND            | 0.028       | 0.10000 | 0.0028 | 0.0028 | 0.0028 |
| 1,2,3,7,8,9-HxCDF   | ND            | 0.042       | 0.10000 | 0.0000 | 0.0021 | 0.0042 |
| Total HxCDF         | 0.094         | 0.038       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | ND            | 0.050       | 0.10000 | 0.0000 | 0.0025 | 0.0050 |
| 1,2,3,6,7,8-HxCDD   | ND            | 0.052       | 0.10000 | 0.0081 | 0.0081 | 0.0081 |
| 1,2,3,7,8,9-HxCDD   | ND            | 0.048       | 0.10000 | 0.0000 | 0.0024 | 0.0048 |
| Total HxCDD         | ND            | 0.050       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | ND            | 0.056       | 0.01000 | 0.0010 | 0.0010 | 0.0010 |
| 1,2,3,4,7,8,9-HpCDF | ND            | 0.047       | 0.01000 | 0.0000 | 0.0002 | 0.0005 |
| Total HpCDF         | ND            | 0.051       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | ND            | 0.054       | 0.01000 | 0.0007 | 0.0007 | 0.0007 |
| Total HpCDD         | 0.100         | 0.054       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | 0.200         | 0.12        | 0.00030 | 0.0001 | 0.0001 | 0.0001 |
| OCDD                | ND            | 0.15        | 0.00030 | 0.0001 | 0.0001 | 0.0001 |

**0.19 ng/Kg      0.20 ng/Kg      0.22 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Blank Analysis Results**

|                        |                           |             |                  |
|------------------------|---------------------------|-------------|------------------|
| Lab Sample ID          | BLANK-39465               | Matrix      | Solid            |
| Filename               | P140301A_05               | Dilution    | NA               |
| Total Amount Extracted | 10.4 g                    | Extracted   | 02/26/2014 21:30 |
| ICAL ID                | P130624                   | Analyzed    | 03/01/2014 10:03 |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Injected By | BAL              |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | -----      | 0.14       | 0.140 IJ | 2,3,7,8-TCDF-13C         | 2.00       | 75               |
| Total TCDF          | ND         | -----      | 0.140    | 2,3,7,8-TCDD-13C         | 2.00       | 82               |
|                     |            |            |          | 1,2,3,7,8-PeCDF-13C      | 2.00       | 75               |
| 2,3,7,8-TCDD        | ND         | -----      | 0.150    | 2,3,4,7,8-PeCDF-13C      | 2.00       | 70               |
| Total TCDD          | ND         | -----      | 0.150    | 1,2,3,7,8-PeCDD-13C      | 2.00       | 73               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 83               |
| 1,2,3,7,8-PeCDF     | ND         | -----      | 0.160    | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 87               |
| 2,3,4,7,8-PeCDF     | ND         | -----      | 0.110    | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 83               |
| Total PeCDF         | ND         | -----      | 0.130    | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 83               |
|                     |            |            |          | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 79               |
| 1,2,3,7,8-PeCDD     | ND         | -----      | 0.160    | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 72               |
| Total PeCDD         | ND         | -----      | 0.160    | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 77               |
|                     |            |            |          | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 74               |
| 1,2,3,4,7,8-HxCDF   | ND         | -----      | 0.110    | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 76               |
| 1,2,3,6,7,8-HxCDF   | ND         | -----      | 0.084    | OCDD-13C                 | 4.00       | 81 Y             |
| 2,3,4,6,7,8-HxCDF   | ND         | -----      | 0.091    |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | ND         | -----      | 0.110    | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | ND         | -----      | 0.099    | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | ND         | -----      | 0.130    | 2,3,7,8-TCDD-37Cl4       | 0.20       | 90               |
| 1,2,3,6,7,8-HxCDD   | ND         | -----      | 0.130    |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | ND         | -----      | 0.130    |                          |            |                  |
| Total HxCDD         | ND         | -----      | 0.130    |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | -----      | 0.20       | 0.120 IJ | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | ND         | -----      | 0.170    | Equivalence: 0.018 ng/Kg |            |                  |
| Total HpCDF         | ND         | -----      | 0.140    | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 0.17       | -----      | 0.140 J  |                          |            |                  |
| Total HpCDD         | 0.44       | -----      | 0.140 J  |                          |            |                  |
|                     |            |            |          |                          |            |                  |
| OCDF                | ND         | -----      | 0.260    |                          |            |                  |
| OCDD                | 0.95       | -----      | 0.330 J  |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

J = Estimated value

I = Interference present

Y = Calculated using average of daily RFs

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |  |
|------------------------|---------------------------|-----------|------------------|--|--|
| Client's Sample ID     |                           |           |                  |  |  |
| Lab Sample ID          | BLANK-39465               |           |                  |  |  |
| Filename               | P140301A_05               |           |                  |  |  |
| Injected By            | BAL                       |           |                  |  |  |
| Total Amount Extracted | 10.4 g                    | Matrix    | Solid            |  |  |
| % Moisture             | 0.0                       | Dilution  | NA               |  |  |
| Dry Weight Extracted   | 10.4 g                    | Collected | 02/25/2014 20:04 |  |  |
| ICAL ID                | P130624                   | Received  | 02/25/2014 20:04 |  |  |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Extracted | 02/26/2014 21:30 |  |  |
| Method Blank ID        |                           | Analyzed  | 03/01/2014 10:03 |  |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | ND            | 0.14        | 0.10000 | 0.0143 | 0.0143 | 0.0143 |
| Total TCDF          | ND            | 0.14        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | ND            | 0.15        | 1.00000 | 0.0000 | 0.0752 | 0.1504 |
| Total TCDD          | ND            | 0.15        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | ND            | 0.16        | 0.03000 | 0.0000 | 0.0023 | 0.0047 |
| 2,3,4,7,8-PeCDF     | ND            | 0.11        | 0.30000 | 0.0000 | 0.0169 | 0.0339 |
| Total PeCDF         | ND            | 0.13        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | ND            | 0.16        | 1.00000 | 0.0000 | 0.0775 | 0.1550 |
| Total PeCDD         | ND            | 0.16        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | ND            | 0.11        | 0.10000 | 0.0000 | 0.0053 | 0.0107 |
| 1,2,3,6,7,8-HxCDF   | ND            | 0.084       | 0.10000 | 0.0000 | 0.0042 | 0.0084 |
| 2,3,4,6,7,8-HxCDF   | ND            | 0.091       | 0.10000 | 0.0000 | 0.0045 | 0.0091 |
| 1,2,3,7,8,9-HxCDF   | ND            | 0.11        | 0.10000 | 0.0000 | 0.0057 | 0.0114 |
| Total HxCDF         | ND            | 0.099       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | ND            | 0.13        | 0.10000 | 0.0000 | 0.0067 | 0.0134 |
| 1,2,3,6,7,8-HxCDD   | ND            | 0.13        | 0.10000 | 0.0000 | 0.0064 | 0.0129 |
| 1,2,3,7,8,9-HxCDD   | ND            | 0.13        | 0.10000 | 0.0000 | 0.0065 | 0.0130 |
| Total HxCDD         | ND            | 0.13        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | ND            | 0.12        | 0.01000 | 0.0020 | 0.0020 | 0.0020 |
| 1,2,3,4,7,8,9-HpCDF | ND            | 0.17        | 0.01000 | 0.0000 | 0.0008 | 0.0017 |
| Total HpCDF         | ND            | 0.14        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | 0.17          | 0.14        | 0.01000 | 0.0017 | 0.0017 | 0.0017 |
| Total HpCDD         | 0.44          | 0.14        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | ND            | 0.26        | 0.00030 | 0.0000 | 0.0000 | 0.0001 |
| OCDD                | 0.95          | 0.33        | 0.00030 | 0.0003 | 0.0003 | 0.0003 |

**0.018 ng/Kg      0.23 ng/Kg      0.44 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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**Method 8290 Blank Analysis Results**

|                        |                           |             |                  |
|------------------------|---------------------------|-------------|------------------|
| Lab Sample ID          | BLANK-39477               | Matrix      | Solid            |
| Filename               | P140304B_01               | Dilution    | NA               |
| Total Amount Extracted | 10.1 g                    | Extracted   | 02/27/2014 19:30 |
| ICAL ID                | P130624                   | Analyzed    | 03/04/2014 17:06 |
| CCal Filename(s)       | P140304A_11 & P140304B_15 | Injected By | SMT              |

| Native Isomers      | Conc ng/Kg | EMPC ng/Kg | RL ng/Kg |    | Internal Standards       | ng's Added | Percent Recovery |
|---------------------|------------|------------|----------|----|--------------------------|------------|------------------|
| 2,3,7,8-TCDF        | -----      | 0.110      | 0.076    | IJ | 2,3,7,8-TCDF-13C         | 2.00       | 72               |
| Total TCDF          | ND         | -----      | 0.076    |    | 2,3,7,8-TCDD-13C         | 2.00       | 86               |
|                     |            |            |          |    | 1,2,3,7,8-PeCDF-13C      | 2.00       | 79               |
| 2,3,7,8-TCDD        | ND         | -----      | 0.110    |    | 2,3,4,7,8-PeCDF-13C      | 2.00       | 78               |
| Total TCDD          | ND         | -----      | 0.110    |    | 1,2,3,7,8-PeCDD-13C      | 2.00       | 87               |
|                     |            |            |          |    | 1,2,3,4,7,8-HxCDF-13C    | 2.00       | 89               |
| 1,2,3,7,8-PeCDF     | -----      | 0.089      | 0.086    | IJ | 1,2,3,6,7,8-HxCDF-13C    | 2.00       | 88               |
| 2,3,4,7,8-PeCDF     | 0.10       | -----      | 0.069    | J  | 2,3,4,6,7,8-HxCDF-13C    | 2.00       | 86               |
| Total PeCDF         | 0.10       | -----      | 0.077    | J  | 1,2,3,7,8,9-HxCDF-13C    | 2.00       | 87               |
|                     |            |            |          |    | 1,2,3,4,7,8-HxCDD-13C    | 2.00       | 87               |
| 1,2,3,7,8-PeCDD     | ND         | -----      | 0.140    |    | 1,2,3,6,7,8-HxCDD-13C    | 2.00       | 75               |
| Total PeCDD         | ND         | -----      | 0.140    |    | 1,2,3,4,6,7,8-HpCDF-13C  | 2.00       | 85               |
|                     |            |            |          |    | 1,2,3,4,7,8,9-HpCDF-13C  | 2.00       | 83               |
| 1,2,3,4,7,8-HxCDF   | -----      | 0.120      | 0.048    | IJ | 1,2,3,4,6,7,8-HpCDD-13C  | 2.00       | 89               |
| 1,2,3,6,7,8-HxCDF   | 0.10       | -----      | 0.041    | J  | OCDD-13C                 | 4.00       | 68               |
| 2,3,4,6,7,8-HxCDF   | 0.17       | -----      | 0.043    | J  |                          |            |                  |
| 1,2,3,7,8,9-HxCDF   | 0.14       | -----      | 0.078    | J  | 1,2,3,4-TCDD-13C         | 2.00       | NA               |
| Total HxCDF         | 0.41       | -----      | 0.053    | J  | 1,2,3,7,8,9-HxCDD-13C    | 2.00       | NA               |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,7,8-HxCDD   | 0.14       | -----      | 0.130    | J  | 2,3,7,8-TCDD-37Cl4       | 0.20       | 83               |
| 1,2,3,6,7,8-HxCDD   | -----      | 0.140      | 0.120    | IJ |                          |            |                  |
| 1,2,3,7,8,9-HxCDD   | -----      | 0.120      | 0.094    | IJ |                          |            |                  |
| Total HxCDD         | 0.14       | -----      | 0.110    | J  |                          |            |                  |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDF | -----      | 0.150      | 0.085    | IJ | Total 2,3,7,8-TCDD       |            |                  |
| 1,2,3,4,7,8,9-HpCDF | -----      | 0.220      | 0.078    | IJ | Equivalence: 0.15 ng/Kg  |            |                  |
| Total HpCDF         | ND         | -----      | 0.082    |    | (Using 2005 WHO Factors) |            |                  |
|                     |            |            |          |    |                          |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 0.23       | -----      | 0.110    | J  |                          |            |                  |
| Total HpCDD         | 0.41       | -----      | 0.110    | J  |                          |            |                  |
|                     |            |            |          |    |                          |            |                  |
| OCDF                | 0.70       | -----      | 0.160    | J  |                          |            |                  |
| OCDD                | 1.30       | -----      | 0.220    | J  |                          |            |                  |

Conc = Concentration (Totals include 2,3,7,8-substituted isomers).  
EMPC = Estimated Maximum Possible Concentration  
RL = Reporting Limit

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

J = Estimated value

I = Interference present

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**2,3,7,8-TCDD Toxic Equivalency (TEQ) Calculations**  
ESN Pacific

|                        |                           |           |                  |  |  |
|------------------------|---------------------------|-----------|------------------|--|--|
| Client's Sample ID     |                           |           |                  |  |  |
| Lab Sample ID          | BLANK-39477               |           |                  |  |  |
| Filename               | P140304B_01               |           |                  |  |  |
| Injected By            | SMT                       |           |                  |  |  |
| Total Amount Extracted | 10.1 g                    | Matrix    | Solid            |  |  |
| % Moisture             | 0.0                       | Dilution  | NA               |  |  |
| Dry Weight Extracted   | 10.1 g                    | Collected | 02/26/2014 19:20 |  |  |
| ICAL ID                | P130624                   | Received  | 02/26/2014 19:20 |  |  |
| CCal Filename(s)       | P140304A_11 & P140304B_15 | Extracted | 02/27/2014 19:30 |  |  |
| Method Blank ID        |                           | Analyzed  | 03/04/2014 17:06 |  |  |

| Parameter           | Conc<br>ng/Kg | RL<br>ng/Kg | WHO2005 | LB     | MB     | UB     |
|---------------------|---------------|-------------|---------|--------|--------|--------|
| 2,3,7,8-TCDF        | ND            | 0.076       | 0.10000 | 0.0115 | 0.0115 | 0.0115 |
| Total TCDF          | ND            | 0.076       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 2,3,7,8-TCDD        | ND            | 0.11        | 1.00000 | 0.0000 | 0.0525 | 0.1050 |
| Total TCDD          | ND            | 0.11        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDF     | ND            | 0.086       | 0.03000 | 0.0027 | 0.0027 | 0.0027 |
| 2,3,4,7,8-PeCDF     | 0.10          | 0.069       | 0.30000 | 0.0308 | 0.0308 | 0.0308 |
| Total PeCDF         | 0.10          | 0.077       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,7,8-PeCDD     | ND            | 0.14        | 1.00000 | 0.0000 | 0.0711 | 0.1422 |
| Total PeCDD         | ND            | 0.14        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDF   | ND            | 0.048       | 0.10000 | 0.0118 | 0.0118 | 0.0118 |
| 1,2,3,6,7,8-HxCDF   | 0.10          | 0.041       | 0.10000 | 0.0104 | 0.0104 | 0.0104 |
| 2,3,4,6,7,8-HxCDF   | 0.17          | 0.043       | 0.10000 | 0.0171 | 0.0171 | 0.0171 |
| 1,2,3,7,8,9-HxCDF   | 0.14          | 0.078       | 0.10000 | 0.0138 | 0.0138 | 0.0138 |
| Total HxCDF         | 0.41          | 0.053       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,7,8-HxCDD   | 0.14          | 0.13        | 0.10000 | 0.0142 | 0.0142 | 0.0142 |
| 1,2,3,6,7,8-HxCDD   | ND            | 0.12        | 0.10000 | 0.0143 | 0.0143 | 0.0143 |
| 1,2,3,7,8,9-HxCDD   | ND            | 0.094       | 0.10000 | 0.0121 | 0.0121 | 0.0121 |
| Total HxCDD         | 0.14          | 0.11        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDF | ND            | 0.085       | 0.01000 | 0.0015 | 0.0015 | 0.0015 |
| 1,2,3,4,7,8,9-HpCDF | ND            | 0.078       | 0.01000 | 0.0022 | 0.0022 | 0.0022 |
| Total HpCDF         | ND            | 0.082       | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| 1,2,3,4,6,7,8-HpCDD | 0.23          | 0.11        | 0.01000 | 0.0023 | 0.0023 | 0.0023 |
| Total HpCDD         | 0.41          | 0.11        | 0.00000 | 0.0000 | 0.0000 | 0.0000 |
| OCDF                | 0.70          | 0.16        | 0.00030 | 0.0002 | 0.0002 | 0.0002 |
| OCDD                | 1.30          | 0.22        | 0.00030 | 0.0004 | 0.0004 | 0.0004 |

**0.15 ng/Kg      0.27 ng/Kg      0.39 ng/Kg**

Final values are valid to only 2 significant figures  
TEQs for Totals classes include contributions from non 2,3,7,8 isomers only  
LB = Lower Bound, Where "ND", TEQ Conc = 0  
MB = Medium Bound, Where "ND", TEQ Conc = (LOD/2) \* (TEF Factor)  
UB = Upper Bound, Where "ND", TEQ Conc = LOD \* (TEF Factor)  
RL = Reporting Limit

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### Method 8290 Laboratory Control Spike Results

|                        |                           |             |                  |
|------------------------|---------------------------|-------------|------------------|
| Lab Sample ID          | LCS-39447                 | Matrix      | Solid            |
| Filename               | F140226B_15               | Dilution    | NA               |
| Total Amount Extracted | 10.0 g                    | Extracted   | 02/24/2014 21:00 |
| ICAL ID                | F131125                   | Analyzed    | 02/27/2014 06:48 |
| CCal Filename(s)       | F140226B_01 & F140226B_17 | Injected By | BAL              |
| Method Blank ID        | BLANK-39446               |             |                  |

| Native Isomers      | Qs (ng) | Qm (ng) | % Rec. | Internal Standards      | ng's Added | Percent Recovery |
|---------------------|---------|---------|--------|-------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 0.20    | 0.23    | 115    | 2,3,7,8-TCDF-13C        | 2.0        | 75               |
| Total TCDF          |         |         |        | 2,3,7,8-TCDD-13C        | 2.0        | 85               |
|                     |         |         |        | 1,2,3,7,8-PeCDF-13C     | 2.0        | 85               |
| 2,3,7,8-TCDD        | 0.20    | 0.18    | 89     | 2,3,4,7,8-PeCDF-13C     | 2.0        | 85               |
| Total TCDD          |         |         |        | 1,2,3,7,8-PeCDD-13C     | 2.0        | 97               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDF-13C   | 2.0        | 81               |
| 1,2,3,7,8-PeCDF     | 1.0     | 1.2     | 117    | 1,2,3,6,7,8-HxCDF-13C   | 2.0        | 89               |
| 2,3,4,7,8-PeCDF     | 1.0     | 1.1     | 109    | 2,3,4,6,7,8-HxCDF-13C   | 2.0        | 87               |
| Total PeCDF         |         |         |        | 1,2,3,7,8,9-HxCDF-13C   | 2.0        | 82               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDD-13C   | 2.0        | 77               |
| 1,2,3,7,8-PeCDD     | 1.0     | 1.0     | 100    | 1,2,3,6,7,8-HxCDD-13C   | 2.0        | 78               |
| Total PeCDD         |         |         |        | 1,2,3,4,6,7,8-HpCDF-13C | 2.0        | 78               |
|                     |         |         |        | 1,2,3,4,7,8,9-HpCDF-13C | 2.0        | 75               |
| 1,2,3,4,7,8-HxCDF   | 1.0     | 1.1     | 110    | 1,2,3,4,6,7,8-HpCDD-13C | 2.0        | 83               |
| 1,2,3,6,7,8-HxCDF   | 1.0     | 1.1     | 109    | OCDD-13C                | 4.0        | 69               |
| 2,3,4,6,7,8-HxCDF   | 1.0     | 1.0     | 100    |                         |            |                  |
| 1,2,3,7,8,9-HxCDF   | 1.0     | 1.1     | 107    | 1,2,3,4-TCDD-13C        | 2.0        | NA               |
| Total HxCDF         |         |         |        | 1,2,3,7,8,9-HxCDD-13C   | 2.0        | NA               |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,7,8-HxCDD   | 1.0     | 1.2     | 115    | 2,3,7,8-TCDD-37Cl4      | 0.20       | 87               |
| 1,2,3,6,7,8-HxCDD   | 1.0     | 1.3     | 128    |                         |            |                  |
| 1,2,3,7,8,9-HxCDD   | 1.0     | 1.2     | 119    |                         |            |                  |
| Total HxCDD         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 1.0     | 1.1     | 106    |                         |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 1.0     | 0.94    | 94     |                         |            |                  |
| Total HpCDF         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 1.0     | 1.0     | 102    |                         |            |                  |
| Total HpCDD         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| OCDF                | 2.0     | 2.0     | 99     |                         |            |                  |
| OCDD                | 2.0     | 2.2     | 110    |                         |            |                  |

Qs = Quantity Spiked  
Qm = Quantity Measured  
Rec. = Recovery (Expressed as Percent)  
R = Recovery outside of target range

Y = RF averaging used in calculations  
Nn = Value obtained from additional analysis  
NA = Not Applicable  
\* = See Discussion

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**Method 8290 Laboratory Control Spike Results**

|                        |                           |             |                  |
|------------------------|---------------------------|-------------|------------------|
| Lab Sample ID          | LCS-39466                 | Matrix      | Solid            |
| Filename               | P140301A_01               | Dilution    | NA               |
| Total Amount Extracted | 10.6 g                    | Extracted   | 02/26/2014 21:30 |
| ICAL ID                | P130624                   | Analyzed    | 03/01/2014 07:12 |
| CCal Filename(s)       | P140228B_19 & P140301A_15 | Injected By | BAL              |
| Method Blank ID        | BLANK-39465               |             |                  |

| Native Isomers      | Qs (ng) | Qm (ng) | % Rec. | Internal Standards      | ng's Added | Percent Recovery |
|---------------------|---------|---------|--------|-------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 0.20    | 0.22    | 112    | 2,3,7,8-TCDF-13C        | 2.0        | 79               |
| Total TCDF          |         |         |        | 2,3,7,8-TCDD-13C        | 2.0        | 85               |
|                     |         |         |        | 1,2,3,7,8-PeCDF-13C     | 2.0        | 75               |
| 2,3,7,8-TCDD        | 0.20    | 0.18    | 89     | 2,3,4,7,8-PeCDF-13C     | 2.0        | 71               |
| Total TCDD          |         |         |        | 1,2,3,7,8-PeCDD-13C     | 2.0        | 75               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDF-13C   | 2.0        | 88               |
| 1,2,3,7,8-PeCDF     | 1.0     | 1.2     | 117    | 1,2,3,6,7,8-HxCDF-13C   | 2.0        | 90               |
| 2,3,4,7,8-PeCDF     | 1.0     | 1.1     | 111    | 2,3,4,6,7,8-HxCDF-13C   | 2.0        | 86               |
| Total PeCDF         |         |         |        | 1,2,3,7,8,9-HxCDF-13C   | 2.0        | 88               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDD-13C   | 2.0        | 80               |
| 1,2,3,7,8-PeCDD     | 1.0     | 0.97    | 97     | 1,2,3,6,7,8-HxCDD-13C   | 2.0        | 75               |
| Total PeCDD         |         |         |        | 1,2,3,4,6,7,8-HpCDF-13C | 2.0        | 82               |
|                     |         |         |        | 1,2,3,4,7,8,9-HpCDF-13C | 2.0        | 77               |
| 1,2,3,4,7,8-HxCDF   | 1.0     | 1.1     | 111    | 1,2,3,4,6,7,8-HpCDD-13C | 2.0        | 80               |
| 1,2,3,6,7,8-HxCDF   | 1.0     | 1.1     | 112    | OCDD-13C                | 4.0        | 100 Y            |
| 2,3,4,6,7,8-HxCDF   | 1.0     | 1.1     | 112    |                         |            |                  |
| 1,2,3,7,8,9-HxCDF   | 1.0     | 1.1     | 106    | 1,2,3,4-TCDD-13C        | 2.0        | NA               |
| Total HxCDF         |         |         |        | 1,2,3,7,8,9-HxCDD-13C   | 2.0        | NA               |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,7,8-HxCDD   | 1.0     | 1.2     | 115    | 2,3,7,8-TCDD-37Cl4      | 0.20       | 85               |
| 1,2,3,6,7,8-HxCDD   | 1.0     | 1.2     | 122    |                         |            |                  |
| 1,2,3,7,8,9-HxCDD   | 1.0     | 1.2     | 118    |                         |            |                  |
| Total HxCDD         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 1.0     | 1.1     | 114    |                         |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 1.0     | 1.0     | 103    |                         |            |                  |
| Total HpCDF         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 1.0     | 1.1     | 106    |                         |            |                  |
| Total HpCDD         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| OCDF                | 2.0     | 2.2     | 108    |                         |            |                  |
| OCDD                | 2.0     | 2.2     | 110    |                         |            |                  |

Qs = Quantity Spiked  
Qm = Quantity Measured  
Rec. = Recovery (Expressed as Percent)  
R = Recovery outside of target range

Y = RF averaging used in calculations  
Nn = Value obtained from additional analysis  
NA = Not Applicable  
\* = See Discussion

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**Method 8290 Laboratory Control Spike Results**

|                        |                           |             |                  |
|------------------------|---------------------------|-------------|------------------|
| Lab Sample ID          | LCS-39478                 | Matrix      | Solid            |
| Filename               | P140304B_14               | Dilution    | NA               |
| Total Amount Extracted | 10.3 g                    | Extracted   | 02/27/2014 19:30 |
| ICAL ID                | P130624                   | Analyzed    | 03/05/2014 02:21 |
| CCal Filename(s)       | P140304A_11 & P140304B_15 | Injected By | SMT              |
| Method Blank ID        | BLANK-39477               |             |                  |

| Native Isomers      | Qs (ng) | Qm (ng) | % Rec. | Internal Standards      | ng's Added | Percent Recovery |
|---------------------|---------|---------|--------|-------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 0.20    | 0.23    | 115    | 2,3,7,8-TCDF-13C        | 2.0        | 81               |
| Total TCDF          |         |         |        | 2,3,7,8-TCDD-13C        | 2.0        | 96               |
|                     |         |         |        | 1,2,3,7,8-PeCDF-13C     | 2.0        | 89               |
| 2,3,7,8-TCDD        | 0.20    | 0.19    | 93     | 2,3,4,7,8-PeCDF-13C     | 2.0        | 87               |
| Total TCDD          |         |         |        | 1,2,3,7,8-PeCDD-13C     | 2.0        | 100              |
|                     |         |         |        | 1,2,3,4,7,8-HxCDF-13C   | 2.0        | 94               |
| 1,2,3,7,8-PeCDF     | 1.0     | 1.2     | 116    | 1,2,3,6,7,8-HxCDF-13C   | 2.0        | 94               |
| 2,3,4,7,8-PeCDF     | 1.0     | 1.1     | 108    | 2,3,4,6,7,8-HxCDF-13C   | 2.0        | 91               |
| Total PeCDF         |         |         |        | 1,2,3,7,8,9-HxCDF-13C   | 2.0        | 94               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDD-13C   | 2.0        | 96               |
| 1,2,3,7,8-PeCDD     | 1.0     | 0.96    | 96     | 1,2,3,6,7,8-HxCDD-13C   | 2.0        | 81               |
| Total PeCDD         |         |         |        | 1,2,3,4,6,7,8-HpCDF-13C | 2.0        | 90               |
|                     |         |         |        | 1,2,3,4,7,8,9-HpCDF-13C | 2.0        | 87               |
| 1,2,3,4,7,8-HxCDF   | 1.0     | 1.1     | 110    | 1,2,3,4,6,7,8-HpCDD-13C | 2.0        | 91               |
| 1,2,3,6,7,8-HxCDF   | 1.0     | 1.1     | 107    | OCDD-13C                | 4.0        | 70               |
| 2,3,4,6,7,8-HxCDF   | 1.0     | 1.1     | 107    |                         |            |                  |
| 1,2,3,7,8,9-HxCDF   | 1.0     | 1.1     | 105    | 1,2,3,4-TCDD-13C        | 2.0        | NA               |
| Total HxCDF         |         |         |        | 1,2,3,7,8,9-HxCDD-13C   | 2.0        | NA               |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,7,8-HxCDD   | 1.0     | 1.1     | 109    | 2,3,7,8-TCDD-37Cl4      | 0.20       | 92               |
| 1,2,3,6,7,8-HxCDD   | 1.0     | 1.2     | 120    |                         |            |                  |
| 1,2,3,7,8,9-HxCDD   | 1.0     | 1.1     | 113    |                         |            |                  |
| Total HxCDD         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 1.0     | 1.1     | 109    |                         |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 1.0     | 0.98    | 98     |                         |            |                  |
| Total HpCDF         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 1.0     | 1.0     | 102    |                         |            |                  |
| Total HpCDD         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| OCDF                | 2.0     | 2.3     | 115    |                         |            |                  |
| OCDD                | 2.0     | 2.2     | 112    |                         |            |                  |

Qs = Quantity Spiked  
Qm = Quantity Measured  
Rec. = Recovery (Expressed as Percent)  
R = Recovery outside of target range

Y = RF averaging used in calculations  
Nn = Value obtained from additional analysis  
NA = Not Applicable  
\* = See Discussion

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**Method 8290 Laboratory Control Spike Results**

|                        |                           |             |                  |
|------------------------|---------------------------|-------------|------------------|
| Lab Sample ID          | LCSD-39448                | Matrix      | Solid            |
| Filename               | F140226B_16               | Dilution    | NA               |
| Total Amount Extracted | 10.5 g                    | Extracted   | 02/24/2014 21:00 |
| ICAL ID                | F131125                   | Analyzed    | 02/27/2014 07:34 |
| CCal Filename(s)       | F140226B_01 & F140226B_17 | Injected By | BAL              |
| Method Blank ID        | BLANK-39446               |             |                  |

| Native Isomers      | Qs (ng) | Qm (ng) | % Rec. | Internal Standards      | ng's Added | Percent Recovery |
|---------------------|---------|---------|--------|-------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 0.20    | 0.24    | 119    | 2,3,7,8-TCDF-13C        | 2.0        | 64               |
| Total TCDF          |         |         |        | 2,3,7,8-TCDD-13C        | 2.0        | 74               |
|                     |         |         |        | 1,2,3,7,8-PeCDF-13C     | 2.0        | 78               |
| 2,3,7,8-TCDD        | 0.20    | 0.18    | 91     | 2,3,4,7,8-PeCDF-13C     | 2.0        | 81               |
| Total TCDD          |         |         |        | 1,2,3,7,8-PeCDD-13C     | 2.0        | 90               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDF-13C   | 2.0        | 78               |
| 1,2,3,7,8-PeCDF     | 1.0     | 1.2     | 119    | 1,2,3,6,7,8-HxCDF-13C   | 2.0        | 90               |
| 2,3,4,7,8-PeCDF     | 1.0     | 1.1     | 109    | 2,3,4,6,7,8-HxCDF-13C   | 2.0        | 87               |
| Total PeCDF         |         |         |        | 1,2,3,7,8,9-HxCDF-13C   | 2.0        | 78               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDD-13C   | 2.0        | 73               |
| 1,2,3,7,8-PeCDD     | 1.0     | 1.0     | 103    | 1,2,3,6,7,8-HxCDD-13C   | 2.0        | 81               |
| Total PeCDD         |         |         |        | 1,2,3,4,6,7,8-HpCDF-13C | 2.0        | 78               |
|                     |         |         |        | 1,2,3,4,7,8,9-HpCDF-13C | 2.0        | 74               |
| 1,2,3,4,7,8-HxCDF   | 1.0     | 1.2     | 115    | 1,2,3,4,6,7,8-HpCDD-13C | 2.0        | 84               |
| 1,2,3,6,7,8-HxCDF   | 1.0     | 1.1     | 107    | OCDD-13C                | 4.0        | 73               |
| 2,3,4,6,7,8-HxCDF   | 1.0     | 1.0     | 102    |                         |            |                  |
| 1,2,3,7,8,9-HxCDF   | 1.0     | 1.1     | 107    | 1,2,3,4-TCDD-13C        | 2.0        | NA               |
| Total HxCDF         |         |         |        | 1,2,3,7,8,9-HxCDD-13C   | 2.0        | NA               |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,7,8-HxCDD   | 1.0     | 1.2     | 124    | 2,3,7,8-TCDD-37Cl4      | 0.20       | 75               |
| 1,2,3,6,7,8-HxCDD   | 1.0     | 1.2     | 123    |                         |            |                  |
| 1,2,3,7,8,9-HxCDD   | 1.0     | 1.2     | 121    |                         |            |                  |
| Total HxCDD         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 1.0     | 1.1     | 108    |                         |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 1.0     | 0.94    | 94     |                         |            |                  |
| Total HpCDF         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 1.0     | 1.0     | 102    |                         |            |                  |
| Total HpCDD         |         |         |        |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| OCDF                | 2.0     | 2.0     | 101    |                         |            |                  |
| OCDD                | 2.0     | 2.2     | 110    |                         |            |                  |

Qs = Quantity Spiked  
Qm = Quantity Measured  
Rec. = Recovery (Expressed as Percent)  
R = Recovery outside of target range

Y = RF averaging used in calculations  
Nn = Value obtained from additional analysis  
NA = Not Applicable  
\* = See Discussion

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**Method 8290**

**Spike Recovery Relative Percent Difference (RPD) Results**

Client ESN Pacific

Spike 1 ID LCS-39447  
Spike 1 Filename F140226B\_15

Spike 2 ID LCSD-39448  
Spike 2 Filename F140226B\_16

| Compound            | Spike 1<br>%REC | Spike 2<br>%REC | %RPD |
|---------------------|-----------------|-----------------|------|
| 2,3,7,8-TCDF        | 115             | 119             | 3.4  |
| 2,3,7,8-TCDD        | 89              | 91              | 2.2  |
| 1,2,3,7,8-PeCDF     | 117             | 119             | 1.7  |
| 2,3,4,7,8-PeCDF     | 109             | 109             | 0.0  |
| 1,2,3,7,8-PeCDD     | 100             | 103             | 3.0  |
| 1,2,3,4,7,8-HxCDF   | 110             | 115             | 4.4  |
| 1,2,3,6,7,8-HxCDF   | 109             | 107             | 1.9  |
| 2,3,4,6,7,8-HxCDF   | 100             | 102             | 2.0  |
| 1,2,3,7,8,9-HxCDF   | 107             | 107             | 0.0  |
| 1,2,3,4,7,8-HxCDD   | 115             | 124             | 7.5  |
| 1,2,3,6,7,8-HxCDD   | 128             | 123             | 4.0  |
| 1,2,3,7,8,9-HxCDD   | 119             | 121             | 1.7  |
| 1,2,3,4,6,7,8-HpCDF | 106             | 108             | 1.9  |
| 1,2,3,4,7,8,9-HpCDF | 94              | 94              | 0.0  |
| 1,2,3,4,6,7,8-HpCDD | 102             | 102             | 0.0  |
| OCDF                | 99              | 101             | 2.0  |
| OCDD                | 110             | 110             | 0.0  |

%REC = Percent Recovered

RPD = The difference between the two values divided by the mean value

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### Method 8290 Spiked Sample Report

Client - ESN Pacific

|                        |                           |             |                  |
|------------------------|---------------------------|-------------|------------------|
| Client's Sample ID     | RHS-DU-20-MS              | Matrix      | Soil             |
| Lab Sample ID          | 10257348006-S-MS          | Dilution    | NA               |
| Filename               | U140304B_03               | Extracted   | 02/27/2014 19:30 |
| Total Amount Extracted | 10.5 g                    | Analyzed    | 03/04/2014 18:10 |
| ICAL ID                | U140224                   | Injected By | SMT              |
| CCal Filename(s)       | U140304B_02 & U140304B_20 |             |                  |
| Method Blank ID        | BLANK-39477               |             |                  |

| Native Isomers      | Qs (ng) | Qm (ng) | % Rec. | Internal Standards      | ng's Added | Percent Recovery |
|---------------------|---------|---------|--------|-------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 0.20    | 0.34    | 169    | 2,3,7,8-TCDF-13C        | 2.00       | 95               |
|                     |         |         |        | 2,3,7,8-TCDD-13C        | 2.00       | 101              |
|                     |         |         |        | 1,2,3,7,8-PeCDF-13C     | 2.00       | 104              |
| 2,3,7,8-TCDD        | 0.20    | 0.19    | 95     | 2,3,4,7,8-PeCDF-13C     | 2.00       | 82 Y             |
|                     |         |         |        | 1,2,3,7,8-PeCDD-13C     | 2.00       | 112              |
|                     |         |         |        | 1,2,3,4,7,8-HxCDF-13C   | 2.00       | 89               |
| 1,2,3,7,8-PeCDF     | 1.00    | 1.29    | 129    | 1,2,3,6,7,8-HxCDF-13C   | 2.00       | 95               |
| 2,3,4,7,8-PeCDF     | 1.00    | 1.38    | 138    | 2,3,4,6,7,8-HxCDF-13C   | 2.00       | 101              |
|                     |         |         |        | 1,2,3,7,8,9-HxCDF-13C   | 2.00       | 92               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDD-13C   | 2.00       | 79               |
| 1,2,3,7,8-PeCDD     | 1.00    | 1.08    | 108    | 1,2,3,6,7,8-HxCDD-13C   | 2.00       | 81               |
|                     |         |         |        | 1,2,3,4,6,7,8-HpCDF-13C | 2.00       | 76               |
|                     |         |         |        | 1,2,3,4,7,8,9-HpCDF-13C | 2.00       | 77               |
| 1,2,3,4,7,8-HxCDF   | 1.00    | 1.32    | 132    | 1,2,3,4,6,7,8-HpCDD-13C | 2.00       | 84               |
| 1,2,3,6,7,8-HxCDF   | 1.00    | 1.32    | 132    | OCDD-13C                | 4.00       | 70               |
| 2,3,4,6,7,8-HxCDF   | 1.00    | 1.28    | 128    |                         |            |                  |
| 1,2,3,7,8,9-HxCDF   | 1.00    | 1.20    | 120    | 1,2,3,4-TCDD-13C        | 2.00       | NA               |
|                     |         |         |        | 1,2,3,7,8,9-HxCDD-13C   | 2.00       | NA               |
| 1,2,3,4,7,8-HxCDD   | 1.00    | 1.26    | 126    | 2,3,7,8-TCDD-37Cl4      | 0.20       | 102              |
| 1,2,3,6,7,8-HxCDD   | 1.00    | 1.28    | 128    |                         |            |                  |
| 1,2,3,7,8,9-HxCDD   | 1.00    | 1.32    | 132    |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 1.00    | 1.80    | 180    |                         |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 1.00    | 1.07    | 107    |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 1.00    | 1.93    | 193    |                         |            |                  |
| OCDF                | 2.00    | 2.87    | 144    |                         |            |                  |
| OCDD                | 2.00    | 8.40    | 420    |                         |            |                  |

Qs = Quantity Spiked                      Qm = Quantity Measured                      Rec. = Recovery (Expressed as Percent)

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

Y = Calculated using average of daily RFs

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### Method 8290 Spiked Sample Report

Client - ESN Pacific

|                        |                           |             |                  |
|------------------------|---------------------------|-------------|------------------|
| Client's Sample ID     | RHS-DU-20-MSD             | Matrix      | Soil             |
| Lab Sample ID          | 10257348006-S-MSD         | Dilution    | NA               |
| Filename               | U140304B_04               | Extracted   | 02/27/2014 19:30 |
| Total Amount Extracted | 10.0 g                    | Analyzed    | 03/04/2014 18:55 |
| ICAL ID                | U140224                   | Injected By | SMT              |
| CCal Filename(s)       | U140304B_02 & U140304B_20 |             |                  |
| Method Blank ID        | BLANK-39477               |             |                  |

| Native Isomers      | Qs (ng) | Qm (ng) | % Rec. | Internal Standards      | ng's Added | Percent Recovery |
|---------------------|---------|---------|--------|-------------------------|------------|------------------|
| 2,3,7,8-TCDF        | 0.20    | 0.36    | 179    | 2,3,7,8-TCDF-13C        | 2.00       | 100              |
|                     |         |         |        | 2,3,7,8-TCDD-13C        | 2.00       | 106              |
|                     |         |         |        | 1,2,3,7,8-PeCDF-13C     | 2.00       | 114              |
| 2,3,7,8-TCDD        | 0.20    | 0.21    | 103    | 2,3,4,7,8-PeCDF-13C     | 2.00       | 88 Y             |
|                     |         |         |        | 1,2,3,7,8-PeCDD-13C     | 2.00       | 121              |
|                     |         |         |        | 1,2,3,4,7,8-HxCDF-13C   | 2.00       | 98               |
| 1,2,3,7,8-PeCDF     | 1.00    | 1.36    | 136    | 1,2,3,6,7,8-HxCDF-13C   | 2.00       | 98               |
|                     |         |         |        | 2,3,4,6,7,8-HxCDF-13C   | 2.00       | 107              |
| 2,3,4,7,8-PeCDF     | 1.00    | 1.47    | 147    | 1,2,3,7,8,9-HxCDF-13C   | 2.00       | 96               |
|                     |         |         |        | 1,2,3,4,7,8-HxCDD-13C   | 2.00       | 88               |
| 1,2,3,7,8-PeCDD     | 1.00    | 1.07    | 107    | 1,2,3,6,7,8-HxCDD-13C   | 2.00       | 84               |
|                     |         |         |        | 1,2,3,4,6,7,8-HpCDF-13C | 2.00       | 82               |
|                     |         |         |        | 1,2,3,4,7,8,9-HpCDF-13C | 2.00       | 84               |
| 1,2,3,4,7,8-HxCDF   | 1.00    | 1.35    | 135    | 1,2,3,4,6,7,8-HpCDD-13C | 2.00       | 93               |
|                     |         |         |        | OCDD-13C                | 4.00       | 78               |
| 1,2,3,6,7,8-HxCDF   | 1.00    | 1.36    | 136    | 1,2,3,4-TCDD-13C        | 2.00       | NA               |
|                     |         |         |        | 2,3,4,6,7,8-HxCDF       | 1.00       | 133              |
| 1,2,3,4,6,7,8-HxCDF | 1.00    | 1.33    | 133    | 1,2,3,7,8,9-HxCDD-13C   | 2.00       | NA               |
|                     |         |         |        | 1,2,3,7,8,9-HxCDF       | 1.00       | 120              |
| 1,2,3,7,8,9-HxCDF   | 1.00    | 1.20    | 120    | 2,3,7,8-TCDD-37Cl4      | 0.20       | 106              |
|                     |         |         |        | 1,2,3,4,7,8-HxCDD       | 1.00       | 124              |
|                     |         |         |        | 1,2,3,6,7,8-HxCDD       | 1.00       | 135              |
| 1,2,3,7,8,9-HxCDD   | 1.00    | 1.32    | 132    |                         |            |                  |
|                     |         |         |        |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDF | 1.00    | 1.92    | 192    |                         |            |                  |
| 1,2,3,4,7,8,9-HpCDF | 1.00    | 1.05    | 105    |                         |            |                  |
| 1,2,3,4,6,7,8-HpCDD | 1.00    | 2.01    | 201    |                         |            |                  |
| OCDF                | 2.00    | 3.01    | 150    |                         |            |                  |
| OCDD                | 2.00    | 9.02    | 451    |                         |            |                  |

Qs = Quantity Spiked                      Qm = Quantity Measured                      Rec. = Recovery (Expressed as Percent)

Results reported on a dry weight basis and are valid to no more than 2 significant figures.

Y = Calculated using average of daily RFs

### REPORT OF LABORATORY ANALYSIS

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### Method 8290 Spike Sample Results

Client - ESN Pacific

|                  |                   |                 |             |                    |        |
|------------------|-------------------|-----------------|-------------|--------------------|--------|
| Client Sample ID | RHS-DU-20         | Sample Filename | U140304B_07 | <u>Dry Weights</u> |        |
| Lab Sample ID    | 10257348006-S     | MS Filename     | U140304B_03 | Sample Amount      | 8.49 g |
| MS ID            | 10257348006-S-MS  | MSD Filename    | U140304B_04 | MS Amount          | 8.9 g  |
| MSD ID           | 10257348006-S-MSD |                 |             | MSD Amount         | 8.5 g  |

| Analyte             | Sample Conc.<br>ng/Kg | MS/MSD Qs<br>(ng) | MS Qm<br>(ng) | MSD Qm<br>(ng) | RPD | Background Subtracted |            |      |
|---------------------|-----------------------|-------------------|---------------|----------------|-----|-----------------------|------------|------|
|                     |                       |                   |               |                |     | MS % Rec.             | MSD % Rec. | RPD  |
| 2,3,7,8-TCDF        | 9.118                 | 0.20              | 0.34          | 0.36           | 5.9 | 129                   | 141        | 9.1  |
| 2,3,7,8-TCDD        | 0.402                 | 0.20              | 0.19          | 0.21           | 8.0 | 94                    | 102        | 8.2  |
| 1,2,3,7,8-PeCDF     | 9.593                 | 1.00              | 1.29          | 1.36           | 5.0 | 120                   | 127        | 5.6  |
| 2,3,4,7,8-PeCDF     | 21.246                | 1.00              | 1.38          | 1.47           | 6.3 | 119                   | 129        | 8.0  |
| 1,2,3,7,8-PeCDD     | 0.000                 | 1.00              | 1.08          | 1.07           | 1.3 | 106                   | 105        | 1.2  |
| 1,2,3,4,7,8-HxCDF   | 16.158                | 1.00              | 1.32          | 1.35           | 2.5 | 117                   | 121        | 3.4  |
| 1,2,3,6,7,8-HxCDF   | 17.658                | 1.00              | 1.32          | 1.36           | 3.1 | 116                   | 121        | 4.1  |
| 2,3,4,6,7,8-HxCDF   | 19.601                | 1.00              | 1.28          | 1.33           | 4.2 | 110                   | 117        | 5.5  |
| 1,2,3,7,8,9-HxCDF   | 3.292                 | 1.00              | 1.20          | 1.20           | 0.2 | 117                   | 117        | 0.1  |
| 1,2,3,4,7,8-HxCDD   | 2.476                 | 1.00              | 1.26          | 1.24           | 2.1 | 124                   | 122        | 2.1  |
| 1,2,3,6,7,8-HxCDD   | 7.186                 | 1.00              | 1.28          | 1.35           | 5.2 | 122                   | 129        | 5.7  |
| 1,2,3,7,8,9-HxCDD   | 4.806                 | 1.00              | 1.32          | 1.32           | 0.6 | 128                   | 127        | 0.4  |
| 1,2,3,4,6,7,8-HpCDF | 72.436                | 1.00              | 1.80          | 1.92           | 6.2 | 116                   | 130        | 11.9 |
| 1,2,3,4,7,8,9-HpCDF | 3.936                 | 1.00              | 1.07          | 1.05           | 2.2 | 104                   | 101        | 2.1  |
| 1,2,3,4,6,7,8-HpCDD | 104.476               | 1.00              | 1.93          | 2.01           | 4.1 | 100                   | 113        | 11.8 |
| OCDF                | 62.070                | 2.00              | 2.87          | 3.01           | 4.6 | 116                   | 124        | 6.7  |
| OCDD                | 667.257               | 2.00              | 8.40          | 9.02           | 7.2 | 122                   | 168        | 31.3 |

**Definitions**

|                                   |                                    |
|-----------------------------------|------------------------------------|
| MS = Matrix Spike                 | CDD = Chlorinated dibenzo-p-dioxin |
| MSD = Matrix Spike Duplicate      | CDF = Chlorinated dibenzo-p-furan  |
| Qm = Quantity Measured            | T = Tetra                          |
| Qs = Quantity Spiked              | Pe = Penta                         |
| % Rec. = Percent Recovery         | Hx = Hexa                          |
| RPD = Relative Percent Difference | Hp = Hepta                         |
| NA = Not Applicable               | O = Octa                           |
| NC = Not Calculated               |                                    |



**APPENDIX D**  
**ASBESTOS DISPOSAL MANIFEST**

Unit 1901SC

87834

**ASBESTOS WASTE SHIPMENT RECORD**

|   |  |               |                    |                                       |  |                    |  |                          |
|---|--|---------------|--------------------|---------------------------------------|--|--------------------|--|--------------------------|
| G<br>E<br>N<br>E<br>R<br>A<br>T<br>O<br>R   | <b>1. Work Site Name and Mailing Address (Generator)</b><br>Radford High School- ACM Disposal<br>4361 Salt Lake Blvd, Honolulu HI 96818  |               |                    |                                       | <b>Owner's Name</b><br>State of Hawaii<br>Dept. of Education |                    | <b>Owners Telephone #</b><br>(808) 543-6132<br>Agent |                          |
|   | <b>2. Remover's Name and Address</b><br>Unitek Insulation, LLC.<br>P.O. Box 29177<br>Honolulu, HI 96820  |               |                    |                                       |  |                    | <b>Remover's Telephone #</b><br>(808) 831-3076       |                          |
|   | <b>3. Waste Disposal site (WDS) Name and Physical Site Location</b><br>PVT Land Company<br>87-2020 Farrington Hwy.<br>Waianae, HI 96792-3749   |               |                    |                                       |  |                    | <b>WDS Telephone No.</b><br>(808) 668-4561           |                          |
|   | <b>4. Name and Address of EPA Office, local, state or regional</b><br>EPA NESHAP Region 9 Noise, Radiation & Indoor Air Branch<br>75 Hawthorne St., A-3-3 or 591 Ala Moana Blvd.<br>San Francisco, CA 94105 Honolulu, HI 96813 |               |                    |                                       |  |                    |  |                          |
| R<br>A<br>T<br>I<br>O<br>N<br>A<br>L  | <b>5. HM</b>   | <b>ID</b>     | <b>Description</b> | <b>Hazard</b>                         | <b>Packing</b>   | <b>6. Packages</b> |  | <b>7. Total Quantity</b> |
|   |  | <b>Number</b> | <b>of Material</b> | <b>Class</b>                          | <b>Group#</b>  | <b>No.</b>         | <b>Type</b>  | <b>m3</b>                |
|   | Asbestos   | NA2212        | ACM Debris         | 9                                     | III  | 5                  | Bags   | ./                       |
| <b>8. Special Handling Instructions and 24 Hour Emergency Telephone Number (provided by Generator)</b><br><br>24 Hour Emergency Contact Person - Darwyn Scott ph: 808-831-3076<br>ERG 171<br><br>GENERATORS' CERTIFICATION: I hereby declare that the contents of this consignment are fully and accurately described above by proper shipping name and are classified, packed, marked, and labeled, and are in all respects in proper condition for transport by highway according to applicable international and government regulations.<br>Note: Generator must retain a copy of this form. |  |               |                    |                                       |  |                    |  |                          |
| T<br>R<br>A<br>N<br>S<br>P<br>O<br>R<br>T<br>E<br>R   | <b>10. TRANSPORTER 1(ACKNOWLEDGEMENT OF RECEIPT OF MATERIALS) Note Transporter must retain a copy of this form.</b>  |               |                    |                                       | <b>Signature (on behalf of)</b>                              |                    | <b>Month/Day/Year</b>                                |                          |
|   | <b>Printed/Typed name and Title</b><br>Sheena Sylva, Waste Manager   |               |                    |                                       | <i>Sheena M. Sylva</i>                                       |                    | 02/19/2014   |                          |
|   | <b>Printed/Typed name and Title</b><br>Unitek Insulation, LLC  |               |                    |                                       | <b>Signature</b><br><i>Mike Ali</i>                          |                    | <b>Month/Day/Year</b><br>2-20-14                     |                          |
| <b>Address</b><br>P.O. Box 29177<br>Honolulu, Hawaii  |  |               |                    | <b>Telephone No.</b><br>(808)831-3076 |  |                    |  |                          |
| P<br>O<br>R<br>T<br>E<br>R  | <b>11. TRANSPORTER 2(ACKNOWLEDGEMENT OF RECEIPT OF MATERIALS) Note Transporter must retain a copy of this form.</b>  |               |                    |                                       |  |                    |  |                          |
|   | <b>Printed/Typed name and Title</b><br>Unitek Insulation, LLC  |               |                    |                                       | <b>Signature</b>   |                    | <b>Month/Day/Year</b>                                |                          |
|   | <b>Address</b><br>P.O. Box 29177<br>Honolulu, Hawaii   |               |                    |                                       | <b>Telephone No.</b><br>(808)831-3076                        |                    |  |                          |
| D<br>I<br>S<br>P<br>O<br>S<br>E<br>R  | <b>12. Problems with Containment or Packaging</b>  |               |                    |                                       |  |                    | <b>Rejected</b><br>yes or no                         |                          |
|   | <b>13. WASTE DISPOSAL SITE OWNER OR OPERATOR</b><br>Certification of receipt of asbestos materials covered by this manifest except as noted in item 12.  |               |                    |                                       |  |                    |  |                          |
|   | <b>Printed/Typed name and Title</b><br>LAHELA ROMENA WEIGHMASTER   |               |                    |                                       | <b>Signature</b><br><i>L. Romena</i>                         |                    | <b>Month/Day/Year</b><br>FEB 20 2014                 |                          |

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