

Final First Quarter 2015 - Quarterly Groundwater Monitoring Report Inside Tunnel Wells

**Red Hill Bulk Fuel Storage Facility
Joint Base Pearl Harbor-Hickam, Oahu, Hawaii**

DOH Facility ID: 9-102271

DOH Release ID: 990051, 010011, 020028, and 140010

March 2015

**Department of the Navy
Naval Facilities Engineering Command, Hawaii
400 Marshall Road
JBPHH HI 96860-3139**



Contract Number N62742-12-D-1853, CTO 0002

This Page Intentionally Left Blank.

Final First Quarter 2015 - Quarterly Groundwater Monitoring Report Inside Tunnel Wells

**Red Hill Bulk Fuel Storage Facility
Joint Base Pearl Harbor-Hickam, Oahu, Hawaii**

DOH Facility ID: 9-102271

DOH Release ID: 990051, 010011, 020028, and 140010

March 2015

Prepared for:



**Department of the Navy
Naval Facilities Engineering Command, Hawaii
400 Marshall Road
JBPHH, HI 96860-3139**

Prepared by:

**Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734**

Prepared under:

Contract Number N62742-12-D-1853, CTO 0002

This Page Intentionally Left Blank.

**FINAL
FIRST QUARTER 2015 - QUARTERLY GROUNDWATER MONITORING REPORT
INSIDE TUNNEL WELLS
RED HILL BULK FUEL STORAGE FACILITY**

Long-Term Groundwater and Soil Vapor Monitoring
Red Hill Bulk Fuel Storage Facility
Joint Base Pearl Harbor-Hickam, Oahu, Hawaii

Prepared for:

Department of the Navy
Commanding Officer, Naval Facilities Engineering Command, Hawaii
400 Marshall Road
JBPHH, HI 96860-3139

Prepared by:

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734
(808) 261-0740

Prepared under:

Contract Number: N62742-12-D-1853
Contract Task Order: 0002

Approval Signature:

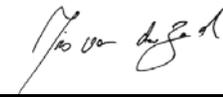


Scott Simmons, ESI Project Manager

3/13/2015

Date

Approval Signature:



Iris van der Zander, ESI QA Manager

3/13/2015

Date

This Page Intentionally Left Blank.

TABLE OF CONTENTS

| <u>Section</u> | <u>Title</u> | <u>Page</u> |
|----------------|--|-------------|
| ES | EXECUTIVE SUMMARY | ES-1 |
| 1.0 | INTRODUCTION | 1-1 |
| 1.1 | Site Description | 1-1 |
| 1.2 | Physical Settings | 1-2 |
| 1.3 | Background | 1-4 |
| 2.0 | GROUNDWATER SAMPLING | 2-1 |
| 2.1 | Groundwater Sampling | 2-1 |
| 2.2 | Analytical Results | 2-2 |
| 2.3 | Groundwater Contaminant Trends | 2-2 |
| 2.4 | Waste Disposal | 2-3 |
| 3.0 | DATA QUALITY ASSESSMENT | 3-1 |
| 3.1 | Data Validation and Assessment | 3-1 |
| 3.2 | Data Assessment and Usability Conclusions | 3-5 |
| 4.0 | SUMMARY, CONCLUSIONS, AND RECOMMENDATIONS | 4-1 |
| 5.0 | FUTURE WORK | 5-1 |
| 6.0 | REFERENCES | 6-1 |

LIST OF TABLES

| <u>Number</u> | <u>Title</u> | <u>Page</u> |
|---------------|--|-------------|
| 1.1 | Current Status of USTs | 1-2 |
| 2.1 | Analytical Results for Groundwater Sampling (January 27 and 28, 2015) | 2-5 |
| 3.1 | Quality Control Results for Groundwater Sampling (January 27 and 28, 2015) | 3-7 |

LIST OF FIGURES

| <u>Number</u> | <u>Title</u> |
|---------------|---------------|
| 1 | Site Location |
| 2 | Site Layout |

APPENDICES

| <u>Appendix</u> | <u>Title</u> |
|------------------------|---|
| A | Groundwater Sampling Logs |
| B | Field Notes |
| C | Laboratory Reports |
| D | Historical Groundwater Exceedance Trends |
| E | Cumulative Groundwater Results (on attached CD) |

ACRONYMS AND ABBREVIATIONS

| ACRONYMS/ ABBREVIATIONS | DEFINITION/MEANING |
|----------------------------|--|
| bgs | below ground surface |
| COPC | Contaminant of Potential Concern |
| DLNR | State of Hawaii Department of Land and Natural Resources |
| DoD | Department of Defense |
| DOH | State of Hawaii Department of Health |
| DON | Department of the Navy |
| EAL | Environmental Action Level |
| EPA | Environmental Protection Agency |
| ESI | Environmental Science International, Inc. |
| F-76 | Marine Diesel Fuel |
| ID | Identification |
| JBPHH | Joint Base Pearl Harbor-Hickam |
| JP-5 | Jet Fuel Propellant-5 |
| JP-8 | Jet Fuel Propellant-8 |
| LCS | Laboratory Control Sample |
| LCSD | Laboratory Control Sample Duplicate |
| LOD | Limit of Detection |
| LOQ | Limit of Quantitation |
| µg/L | micrograms per liter |
| MS | Matrix Spike |
| MSD | Matrix Spike Duplicate |
| NAVFAC | Naval Facilities Engineering Command |
| NAVSUP FLC | Naval Supply Systems Command Fleet Logistics Center |
| PAH | Polycyclic Aromatic Hydrocarbons |
| PARCCS | Precision, Accuracy, Representativeness, Completeness, Comparability, and Sensitivity |
| pH | hydrogen activity |
| RHSF | Red Hill Bulk Fuel Storage Facility |
| RPD | Relative Percent Difference |
| SAP | Sampling and Analysis Plan |
| SSRBL | Site-Specific Risk-Based Level |
| TEC | The Environmental Company, Inc. |
| TPH | Total Petroleum Hydrocarbons |
| TPH-d | Total Petroleum Hydrocarbons as diesel fuel |
| TPH-g | Total Petroleum Hydrocarbons as gasoline |
| U.S. | United States of America |
| UST | Underground Storage Tank |
| VOC | Volatile Organic Compound |
| WP | Work Plan |

This Page Intentionally Left Blank.

EXECUTIVE SUMMARY

This quarterly groundwater monitoring report presents the results of the first quarter 2015 groundwater sampling event, conducted on January 27 and 28, 2015, at the Red Hill Bulk Fuel Storage Facility [RHSF], Joint Base Pearl Harbor-Hickam [JBPHH], Hawaii. The RHSF is located in Halawa Heights on the Island of Oahu. There are 18 active and 2 inactive Underground Storage Tanks [USTs] located at the RHSF. The State of Hawaii Department of Health [DOH] Facility Identification [ID] number is 9-102271. The DOH Release ID numbers are 990051, 010011, 020028, and 140010.

The groundwater sampling was conducted as part of the long-term groundwater and soil vapor monitoring program at the RHSF and concurrent with release response activities initiated at Tank 5 in January 2014, under Naval Facilities Engineering Command [NAVFAC] Contract Number N62742-12-D-1853. The sampling was conducted in accordance with the approved 2012 Work Plan [WP]/Sampling and Analysis Plan [SAP] prepared by Environmental Science International, Inc. [ESI].

On January 27 and 28, 2015, ESI personnel collected groundwater samples from four monitoring wells at the RHSF (wells RHMW01, RHMW02, RHMW03, and RHMW05) and one sampling point at Red Hill Shaft (RHMW2254-01). A primary and duplicate sample were collected from well RHMW02. A summary of the analytical results is provided below.

- **RHMW01** – Total Petroleum Hydrocarbons as diesel fuel [TPH-d] (33 micrograms per liter [$\mu\text{g/L}$]) and lead (0.631 $\mu\text{g/L}$) were the only analytes detected. The concentrations did not exceed the DOH Environmental Action Level [EAL] or the site-specific risk-based level [SSRBL] of 4,500 $\mu\text{g/L}$ for Total Petroleum Hydrocarbons [TPH]. TPH-d concentrations in this well have shown an overall decreasing trend from a high of 1,500 $\mu\text{g/L}$ in February 2005.
- **RHMW02** – TPH-d (1,100 and 1,700 $\mu\text{g/L}$), Total Petroleum Hydrocarbons as gasoline [TPH-g] (54 and 59 $\mu\text{g/L}$), xylenes (0.35 $\mu\text{g/L}$ in both primary and duplicate samples), acenaphthene (0.59 and 0.55 $\mu\text{g/L}$), ethylbenzene (0.16 and 0.17 $\mu\text{g/L}$), fluorene (0.30 and 0.22 $\mu\text{g/L}$), 1-methylnaphthalene (34 and 25 $\mu\text{g/L}$), 2-methylnaphthalene (7.6 and 2.7 $\mu\text{g/L}$), and naphthalene (90 and 63 $\mu\text{g/L}$) were detected. TPH-d was detected at concentrations above the DOH EALs for both drinking water toxicity and gross contamination, but below half the SSRBL. The polycyclic aromatic hydrocarbons [PAHs] 1-methylnaphthalene and naphthalene were detected at concentrations above the DOH EALs for both drinking water toxicity and gross contamination. The concentrations of TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene decreased from the previous sampling event in October 2014.
- **RHMW03** – TPH-d (39 $\mu\text{g/L}$) was the only analyte detected. The concentration did not exceed the DOH EALs or the SSRBL.
- **RHMW05** – None of the chemical constituents analyzed for were detected at a concentration at or above the limit of detection [LOD].

- **RHMW2254-01** – None of the chemical constituents analyzed for were detected at a concentration at or above the LOD.

During this quarterly event, the concentrations of TPH-d, 1-methylnaphthalene, and naphthalene in RHMW02 exceeded the DOH EALs for both gross contamination and drinking water toxicity. The concentration of TPH-d in RHMW01 decreased from the previous event in October 2014 to a concentration below the DOH EALs. Groundwater contaminant concentrations in RHMW03, RHMW05, and RHMW2254-01 remained at low concentrations and did not change significantly from the previous event, or were not detected.

Concentrations of 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene in RHMW02 had shown an increasing trend between March and October 2014; however, concentrations of these three analytes and TPH-d all decreased since the previous event in October 2014. All other analytical results were generally consistent with historical data.

Based on the groundwater monitoring results and the reported release at Tank 5 in January 2014, continued groundwater monitoring at the wells inside the RHSF tunnel is recommended. The next quarterly event is tentatively scheduled for April 2015.

SECTION 1 – INTRODUCTION

This quarterly groundwater monitoring report presents the results of the first quarter 2015 groundwater sampling conducted on January 27 and 28, 2015, at the RHSF, JBPHH. The RHSF is located in Halawa Heights on the Island of Oahu. The purpose of the sampling is to (1) assess the condition of groundwater beneath the RHSF with respect to chemical constituents associated with jet fuel propellant and marine diesel fuel, and (2) to ensure the Navy remains in compliance with DOH UST release response requirements as described in Hawaii Administrative Rules Chapter 11-281 Subchapter 7, *Release Response Action* (DOH, 2013). The DOH Facility ID number for the RHSF is 9-102271. The DOH Release ID numbers are 990051, 010011, 020028, and 140010.

The groundwater sampling was conducted as part of the long-term groundwater and soil vapor monitoring program at the RHSF, under NAVFAC Contract Number N62742-12-D-1853. The sampling was conducted in accordance with the approved WP/SAP prepared by ESI (ESI, 2012).

1.1 SITE DESCRIPTION

The RHSF is located on federal government land (zoned F1- Military and Federal), located in Halawa Heights, approximately 2.5 miles northeast of Pearl Harbor (Figure 1). It is located on a low ridge on the western edge of the Koolau Mountain Range that divides Halawa Valley from Moanalua Valley. The RHSF is bordered on the north by Halawa Correctional Facility and private businesses, on the southwest by the United States of America [U.S.] Coast Guard reservation, on the south by residential neighborhoods, and on the east by Moanalua Valley. A quarry is located less than a quarter mile away to the northwest. The RHSF occupies 144 acres of land and the majority of the site is at an elevation of approximately 200 to 500 feet above mean sea level.

The RHSF contains 18 active and 2 inactive USTs that are operated by Naval Supply Systems Command Fleet Logistics Center [NAVSUP FLC] Pearl Harbor (formerly Fleet and Industrial Supply Center). Each UST has a capacity of approximately 12.5 million gallons. The RHSF is located approximately 100 feet above the basal aquifer. The USTs contain Jet Fuel Propellant-5 [JP-5], Jet Fuel Propellant-8 [JP-8], and Marine Diesel Fuel [F-76]. The current status of the USTs are summarized in Table 1.1.

Four groundwater monitoring wells (wells RHMW01, RHMW02, RHMW03, and RHMW05) and one sampling point at Red Hill Shaft (RHMW2254-01) are located within the RHSF lower access tunnel (Figure 2). Three groundwater monitoring wells (wells HDMW2253-03, OWDFMW01, and RHMW04) are located outside of the RHSF tunnel system. Monitoring data for the three wells located outside the tunnel are included in a separate report.

As noted, monitoring wells RHMW01, RHMW02, RHMW03, and RHMW05 are located inside the underground tunnels. Sampling point RHMW2254-01 is located inside the infiltration gallery of the Department of the Navy [DON] drinking water supply Well 2254-01. The DON Well 2254-01 is located approximately 2,400 feet downgradient of the USTs and provides potable water to the JBPHH Water System, which serves approximately 65,200 military customers. NAVFAC Public Works Department operates the infiltration gallery and DON Well 2254-01.

TABLE 1.1
Current Status of the USTs
Red Hill Bulk Fuel Storage Facility
January 2015 Quarterly Monitoring Report

| Tank Identification | Fuel Type | Status | Capacity |
|---------------------|-----------|----------|----------------------|
| F-1 | None | Inactive | 12.5 million gallons |
| F-2 | JP-8 | Active | 12.5 million gallons |
| F-3 | JP-8 | Active | 12.5 million gallons |
| F-4 | JP-8 | Active | 12.5 million gallons |
| F-5 | JP-8 | Active | 12.5 million gallons |
| F-6 | JP-8 | Active | 12.5 million gallons |
| F-7 | JP-5 | Active | 12.5 million gallons |
| F-8 | JP-5 | Active | 12.5 million gallons |
| F-9 | JP-5 | Active | 12.5 million gallons |
| F-10 | JP-5 | Active | 12.5 million gallons |
| F-11 | JP-5 | Active | 12.5 million gallons |
| F-12 | JP-5 | Active | 12.5 million gallons |
| F-13 | F-76 | Active | 12.5 million gallons |
| F-14 | F-76 | Active | 12.5 million gallons |
| F-15 | F-76 | Active | 12.5 million gallons |
| F-16 | F-76 | Active | 12.5 million gallons |
| F-17 | JP-5 | Active | 12.5 million gallons |
| F-18 | JP-5 | Active | 12.5 million gallons |
| F-19 | None | Inactive | 12.5 million gallons |
| F-20 | JP-5 | Active | 12.5 million gallons |

F-76 Marine Diesel Fuel
JP-5 Jet Fuel Propellant-5
JP-8 Jet Fuel Propellant-8

1.2 PHYSICAL SETTING

Climatological conditions in the area of the RHSF consist of warm to moderate temperatures and low to moderate rainfall. The RHSF is leeward of the prevailing northeasterly trade winds. The average annual precipitation is approximately 40 inches, which occurs mainly between November and April (State of Hawaii Department of Land and Natural Resources [DLNR], 1986). Annual pan evaporation is approximately 75 inches (DLNR, 1985). Average temperatures range from the low 60's to high 80's (degrees Fahrenheit) (Atlas of Hawaii, 1983).

Oahu consists of the eroded remnants of two shield volcanoes, Waianae and Koolau. The RHSF is located on the southwest flank of the Koolau volcanic shield. Lavas erupted during the shield-building phase of the volcano belong to the *Koolau Volcanic Series* (Stearns and Vaksvik,

1935). Following formation of the Koolau shield, a long period of volcanic quiescence occurred, during which the shield was deeply eroded. Following this erosional period, eruptive activity resumed. Lavas and pyroclastic material erupted during this period belong to the *Honolulu Volcanic Series* (Stearns and Vaksvik, 1935).

In the immediate area of the RHSF, Koolau Volcanic Series lavas dominate, although there are consolidated and unconsolidated non-calcareous deposits in the vicinity that consist of alluvium generated during erosion of the Koolau volcanic shield. South-southwest of the RHSF, and in isolated exposures to the west, are pyroclastic deposits formed during eruptions from three Honolulu Volcanic Series vents, Salt Lake, Aliamanu, and Makalapa (Stearns and Vaksvik, 1935). Based on established geology and records of wells drilled at the RHSF (Stearns and Vaksvik, 1938), the RHSF is underlain by Koolau Volcanic Series basalts. The area of the RHSF is classified as *Rock Land*, where 25-90% of the land surface is covered by exposed rock and there are only shallow soils (Foote, et al., 1972).

Groundwater in Hawaii exists in two principal aquifer types. The first and most important type, in terms of drinking water resources, is the basal aquifer. The basal aquifer exists as a lens of fresh water floating on and displacing seawater within the pore spaces, fractures, and voids of the basalt that forms the underlying mass of each Hawaiian island. In parts of Oahu, groundwater in the basal aquifer is confined by the overlying caprock and is under pressure. Waters that flow freely to the surface from wells that tap the basal aquifer are referred to as *artesian*.

The second type of aquifer is the caprock aquifer, which consists of various kinds of unconfined and semi-confined groundwater. Commonly, the caprock consists of a thick sequence of nearly impermeable clays, coral, and basalt, that separates the caprock aquifer from the basal aquifer. The impermeable nature of these materials and the artesian nature of the basal aquifer severely restrict the downward migration of groundwater from the upper caprock aquifer. However, in the area of the RHSF, there is no discernible caprock.

Groundwater in the area of the RHSF is part of the *Waimalu Aquifer System* of the *Pearl Harbor Aquifer Sector*. The aquifer is classified as a basal, unconfined, flank-type; and is currently used as a drinking water source. The aquifer is considered fresh, with less than 250 milligrams per liter of chloride, and is considered an irreplaceable resource with a high vulnerability to contamination (Mink and Lau, 1990).

The nearest drinking water supply well is DON Well 2254-01, located in the infiltration gallery within the RHSF lower tunnel. The DON Well 2254-01 is located approximately 2,400 feet hydraulically and topographically downgradient of the USTs (Figure 2).

1.3 BACKGROUND

The RHSF was constructed by the U.S. Government in the early 1940s. Twenty USTs and a series of tunnels were constructed. The USTs were constructed of steel, and in the past have stored DON special fuel oil, DON distillate, aviation gasoline, and motor gasoline (Environet, 2010). The tanks currently contain JP-5, JP-8, and F-76. The fueling system is a self-contained underground unit that was installed into native rock comprised primarily of basalt with some interbedded tuffs and breccias (Environet, 2010). Each UST measures approximately 250 feet in height and 100 feet in diameter. The upper domes of the tanks lie at a depth varying between 100 feet and 200 feet below ground surface [bgs].

In 1998, Earth Tech conducted a Phase II Remedial Investigation/Feasibility Study for the Oily Waste Disposal Facility located within the RHSF. The study included the installation of well OWDFMW01 (which was originally MW08) (Earth Tech, 1999).

In February 2001, the DON installed groundwater monitoring well RHMW01 to monitor for contamination in the basal aquifer beneath the RHSF. Well RHMW01 was installed approximately 100 feet below grade within the lower access tunnel. The depth to water was measured at 86 feet below the tunnel floor at the time of the well completion. In February 2001, a groundwater sample was collected from the well. TPH and total lead were detected in the sample. Total lead was detected at a concentration above the DOH Tier 1 groundwater action level of 5.6 µg/L (The Environmental Company, Inc. [TEC], 2009; DOH, 2000).

In 2005, the RHSF groundwater monitoring program was initiated. It involved routine groundwater sampling of well RHMW01 and sampling point RHMW2254-01. Samples were collected in February, June, September, and December of 2005. Lead was detected at concentrations above the DOH Tier 1 action level of 5.6 µg/L in samples collected in February and June. The samples collected in February and June were not filtered prior to analysis, whereas the samples collected in September and December were filtered prior to analysis. Since the samples collected in February and June were not filtered prior to analysis, the lead results were not considered appropriate for a risk assessment (TEC, 2008).

Between June and September 2005, TEC installed three additional groundwater monitoring wells (wells RHMW02, RHMW03, and RHMW04) (TEC, 2008). Well RHMW04 was installed hydraulically upgradient of the USTs to provide background geochemistry information for water moving through the basal aquifer beneath the RHSF. Wells RHMW02 and RHMW03 were installed approximately 125 feet below grade within the RHSF lower tunnel and well RHMW04 was installed to a depth of approximately 300 feet bgs outside of the RHSF tunnels. In September 2005, groundwater samples were collected from the three newly installed groundwater monitoring wells (wells RHMW02, RHMW03, and RHMW04) along with the existing well RHMW01 and sampling point RHMW2254-01. The contaminants of potential concern [COPCs] with concentrations exceeding current DOH EALs are summarized below.

- **RHWM01** – TPH-d was detected at concentrations above the DOH EALs.
- **RHWM02** – TPH-g, TPH-d, naphthalene, trichloroethylene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above their respective DOH EAL.
- **RHWM03** – TPH-d was detected at concentrations above the DOH EALs.

In 2006, TEC installed dedicated sampling pumps in the four wells (wells RHWM01, RHWM02, RHWM03, and RHWM04) and one sampling point (RHWM2254-01). In July and December of 2006, groundwater samples were collected from the four wells and the sampling point. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- **RHWM01** – TPH-d and naphthalene were detected at concentrations above their respective DOH EAL.
- **RHWM02** – TPH-g, TPH-d, and naphthalene were detected at concentrations above their respective DOH EAL.
- **RHWM03** – TPH-d was detected at concentrations above the DOH EALs.

In 2007, SSRBLs were established for TPH (4,500 µg/L) and benzene (750 µg/L) based on the solubility in water of JP-5 and JP-8 (TEC, 2007). Groundwater samples were collected from wells RHWM01, RHWM02, and RHWM03, and sampling point RHWM2254-01. Samples were collected in March, June, and September of 2007. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- **RHWM01** – TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- **RHWM02** – TPH-g, TPH-d, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.
- **RHWM03** – TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.

In 2008, groundwater samples were collected from wells RHWM01, RHWM02, and RHWM03, and sampling point RHWM2254-01. Samples were collected in January, April, July, and October of 2008. The COPCs with concentrations exceeding current DOH EALs are summarized below. In addition, a groundwater protection plan (TEC, 2008) was prepared.

- **RHWM01** – TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- **RHWM02** – TPH-d, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations detected in October 2008 were also above the SSRBL.

- **RHMW03** – TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.

In April 2009, groundwater monitoring well RHMW05 was installed downgradient of the USTs, within the lower access tunnel between RHMW01 and RHMW2254-01. It was installed to identify the extent of contamination hydraulically downgradient of the USTs. Well RHMW05 was added to the quarterly groundwater sampling program. In 2009, quarterly groundwater samples were collected from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in February, May, July, and October of 2009. The COPCs with concentrations exceeding current DOH EALs are summarized below. In addition, the Groundwater Protection Plan was revised to include well RHMW05.

- **RHMW01** – TPH-d and 1-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.
- **RHMW02** – TPH-d, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.
- **RHMW03** – TPH-d was detected at a concentration above the DOH EALs, but below the SSRBL.
- **RHMW05** – TPH-d was detected at a concentration above the DOH EALs, but below the SSRBL.

In 2010, groundwater samples were collected from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in January, April, July, and October. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- **RHMW01** – TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- **RHMW02** – TPH-g, TPH-d, naphthalene, and 1-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.
- **RHMW03** – TPH-d was detected at a concentration above the DOH EALs, but below the SSRBL.
- **RHMW05** – TPH-d was detected at a concentration above the DOH EALs, but below the SSRBL.

In 2011, quarterly groundwater samples were collected from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in January, April, July, and October. In a Fall 2011 update, the DOH EALs were revised. The drinking water toxicity EAL for TPH-d decreased from 210 to 190 µg/L (DOH, 2011). The COPCs with concentrations exceeding current DOH EALs are summarized below.

- **RHMW01** – TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- **RHMW02** – TPH-d, naphthalene, indeno[1,2,3-cd]pyrene, and 1-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.

In 2012, quarterly groundwater samples were collected from wells RHMW01, RHMW02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in February, April, July, and November. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- **RHMW01** – TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- **RHMW02** – TPH-d, TPH-g, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.

In 2013, quarterly groundwater samples were collected from wells RHMW01, RHMW02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in January, April, July, and October. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- **RHMW01** – TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- **RHMW02** – TPH-d, TPH-g, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.

In 2014, quarterly groundwater samples were collected from wells RHMW01, RHMW02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in January, April, July, and October. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- **RHMW01** – TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- **RHMW02** – TPH-d, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.

Between January and June 2014, additional groundwater sampling was conducted at wells RHMW01, RHMW02, RHMW05, and sampling point RHMW2254-01 in response to a reported

release from Tank 5. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- **RHMW01** – TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- **RHMW02** – TPH-d, 1-methylnaphthalene, and naphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.

1.3.1 Previous Reports

The following groundwater monitoring reports were previously submitted to the DOH:

1. Groundwater Sampling Report, First Quarter 2005 (submitted April 2005).
2. Groundwater Sampling Report, Second Quarter 2005 (submitted August 2005).
3. Groundwater Sampling Report, Third Quarter 2005 (submitted November 2005).
4. Groundwater Sampling Report, Fourth Quarter 2005 (submitted February 2006).
5. Groundwater Monitoring Results, July 2006 (submitted September 2006).
6. Groundwater Monitoring Results, December 2006 (submitted January 2007).
7. Groundwater Monitoring Results, March 2007 (submitted May 2007).
8. Groundwater Monitoring Results, June 2007 (submitted August 2007).
9. Groundwater Monitoring Results, September 2007 (submitted October 2007).
10. Groundwater Monitoring Report, January 2008 (submitted March 2008).
11. Groundwater Monitoring Report, April 2008 (submitted May 2008).
12. Groundwater Monitoring Report, July 2008 (submitted October 2008).
13. Groundwater Monitoring Report, October and December 2008 (submitted February 2009).
14. Groundwater Monitoring Report, February 2009 (submitted May 2009).
15. Groundwater Monitoring Report, May 2009 (submitted July 2009).
16. Groundwater Monitoring Report, July 2009 (submitted September 2009).
17. Groundwater Monitoring Report, October 2009 (submitted December 2009).
18. Groundwater Monitoring Report, January, February, and March 2010 (submitted April 2010).
19. Groundwater Monitoring Report, April 2010 (submitted May 2010).
20. Groundwater Monitoring Report, July 2010 (submitted August 2010).
21. Groundwater Monitoring Report, October 2010 (submitted December 2010).
22. Groundwater Monitoring Report, January 2011 (submitted March 2011).
23. Groundwater Monitoring Report, April 2011 (submitted June 2011).

24. Groundwater Monitoring Report, July 2011 (submitted September 2011).
25. Groundwater Monitoring Report, October 2011 (submitted December 2011).
26. Groundwater Monitoring Report, January-February 2012 (submitted March 2012).
27. Groundwater Monitoring Report, April 2012 (Submitted July 2012).
28. Groundwater Monitoring Report, October 2012 (Submitted January 2013).
29. Groundwater Monitoring Report, January 2013 (Submitted April 2013).
30. Groundwater Monitoring Report, April 2013 (Submitted July 2013).
31. Groundwater Monitoring Report, July 2013 (Submitted September 2013).
32. Groundwater Monitoring Report, October 2013 (Submitted January 2014).
33. Groundwater Sampling Report for Additional Sampling, January 2014 (submitted January 2014).
34. Groundwater Monitoring Report, January 2014 (Submitted April 2014).
35. Groundwater Sampling Report for Tank 5 Release Response on March 5 and 6, 2014 (submitted March 2014).
36. Groundwater Sampling Report for Tank 5 Release Response on March 10, 2014 (submitted March 2014).
37. Groundwater Sampling Report for Tank 5 Release Response on March 25 and 26, 2014 (submitted April 2014).
38. Groundwater Sampling Report for Tank 5 Release Response on April 7, 2014 (submitted April 2014).
39. Groundwater Monitoring Report, April 2014 (Submitted June 2014).
40. Groundwater Sampling Report for Tank 5 Release Response on May 27 and 28, 2014 (submitted June 2014).
41. Groundwater Sampling Report for Tank 5 Release Response on June 23 and 24, 2014 (submitted July 2014).
42. Groundwater Monitoring Report, July 2014 (Submitted September 2014).
43. Groundwater Monitoring Report, October 2014 (Submitted January 2015).

This Page Intentionally Left Blank.

SECTION 2 – GROUNDWATER SAMPLING

On January 27 and 28, 2015, ESI personnel collected groundwater samples from four monitoring wells at the RHSF (wells RHMW01, RHMW02, RHMW03, and RHMW05) and one sampling point at Red Hill Shaft (RHMW2254-01). The samples were collected in accordance with the 2012 WP/SAP. The WP/SAP is consistent with DOH UST release response requirements (DOH, 2000); DON Procedure I-C-3, *Monitoring Well Sampling* (DON, 2007); and the RHSF Final Groundwater Protection Plan (TEC, 2008). Prior to purging and sampling, the depth to groundwater and the depth to the bottoms of the wells were measured using a Geotech oil/water interface probe. No measurable product, sheen, or petroleum hydrocarbon odor was detected in any of the wells.

2.1 GROUNDWATER SAMPLING

Prior to collecting groundwater samples, the monitoring wells were purged of standing water in the well casings. Each well contains a dedicated bladder pump which was used to purge the well and to collect samples. To operate the pump, a portable air compressor with an in-line filter was connected to a QED MP50 MicroPurge[®] Basics Controller box, which was then connected to the pump. The compressor was turned on to power the pump and the controller was used to adjust the pumping rate to less than one liter of water per minute.

Water quality parameters were monitored on a periodic basis during well purging. The water quality parameters that were measured included hydrogen activity [pH], temperature, conductivity, dissolved oxygen, and oxidation reduction potential. The water quality parameters were evaluated to assess whether the natural characteristics of the aquifer formation water were present within the monitoring wells before collecting the samples. At least four readings were collected during the purging process. Purging was considered complete when at least three consecutive water quality measurements stabilized within approximately 10%. The readings were recorded on groundwater monitoring logs. The groundwater monitoring logs are included in Appendix A. In addition, field notes were taken to document the sampling event. The field notes are included in Appendix B.

When the water quality parameters stabilized, groundwater samples were collected from the wells using the bladder pumps. The groundwater samples were collected no more than two hours after purging was completed to decrease groundwater interaction with the monitoring well casing and atmosphere. Prior to collecting the sample, the water level in the monitoring wells was measured and recorded to ensure that excessive drawn down had not occurred. The groundwater samples were collected at flow rates of approximately 0.17 to 0.5 liters per minute. Samples collected for dissolved lead analysis were filtered in the field using new, 0.45-micron filters.

All samples were labeled and logged on the Sample Inventory Log, placed in Ziploc[™] bags and sealed, custody sealed, sealed with tape, placed in a cooler with wet ice, and logged onto the Chain-of-Custody form. The samples were labeled and logged in accordance with DON

Procedure III-E, *Record Keeping, Sample Labeling, and Chain-of-Custody Procedures* (DON, 2007). All samples were shipped under Chain-of-Custody to the analytical laboratory and analyzed for the COPCs as described in Section 2.2.

2.2 ANALYTICAL RESULTS

The samples were analyzed for TPH-d using U.S. Environmental Protection Agency [EPA] Method 8015M, TPH-g and volatile organic compounds [VOCs] using EPA Method 8260B, PAHs using EPA Method 8270C SIM, dissolved lead using EPA Method 6020, and total lead using EPA Method 200.8. The sample collected from sampling point RHMW2254-01 was analyzed for total lead (unfiltered) as the sampling point is a drinking water supply infiltration shaft. A copy of the laboratory report is included as Appendix C.

Analytical results were compared to the DOH EALs for drinking water toxicity and gross contamination. Analytical results for wells RHMW01, RHMW02, RHMW03, and RHMW05 were also compared to the SSRBLs for TPH (4,500 µg/L) and benzene (750 µg/L), established in the RHSF Final Groundwater Protection Plan (TEC, 2008). The results of the first quarter groundwater sampling event are summarized in Table 2.1 and described below.

- **RHMW01** – TPH-d (33 µg/L) and lead (0.0631 µg/L) were the only analytes detected. The concentrations did not exceed the DOH EALs or the SSRBL.
- **RHMW02** – TPH-d (1,100 and 1,700 µg/L), TPH-g (54 and 59 µg/L), xylenes (0.35 µg/L in both primary and duplicate samples), acenaphthene (0.59 and 0.55 µg/L), ethylbenzene (0.16 and 0.17 µg/L), fluorene (0.30 and 0.22 µg/L), 1-methylnaphthalene (34 and 25 µg/L), 2-methylnaphthalene (7.6 and 2.7 µg/L), and naphthalene (90 and 63 µg/L) were detected in both the primary and duplicate samples collected. TPH-d, 1-methylnaphthalene, and naphthalene were detected at concentrations above their respective DOH EALs for both drinking water toxicity and gross contamination. However, the TPH-d concentrations did not exceed the SSRBL.
- **RHMW03** – TPH-d (39 µg/L) was the only analyte detected. The concentration did not exceed the DOH EALs or the SSRBL.
- **RHMW05** – None of the chemical constituents analyzed for were detected at a concentration at or above the LOD.
- **RHMW2254-01** – None of the chemical constituents analyzed for were detected at a concentration at or above the LOD.

2.3 GROUNDWATER CONTAMINANT TRENDS

The historical groundwater contaminant concentration trends for COPCs that exceeded the DOH EALs or SSRBLs are illustrated in Appendix D. A table of cumulative historical groundwater results is included as Appendix E. A summary of groundwater contaminant trends is provided below.

- **RHMW01** – COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW01. TPH-d has historically been detected at concentrations above the DOH EAL for both drinking water toxicity and gross contamination. TPH-d concentrations have shown an overall decreasing trend from a high of 1,500 µg/L in February 2005.
- **RHMW02** – COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW02. TPH-g, TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene have historically been detected at concentrations above the DOH EALs. During the January 2015 event, concentrations of TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene decreased from the previous event in October 2014, with the concentrations of 2-methylnaphthalene decreasing to levels below the DOH EALs. The concentrations of TPH-g remained below the DOH EALs for gross contamination and drinking water toxicity and were comparable to the concentrations detected during the previous event. Trichloroethylene was detected once in RHMW02 in September 2005 in the primary sample at a concentration above the DOH EAL for drinking water toxicity; however, trichloroethylene was not detected in the duplicate sample, and this may have been an anomalous result.
- **RHMW03** – COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW03. TPH-d has historically been detected at concentrations above the DOH EALs; however, it has not been detected at concentrations above the DOH EALs since October 2010.
- **RHMW05** – COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW05. TPH-d has historically been detected in RHMW05 at concentrations above the DOH EAL for both drinking water toxicity and gross contamination; however, it has not been detected at concentrations above the DOH EALs since January 2010.
- **RHMW2254-01** – COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW2254-01. Although the method reporting limits for TPH-d exceeded one or both DOH EALs for drinking water toxicity and gross contamination between May 2009 and July 2010, TPH-d was last detected in RHMW2254-01 at a concentration above the DOH EAL for gross contamination in January 2008.

2.4 WASTE DISPOSAL

The purged groundwater and decontamination water generated during sampling of the inside tunnel were placed in a 55-gallon drum along with the purged water and decontamination water from the outside tunnel wells. The drum is currently stored onsite at ADIT 3 on top of a secondary containment spill pallet and covered by a tarp. There is a non-hazardous label affixed to the drum with all pertinent information relating to its generation. The drum is nearly full and is currently awaiting disposal.

This Page Intentionally Left Blank.

TABLE 2.1
Analytical Results for Groundwater Sampling (January 27 and 28, 2015)
Red Hill Bulk Fuel Storage Facility
January 2015 Quarterly Monitoring Report

| Method | Chemical | DOH EALs | | RHMW2254-01 (ES125) | | | | | RHMW01 (ES120X) | | | | | RHMW02 (ES126) | | | | | RHMW03 (ES123) | | | | | RHMW05 (ES124) | | | | | |
|---|-------------------------------------|--|---------------------|---------------------|-------|------|-------|-------|-----------------|-------|------|-------|-------|----------------|-------|------|-------|-------|----------------|-------|------|-------|-------|----------------|-------|------|-------|-------|------|
| | | Drinking Water Toxicity | Gross Contamination | Results | Q | LOQ | LOD | DL | Results | Q | LOQ | LOD | DL | Results | Q | LOQ | LOD | DL | Results | Q | LOQ | LOD | DL | Results | Q | LOQ | LOD | DL | |
| EPA 8015B | TPH-d | 190 | 100 | N.D. | U | 25 | 12 | 11 | 33 | HD | 25 | 12 | 11 | 1,100 | HD | 25 | 10 | 2.9 | 39 | HD | 25 | 10 | 2.9 | N.D. | U | 26 | 13 | 12 | |
| EPA 8260B | TPH-g | 100 | 100 | N.D. | U | 50 | 30 | 26 | N.D. | U | 50 | 30 | 26 | 54 | | 50 | 30 | 26 | N.D. | U | 50 | 30 | 26 | N.D. | U | 50 | 30 | 26 | |
| EPA 8270C | Acenaphthene | 370 | 20 | N.D. | U | 0.20 | 0.050 | 0.027 | N.D. | U | 0.22 | 0.054 | 0.029 | 0.59 | | 0.20 | 0.050 | 0.027 | N.D. | U | 0.19 | 0.048 | 0.026 | N.D. | U | 0.19 | 0.048 | 0.026 | |
| | Acenaphthylene | 240 | 2,000 | N.D. | U | 0.20 | 0.050 | 0.044 | N.D. | U | 0.22 | 0.054 | 0.048 | N.D. | U | 0.20 | 0.050 | 0.044 | N.D. | U | 0.19 | 0.048 | 0.043 | N.D. | U | 0.19 | 0.048 | 0.043 | |
| | Anthracene | 1,800 | 22 | N.D. | U | 0.20 | 0.050 | 0.029 | N.D. | U | 0.22 | 0.054 | 0.031 | N.D. | U | 0.20 | 0.050 | 0.029 | N.D. | U | 0.19 | 0.048 | 0.028 | N.D. | U | 0.19 | 0.048 | 0.028 | |
| | Benzo[a]anthracene | 0.092 | 4.7 | N.D. | U | 0.20 | 0.050 | 0.033 | N.D. | U | 0.22 | 0.054 | 0.035 | N.D. | U | 0.20 | 0.050 | 0.032 | N.D. | U | 0.19 | 0.048 | 0.032 | N.D. | U | 0.19 | 0.048 | 0.031 | |
| | Benzo[g,h,i]perylene | 1,500 | 0.13 | N.D. | U | 0.20 | 0.10 | 0.082 | N.D. | U | 0.22 | 0.11 | 0.089 | N.D. | U | 0.20 | 0.099 | 0.081 | N.D. | U | 0.19 | 0.097 | 0.079 | N.D. | U | 0.19 | 0.096 | 0.079 | |
| | Benzo[a]pyrene | 0.2 | 0.81 | N.D. | U | 0.20 | 0.050 | 0.022 | N.D. | U | 0.22 | 0.054 | 0.024 | N.D. | U | 0.20 | 0.050 | 0.022 | N.D. | U | 0.19 | 0.048 | 0.022 | N.D. | U | 0.19 | 0.048 | 0.021 | |
| | Benzo[b]fluoranthene | 0.092 | 0.75 | N.D. | U | 0.20 | 0.050 | 0.018 | N.D. | U | 0.22 | 0.054 | 0.019 | N.D. | U | 0.20 | 0.050 | 0.017 | N.D. | U | 0.19 | 0.048 | 0.017 | N.D. | U | 0.19 | 0.048 | 0.017 | |
| | Benzo[k]fluoranthene | 0.92 | 0.4 | N.D. | U | 0.20 | 0.050 | 0.031 | N.D. | U | 0.22 | 0.054 | 0.034 | N.D. | U | 0.20 | 0.050 | 0.031 | N.D. | U | 0.19 | 0.048 | 0.030 | N.D. | U | 0.19 | 0.048 | 0.030 | |
| | Chrysene | 9.2 | 1 | N.D. | U | 0.20 | 0.050 | 0.025 | N.D. | U | 0.22 | 0.054 | 0.027 | N.D. | U | 0.20 | 0.050 | 0.025 | N.D. | U | 0.19 | 0.048 | 0.024 | N.D. | U | 0.19 | 0.048 | 0.024 | |
| | Dibenzo[a,h]anthracene ¹ | 0.0092 | 0.52 | N.D. | U | 0.20 | 0.050 | 0.047 | N.D. | U | 0.22 | 0.054 | 0.052 | N.D. | U | 0.20 | 0.050 | 0.047 | N.D. | U | 0.19 | 0.048 | 0.046 | N.D. | U | 0.19 | 0.048 | 0.046 | |
| | Fluoranthene | 1,500 | 130 | N.D. | U | 0.20 | 0.050 | 0.047 | N.D. | U | 0.22 | 0.054 | 0.051 | N.D. | U | 0.20 | 0.050 | 0.046 | N.D. | U | 0.19 | 0.048 | 0.045 | N.D. | U | 0.19 | 0.048 | 0.045 | |
| | Fluorene | 240 | 950 | N.D. | U | 0.20 | 0.050 | 0.042 | N.D. | U | 0.22 | 0.054 | 0.046 | 0.30 | | 0.20 | 0.050 | 0.042 | N.D. | U | 0.19 | 0.048 | 0.041 | N.D. | U | 0.19 | 0.048 | 0.041 | |
| | Indeno[1,2,3-cd]pyrene | 0.092 | 0.095 | N.D. | U | 0.20 | 0.050 | 0.021 | N.D. | U | 0.22 | 0.054 | 0.023 | N.D. | U | 0.20 | 0.050 | 0.021 | N.D. | U | 0.19 | 0.048 | 0.020 | N.D. | U | 0.19 | 0.048 | 0.020 | |
| | 1-Methylnaphthalene | 4.7 | 10 | N.D. | U | 0.20 | 0.10 | 0.051 | N.D. | U | 0.22 | 0.11 | 0.056 | 34 | | 2.0 | 0.99 | 0.51 | N.D. | U | 0.19 | 0.097 | 0.050 | N.D. | U | 0.19 | 0.096 | 0.050 | |
| | 2-Methylnaphthalene | 24 | 10 | N.D. | U | 0.20 | 0.050 | 0.046 | N.D. | U | 0.22 | 0.054 | 0.050 | 7.6 | J | 0.20 | 0.050 | 0.046 | N.D. | U | 0.19 | 0.048 | 0.045 | N.D. | U | 0.19 | 0.048 | 0.045 | |
| | Naphthalene | 17 | 21 | N.D. | U | 0.20 | 0.050 | 0.034 | N.D. | U | 0.22 | 0.054 | 0.037 | 90 | | 2.0 | 0.50 | 0.34 | N.D. | U | 0.19 | 0.048 | 0.033 | N.D. | U | 0.19 | 0.048 | 0.033 | |
| | Phenanthrene | 240 | 410 | N.D. | U | 0.20 | 0.050 | 0.027 | N.D. | U | 0.22 | 0.054 | 0.029 | N.D. | U | 0.20 | 0.050 | 0.027 | N.D. | U | 0.19 | 0.048 | 0.026 | N.D. | U | 0.19 | 0.048 | 0.026 | |
| | Pyrene | 180 | 68 | N.D. | U | 0.20 | 0.050 | 0.020 | N.D. | U | 0.22 | 0.054 | 0.022 | N.D. | U | 0.20 | 0.050 | 0.020 | N.D. | U | 0.19 | 0.048 | 0.020 | N.D. | U | 0.19 | 0.048 | 0.020 | |
| | EPA 8260B | 1,1,1,2-Tetrachloroethane | 0.52 | 50,000 | N.D. | U | 1 | 0.5 | 0.4 | N.D. | U | 1 | 0.5 | 0.4 | N.D. | U | 1 | 0.5 | 0.4 | N.D. | U | 1 | 0.5 | 0.4 | N.D. | U | 1 | 0.5 | 0.4 |
| | | 1,1,2,2-Tetrachloroethane ¹ | 0.067 | 500 | N.D. | U | 1 | 0.5 | 0.41 | N.D. | U | 1 | 0.5 | 0.41 | N.D. | U | 1 | 0.5 | 0.41 | N.D. | U | 1 | 0.5 | 0.41 | N.D. | U | 1 | 0.5 | 0.41 |
| 1,1,1-Trichloroethane | | 200 | 970 | N.D. | U | 5 | 0.5 | 0.3 | N.D. | U | 5 | 0.5 | 0.3 | N.D. | U | 5 | 0.5 | 0.3 | N.D. | U | 5 | 0.5 | 0.3 | N.D. | U | 5 | 0.5 | 0.3 | |
| 1,1,2-Trichloroethane | | 5 | 50,000 | N.D. | U | 1 | 0.5 | 0.38 | N.D. | U | 1 | 0.5 | 0.38 | N.D. | U | 1 | 0.5 | 0.38 | N.D. | U | 1 | 0.5 | 0.38 | N.D. | U | 1 | 0.5 | 0.38 | |
| 1,1-Dichloroethane | | 2.4 | 50,000 | N.D. | U | 5 | 0.5 | 0.28 | N.D. | U | 5 | 0.5 | 0.28 | N.D. | U | 5 | 0.5 | 0.28 | N.D. | U | 5 | 0.5 | 0.28 | N.D. | U | 5 | 0.5 | 0.28 | |
| 1,1-Dichloroethylene | | 7 | 1,500 | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.43 | |
| 1,2,3-Trichloropropane ¹ | | 0.6 | 50,000 | N.D. | U | 5 | 1 | 0.64 | N.D. | U | 5 | 1 | 0.64 | N.D. | U | 5 | 1 | 0.64 | N.D. | U | 5 | 1 | 0.64 | N.D. | U | 5 | 1 | 0.64 | |
| 1,2,4-Trichlorobenzene | | 70 | 3,000 | N.D. | U | 5 | 1 | 0.5 | N.D. | U | 5 | 1 | 0.5 | N.D. | U | 5 | 1 | 0.5 | N.D. | U | 5 | 1 | 0.5 | N.D. | U | 5 | 1 | 0.5 | |
| 1,2-Dibromo-3-chloropropane ¹ | | 0.04 | 10 | N.D. | U | 10 | 2 | 1.2 | N.D. | U | 10 | 2 | 1.2 | N.D. | U | 10 | 2 | 1.2 | N.D. | U | 10 | 2 | 1.2 | N.D. | U | 10 | 2 | 1.2 | |
| 1,2-Dibromoethane ¹ | | 0.04 | 50,000 | N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 | |
| 1,2-Dichlorobenzene | | 600 | 10 | N.D. | U | 1 | 0.5 | 0.46 | N.D. | U | 1 | 0.5 | 0.46 | N.D. | U | 1 | 0.5 | 0.46 | N.D. | U | 1 | 0.5 | 0.46 | N.D. | U | 1 | 0.5 | 0.46 | |
| 1,2-Dichloroethane ¹ | | 0.15 | 7,000 | N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 | |
| 1,2-Dichloropropane | | 5 | 10 | N.D. | U | 5 | 0.5 | 0.42 | N.D. | U | 5 | 0.5 | 0.42 | N.D. | U | 5 | 0.5 | 0.42 | N.D. | U | 5 | 0.5 | 0.42 | N.D. | U | 5 | 0.5 | 0.42 | |
| 1,3-Dichlorobenzene | | 180 | 5 | N.D. | U | 1 | 0.5 | 0.4 | N.D. | U | 1 | 0.5 | 0.4 | N.D. | U | 1 | 0.5 | 0.4 | N.D. | U | 1 | 0.5 | 0.4 | N.D. | U | 1 | 0.5 | 0.4 | |
| 1,3-Dichloropropene (total of cis/trans) ¹ | | 0.43 | 50,000 | N.D. | U | 1 | 0.5 | 0.25 | N.D. | U | 1 | 0.5 | 0.25 | N.D. | U | 1 | 0.5 | 0.25 | N.D. | U | 1 | 0.5 | 0.25 | N.D. | U | 1 | 0.5 | 0.25 | |
| 1,4-Dichlorobenzene | | 75 | 5 | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.43 | |
| Acetone | | 22,000 | 20,000 | N.D. | U | 20 | 10 | 6 | N.D. | U | 20 | 10 | 6 | N.D. | U,I,J | 20 | 10 | 6 | N.D. | U,I,J | 20 | 10 | 6 | N.D. | U | 20 | 10 | 6 | |
| Benzene | | 5 | 170 | N.D. | U | 1 | 0.5 | 0.14 | N.D. | U | 1 | 0.5 | 0.14 | N.D. | U | 1 | 0.5 | 0.14 | N.D. | U | 1 | 0.5 | 0.14 | N.D. | U | 1 | 0.5 | 0.14 | |
| Bromodichloromethane ¹ | | 0.12 | 50,000 | N.D. | U | 5 | 0.5 | 0.21 | N.D. | U | 5 | 0.5 | 0.21 | N.D. | U | 5 | 0.5 | 0.21 | N.D. | U | 5 | 0.5 | 0.21 | N.D. | U | 5 | 0.5 | 0.21 | |
| Bromoform | | 80 | 510 | N.D. | U | 10 | 1 | 0.5 | N.D. | U | 10 | 1 | 0.5 | N.D. | U | 10 | 1 | 0.5 | N.D. | U | 10 | 1 | 0.5 | N.D. | U | 10 | 1 | 0.5 | |
| Bromomethane | | 8.7 | 50,000 | N.D. | U,I,H | 20 | 5 | 3.9 | N.D. | U,I,H | 20 | 5 | 3.9 | N.D. | U,I,H | 20 | 5 | 3.9 | N.D. | U,I,H | 20 | 5 | 3.9 | N.D. | U,I,H | 20 | 5 | 3.9 | |
| Carbon Tetrachloride | | 5 | 520 | N.D. | U | 1 | 0.5 | 0.23 | N.D. | U | 1 | 0.5 | 0.23 | N.D. | U | 1 | 0.5 | 0.23 | N.D. | U | 1 | 0.5 | 0.23 | N.D. | U | 1 | 0.5 | 0.23 | |
| Chlorobenzene | | 100 | 50 | N.D. | U | 5 | 0.5 | 0.17 | N.D. | U | 5 | 0.5 | 0.17 | N.D. | U | 5 | 0.5 | 0.17 | N.D. | U | 5 | 0.5 | 0.17 | N.D. | U | 5 | 0.5 | 0.17 | |
| Chloroethane | | 21,000 | 16 | N.D. | U | 10 | 5 | 2.3 | N.D. | U | 10 | 5 | 2.3 | N.D. | U | 10 | 5 | 2.3 | N.D. | U | 10 | 5 | 2.3 | N.D. | U | 10 | 5 | 2.3 | |
| Chloroform | | 70 | 2,400 | N.D. | U | 5 | 0.5 | 0.46 | N.D. | U | 5 | 0.5 | 0.46 | N.D. | U | 5 | 0.5 | 0.46 | N.D. | U | 5 | 0.5 | 0.46 | N.D. | U | 5 | 0.5 | 0.46 | |
| Chloromethane ¹ | | 1.8 | 50,000 | N.D. | U | 10 | 2 | 1.8 | N.D. | U | 10 | 2 | 1.8 | N.D. | U | 10 | 2 | 1.8 | N.D. | U | 10 | 2 | 1.8 | N.D. | U | 10 | 2 | 1.8 | |
| cis-1,2-Dichloroethylene | | 70 | 50,000 | N.D. | U | 1 | 0.5 | 0.48 | N.D. | U | 1 | 0.5 | 0.48 | N.D. | U | 1 | 0.5 | 0.48 | N.D. | U | 1 | 0.5 | 0.48 | N.D. | U | 1 | 0.5 | 0.48 | |

This Page Intentionally Left Blank.

SECTION 3 – DATA QUALITY ASSESSMENT

A data quality assessment, which consists of a review of the overall groundwater sample collection and analysis process, was performed in order to determine whether the analytical data generated met the quality objectives for the project. The data quality assessment was performed in accordance with the approved WP/SAP (ESI, 2012). The field quality control program consisted of standardized sample collection and management procedures, and the collection of field duplicate samples, matrix spike samples, and trip blank samples. The laboratory quality assurance program consisted of the use of standard analytical methods and the preparation and analyses of Matrix Spike [MS]/Matrix Spike Duplicate [MSD] samples, surrogate spikes, blanks, and Laboratory Control Samples [LCSs]/Laboratory Control Sample Duplicates [LCSDs].

3.1 DATA VALIDATION AND ASSESSMENT

The objective of data validation is to provide data of known quality for project decisions. Data quality is judged in terms of Precision, Accuracy, Representativeness, Completeness, Comparability, and Sensitivity [PARCCS]. A number of factors may affect the quality of data, including: sample collection methods, sample analysis methods, and adherence to established procedures for sample collection, preservation, management, shipment, and analysis.

Precision

Precision is defined as the reproducibility of replicate measurements. Precision is evaluated by Relative Percentage Difference [RPD] of field duplicates, LCS/LCSD, and MS/MSD results. Field duplicate and MS/MSD samples were collected at a rate of approximately 10% of primary samples. Field duplicates were sent to the laboratory along with the primary samples.

The RPDs of detected analytes for the primary and field duplicate samples (ES126 and ES127) are provided in Table 3.1. A precision of less than 50% for duplicate pairs is required by the DON Project Procedures Manual to be considered acceptable (DON, 2007). All duplicate RPDs were below the acceptable maximum, except for 2-methylnaphthalene (95%). Both 2-methylnaphthalene results were flagged “J” to indicate a lack of precision. 2-Methylnaphthalene results for both samples were below the DOH EALs but have been detected at concentrations above DOH EALs during previous events. Judged solely by the magnitude of the imprecision associated with the samples it is unlikely that the actual concentration of 2-methylnaphthalene in the RHMW02 samples collected during the first quarter 2015 event exceeded the EALs (gross contamination: 10 µg/l; or drinking water toxicity: 24 µg/l). However, the imprecision is evaluated in the context of additional errors that may have a negative impact on data quality and the ability to use the data for decision making. The potential low bias in the data described in the accuracy section below is one of these additional errors. 2-Methylnaphthalene was not detected in any other sample analyzed in the course of the January 2015 sampling event; and because the LOQs associated with these results (non-detections) were sufficiently low, the negative impact of the lack of precision on data usability (for non-detections) was negligible.

Similarly, the RPD for 1,1,2,2-tetrachloroethane in the MS/MSD was out of control indicating a negative impact on precision due to matrix effects. Although 1,1,2,2-tetrachloroethane was not detected in any samples, the LOQs and LODs for this analyte already exceed the EAL (for drinking water toxicity). Thus, the additional error may have a negative impact on the ability to identify potential contamination with this analyte at concentrations close to the LODs. However, the analyte was not detected previously in any of the samples at levels above the detection limit. When the datum is considered in the context of all previous data obtained in the course of the LTM project, the additional impact on project decision making due to the lack of precision in the current datum is insignificant. RPDs for MS/MSD and LCS/LCSD pairs for all other analytes were within the control limits, and the data precision is considered acceptable.

Accuracy

Accuracy is defined as the degree of conformity of a measurement to a standard or true value. Accuracy is evaluated through measurement of the percent recovery of an analyte in a reference standard or spiked sample. Accuracy limits for surrogates, laboratory control spike, MS, and MSD samples are either prescribed by the Department of Defense [DoD] or established by the individual laboratory. The acceptance criteria for accuracy are dependent on the analytical method and are based on historical laboratory or DoD data.

Between July 2006 and July 2010, naphthalene was analyzed by both EPA Methods 8260B and 8270C, and both results were reported. In September 2005 and in all data beginning in October 2010, only results using EPA Method 8270C were reported. Naphthalene has historically only been detected at concentrations above the DOH EALs in well RHMW02. In this well, concentrations of naphthalene detected in each sample by EPA Method 8260B were generally two to three times higher than those detected by EPA Method 8270C. We assume this is due to the better preservation of VOCs associated with the use of EPA Method 8260B. This suggests that the naphthalene results provided by EPA Method 8270C may be biased low. Since March 2014, naphthalene concentrations in RHMW02 have exceeded DOH EALs for both gross contamination and drinking water toxicity. Therefore, a low bias is unlikely to affect project decisions.

Results for TPH-d in samples ES120X, ES123, ES126, and ES127 were flagged "HD." The laboratory indicated a mismatch between the calibration standard and the TPH-d chromatographic profile. Mismatches of this type are not uncommon. Even though chromatograms are not part of the standard laboratory package, ESI was able to review the chromatograms from RHMW02 dating back to October 2012. The chromatograms of groundwater samples from RHMW02 did not significantly differ between each event, and did not match a standard chromatogram of JP-8 in water.

The MS and MSD recoveries were above the control limits for acetone and trichloroethene and the associated sample results may be biased high; however, neither of these analytes were detected in any samples. The MS and MSD recoveries were below the control limits for 1,1,2,2-tetrachloroethane, 1,2,3-trichloropropane, naphthalene, 1-methylnaphthalene, and

2-methylnaphthalene. 1,1,2,2-Tetrachloroethane and 1,2,3-trichloropropane have not historically been detected in the wells. However, the LOQs and LODs of both compounds exceed the EALs (for drinking water toxicity), with additional issues of low precision for 1,1,2,2-tetrachloroethane data thus potentially increasing the range (between EAL and LOQs) of concentrations at which the compounds may be present in samples at a concentration above the EALs but not detected. Naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene concentrations for ES126, the primary sample on which the MS/MSD were performed, were significantly higher than the added spike concentration, which prevented an accurate evaluation of the MS/MSD recovery for these analytes.

An air bubble was present in one of the VOA containers for sample ES124. However, this VOA container was not used in sample analysis, and this does not affect data usability.

The surrogate spike recovery for nitrobenzene-d5 in samples ES126 and ES127 was below the control limits (0%) for naphthalene and 1-methylnaphthalene results. This is likely a result of the sample dilution, since nitrobenzene-d5 was in control for the PAH results from undiluted sample runs. However, surrogate recoveries for PAH results were generally low (mostly <70%), indicating that a slight low bias may be associated with the data. All other MS/MSD, LCS/LCSD, and surrogate spike recoveries were within acceptable recovery limits; therefore, the data accuracy for this monitoring event is considered acceptable with the exceptions noted above.

Representativeness

Representativeness is the degree that data accurately and precisely represents a characteristic of a population, parameter variations at a sampling point, or an environmental condition. Representativeness was achieved by conducting sampling in accordance with the sample collection procedures described in the project WP/SAP, including standardized sample collection methods (ESI, 2012).

Representativeness is also evaluated through the compliance with the standardized sample holding time and sample preservation methods, and through the analysis of blank samples, including method blank and trip blank samples. For this sampling event, all sample holding times and sample preservation were consistent with EPA guidance.

For this sampling event, one trip blank was included in every cooler containing samples for VOC and TPH-g analysis to assess the potential for contamination during sample transport. Two trip blanks were collected. No analytes were detected in either trip blank. Based on the assessment of representativeness, the groundwater sample data are considered representative of the groundwater quality on site. The trip blank results are provided in Table 3.1.

Completeness

Completeness is defined as the overall percentage of valid analytical results (including estimated results) compared to the total number of analytical results reported by the analytical

laboratory. No data were rejected for this project, and therefore the completeness goal for this project (90%), was successfully met.

Comparability

Comparability expresses the confidence with which one data set can be compared to another data set. Comparability can be related to accuracy and precision because these quantities are measures of data reliability. Data with acceptable precision and accuracy are considered comparable if collection techniques, analytical procedures, methods and reporting are equivalent.

As noted above, between July 2006 and July 2010, naphthalene was analyzed for using both EPA Methods 8260B and 8270C, and in September 2005 and between October 2010 and the most recent event, only results using EPA Method 8270C were reported. In general, EPA Method 8260B resulted in higher, and as discussed above, likely more accurate, results than EPA Method 8270C. However, for the sake of comparability with results from recent events, EPA Method 8270C was used for naphthalene analysis in this event. Consequently, the low bias associated with Method 8270C should be considered when making project decisions.

All project samples for TPH-g analysis through July 2010 were analyzed by EPA Method 8015; beginning in October 2010, EPA Method 8260B was used. There was no event where both methods were used; consequently, there is no way to directly compare the results obtained by method and to assess potential bias. However, there is no reason to believe that using either method should bias the data, and the TPH-g data for all events should be comparable.

Other than the naphthalene bias discussed above, no issues with comparability were identified. The results are considered comparable within this data set and with the data collected from recent sampling events.

Sensitivity

The limits of quantitation [LOQs] are established by the laboratory based on the LODs or instrument detection limits, historical data, and EPA limits established for the various methods. The LOQs for samples may require adjustment by the laboratory due to matrix interference or if high levels of target analytes necessitate dilution before analysis. Matrix interference and sample dilutions have the effect of decreasing sensitivity and increasing the LOQs. Laboratory LODs and LOQs for several analytes (EPA Methods 8260 and 8270) for this event differed from the LODs and LOQs in the WP/SAP because the laboratory updates them quarterly and in some cases, dilution was necessary due to the presence of high concentrations of analytes.

For this event, LODs and LOQs for several analytes were greater than the DOH EALs (as stated in the WP/SAP), negatively impacting the ability to detect analytes at concentrations greater than the DOH EALs but below the LODs. The lack of the required sensitivity should be considered when making project decisions. The affected analytes for this monitoring event are 1,2,3-trichloropropane, 1,2-dibromo-3-chloropropane, 1,2-dibromoethane, 1,2-dichloroethane,

1,3-dichloropropene, bromodichloromethane, chloromethane, dibromochloromethane, 1,1,2,2-tetrachloroethane, and dibenzo[a,h]anthracene. Additional uncertainty associated with the data may reduce the capability of detecting these compounds at concentrations between the LODs and the EALs even further. Issues with accuracy and precision for these data that may have contributed to decreased sensitivity are discussed in their respective sections.

3.2 DATA ASSESSMENT AND USABILITY CONCLUSIONS

The PARCCS criteria were evaluated, and with a few exceptions, all criteria were met. Based on the high RPD of 2-methylnaphthalene results in duplicate samples ES126 and ES127, it is possible that there may be lower precision in the 2-methylnaphthalene results. This decreased precision should be kept in mind when comparing 2-methylnaphthalene results from this sampling event to those of previous events. During this event, Well RHMW02 is the only well where 2-methylnaphthalene was detected at concentrations near the DOH EAL. Because concentrations of 1-methylnaphthalene and naphthalene were detected at concentrations above DOH EALs during this event, the lower precision of 2-methylnaphthalene results should not significantly affect project decisions.

The surrogate spike recovery for nitrobenzene-d5 was below the control limits for naphthalene and 1-methylnaphthalene results, and the surrogate recoveries for other PAH results were generally low (mostly <70%), indicating that a slight low bias may be associated with the PAH data. However, all PAHs except 2-methylnaphthalene (discussed above) were either not detected, detected at concentrations well below DOH EALs, or detected at concentrations above DOH EALs. Since no PAHs were detected at concentrations slightly below DOH EALs, it is unlikely that a slight low bias would have an impact on project decisions.

Additionally, LODs and LOQs for several analytes were greater than the DOH EALs. These analytes have not historically been detected in any of the wells, and with the exception of 1,2-dichloroethane, are not likely to be related to a fuel release. Therefore, the lack of required sensitivity should not have a significant impact on project decisions.

Other than these issues, the data assessment concludes that all data generated during this event are usable for the intended purpose.

This Page Intentionally Left Blank

TABLE 3.1
Quality Control Results for Groundwater Sampling (January 27 and 28, 2015)
Red Hill Bulk Fuel Storage Facility
January 2015 Quarterly Monitoring Report

| Method | Chemical Constituent | DOH EALs | | RHMW02 (ES126) | | | | | RHMW02 (ES127) (DUP) | | | | | RPD Duplicate (%) | ES Trip (1/27/2015) | | | | | ES Trip (1/28/2015) | | | | |
|-----------|---|-------------------------|---------------------|----------------|------|-------|-------|-------|----------------------|------|-------|-------|-------|-------------------|---------------------|------|-----|-----|------|---------------------|------|-----|-----|------|
| | | Drinking Water Toxicity | Gross Contamination | Result | Q | LOQ | LOD | DL | Result | Q | LOQ | LOD | DL | | Result | Q | LOQ | LOD | DL | Result | Q | LOQ | LOD | DL |
| EPA 8015B | TPH-d | 190 | 100 | 1,100 | HD | 25 | 10 | 2.9 | 1,700 | HD | 25 | 10 | 2.9 | 42.86 | - | - | - | - | - | - | - | - | - | - |
| EPA 8260B | TPH-g | 100 | 100 | 54 | | 50 | 30 | 26 | 59 | | 50 | 30 | 26 | 8.85 | N.D. | U | 50 | 30 | 26 | N.D. | U | 50 | 30 | 26 |
| EPA 8270C | Acenaphthene | 370 | 20 | 0.59 | | 0.20 | 0.050 | 0.027 | 0.55 | | 0.20 | 0.049 | 0.026 | 7.02 | - | - | - | - | - | - | - | - | - | - |
| | Acenaphthylene | 240 | 2,000 | N.D. | U | 0.20 | 0.050 | 0.044 | N.D. | U | 0.20 | 0.049 | 0.044 | NA | - | - | - | - | - | - | - | - | - | - |
| | Anthracene | 1,800 | 22 | N.D. | U | 0.20 | 0.050 | 0.029 | N.D. | U | 0.20 | 0.049 | 0.028 | NA | - | - | - | - | - | - | - | - | - | - |
| | Benzo[a]anthracene | 0.092 | 4.7 | N.D. | U | 0.20 | 0.050 | 0.032 | N.D. | U | 0.20 | 0.049 | 0.032 | NA | - | - | - | - | - | - | - | - | - | - |
| | Benzo[g,h,i]perylene | 1,500 | 0.13 | N.D. | U | 0.20 | 0.099 | 0.081 | N.D. | U | 0.20 | 0.098 | 0.080 | NA | - | - | - | - | - | - | - | - | - | - |
| | Benzo[a]pyrene | 0.2 | 0.81 | N.D. | U | 0.20 | 0.050 | 0.022 | N.D. | U | 0.20 | 0.049 | 0.022 | NA | - | - | - | - | - | - | - | - | - | - |
| | Benzo[b]fluoranthene | 0.092 | 0.75 | N.D. | U | 0.20 | 0.050 | 0.017 | N.D. | U | 0.20 | 0.049 | 0.017 | NA | - | - | - | - | - | - | - | - | - | - |
| | Benzo[k]fluoranthene | 0.92 | 0.4 | N.D. | U | 0.20 | 0.050 | 0.031 | N.D. | U | 0.20 | 0.049 | 0.030 | NA | - | - | - | - | - | - | - | - | - | - |
| | Chrysene | 9.2 | 1 | N.D. | U | 0.20 | 0.050 | 0.025 | N.D. | U | 0.20 | 0.049 | 0.024 | NA | - | - | - | - | - | - | - | - | - | - |
| | Dibenzo[a,h]anthracene | 0.0092 | 0.52 | N.D. | U | 0.20 | 0.050 | 0.047 | N.D. | U | 0.20 | 0.049 | 0.047 | NA | - | - | - | - | - | - | - | - | - | - |
| | Fluoranthene | 1,500 | 130 | N.D. | U | 0.20 | 0.050 | 0.046 | N.D. | U | 0.20 | 0.049 | 0.046 | NA | - | - | - | - | - | - | - | - | - | - |
| | Fluorene | 240 | 950 | 0.30 | | 0.20 | 0.050 | 0.042 | 0.22 | | 0.20 | 0.049 | 0.042 | 30.77 | - | - | - | - | - | - | - | - | - | - |
| | Indeno[1,2,3-cd]pyrene | 0.092 | 0.095 | N.D. | U | 0.20 | 0.050 | 0.021 | N.D. | U | 0.20 | 0.049 | 0.021 | NA | - | - | - | - | - | - | - | - | - | - |
| | 1-Methylnaphthalene | 4.7 | 10 | 34 | | 2.0 | 0.99 | 0.51 | 25 | | 2.0 | 0.98 | 0.51 | 30.51 | - | - | - | - | - | - | - | - | - | - |
| | 2-Methylnaphthalene | 24 | 10 | 7.6 | J | 0.20 | 0.050 | 0.046 | 2.7 | J | 0.20 | 0.049 | 0.046 | 95.15 | - | - | - | - | - | - | - | - | - | - |
| | Naphthalene | 17 | 21 | 90 | | 2.0 | 0.50 | 0.34 | 63 | | 2.0 | 0.49 | 0.33 | 35.29 | - | - | - | - | - | - | - | - | - | - |
| | Phenanthrene | 240 | 410 | N.D. | U | 0.20 | 0.050 | 0.027 | N.D. | U | 0.20 | 0.049 | 0.027 | NA | - | - | - | - | - | - | - | - | - | - |
| Pyrene | 180 | 68 | N.D. | U | 0.20 | 0.050 | 0.020 | N.D. | U | 0.20 | 0.049 | 0.020 | NA | - | - | - | - | - | - | - | - | - | - | |
| EPA 8260B | 1,1,1,2-Tetrachloroethane | 0.52 | 50,000 | N.D. | U | 1 | 0.5 | 0.4 | N.D. | U | 1 | 0.5 | 0.4 | NA | N.D. | U | 1 | 0.5 | 0.4 | N.D. | U | 1 | 0.5 | 0.4 |
| | 1,1,2,2-Tetrachloroethane | 0.067 | 500 | N.D. | U | 1 | 0.5 | 0.41 | N.D. | U | 1 | 0.5 | 0.41 | NA | N.D. | U | 1 | 0.5 | 0.41 | N.D. | U | 1 | 0.5 | 0.41 |
| | 1,1,1-Trichloroethane | 200 | 970 | N.D. | U | 5 | 0.5 | 0.3 | N.D. | U | 5 | 0.5 | 0.3 | NA | N.D. | U | 5 | 0.5 | 0.3 | N.D. | U | 5 | 0.5 | 0.3 |
| | 1,1,2-Trichloroethane | 5 | 50,000 | N.D. | U | 1 | 0.5 | 0.38 | N.D. | U | 1 | 0.5 | 0.38 | NA | N.D. | U | 1 | 0.5 | 0.38 | N.D. | U | 1 | 0.5 | 0.38 |
| | 1,1-Dichloroethane | 2.4 | 50,000 | N.D. | U | 5 | 0.5 | 0.28 | N.D. | U | 5 | 0.5 | 0.28 | NA | N.D. | U | 5 | 0.5 | 0.28 | N.D. | U | 5 | 0.5 | 0.28 |
| | 1,1-Dichloroethylene | 7 | 1,500 | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.43 | NA | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.43 |
| | 1,2,3-Trichloropropane | 0.6 | 50,000 | N.D. | U | 5 | 1 | 0.64 | N.D. | U | 5 | 1 | 0.64 | NA | N.D. | U | 5 | 1 | 0.64 | N.D. | U | 5 | 1 | 0.64 |
| | 1,2,4-Trichlorobenzene | 70 | 3,000 | N.D. | U | 5 | 1 | 0.5 | N.D. | U | 5 | 1 | 0.5 | NA | N.D. | U | 5 | 1 | 0.5 | N.D. | U | 5 | 1 | 0.5 |
| | 1,2-Dibromo-3-chloropropane | 0.04 | 10 | N.D. | U | 10 | 2 | 1.2 | N.D. | U | 10 | 2 | 1.2 | NA | N.D. | U | 10 | 2 | 1.2 | N.D. | U | 10 | 2 | 1.2 |
| | 1,2-Dibromoethane | 0.04 | 50,000 | N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 | NA | N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 |
| | 1,2-Dichlorobenzene | 600 | 10 | N.D. | U | 1 | 0.5 | 0.46 | N.D. | U | 1 | 0.5 | 0.46 | NA | N.D. | U | 1 | 0.5 | 0.46 | N.D. | U | 1 | 0.5 | 0.46 |
| | 1,2-Dichloroethane | 0.15 | 7,000 | N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 | NA | N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 |
| | 1,2-Dichloropropane | 5 | 10 | N.D. | U | 5 | 0.5 | 0.42 | N.D. | U | 5 | 0.5 | 0.42 | NA | N.D. | U | 5 | 0.5 | 0.42 | N.D. | U | 5 | 0.5 | 0.42 |
| | 1,3-Dichlorobenzene | 180 | 5 | N.D. | U | 1 | 0.5 | 0.4 | N.D. | U | 1 | 0.5 | 0.4 | NA | N.D. | U | 1 | 0.5 | 0.4 | N.D. | U | 1 | 0.5 | 0.4 |
| | 1,3-Dichloropropene (total of cis/trans) | 0.43 | 50,000 | N.D. | U | 1 | 0.5 | 0.25 | N.D. | U | 1 | 0.5 | 0.25 | NA | N.D. | U | 1 | 0.5 | 0.25 | N.D. | U | 1 | 0.5 | 0.25 |
| | 1,4-Dichlorobenzene | 75 | 5 | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.43 | NA | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.43 |
| | Acetone | 22,000 | 20,000 | N.D. | U,IJ | 20 | 10 | 6 | N.D. | U,IJ | 20 | 10 | 6 | NA | N.D. | U | 20 | 10 | 6 | N.D. | U,IJ | 20 | 10 | 6 |
| | Benzene | 5 | 170 | N.D. | U | 1 | 0.5 | 0.14 | N.D. | U | 1 | 0.5 | 0.14 | NA | N.D. | U | 1 | 0.5 | 0.14 | N.D. | U | 1 | 0.5 | 0.14 |
| | Bromodichloromethane | 0.12 | 50,000 | N.D. | U | 5 | 0.5 | 0.21 | N.D. | U | 5 | 0.5 | 0.21 | NA | N.D. | U | 5 | 0.5 | 0.21 | N.D. | U | 5 | 0.5 | 0.21 |
| | Bromoform | 80 | 510 | N.D. | U | 10 | 1 | 0.5 | N.D. | U | 10 | 1 | 0.5 | NA | N.D. | U | 10 | 1 | 0.5 | N.D. | U | 10 | 1 | 0.5 |
| | Bromomethane | 8.7 | 50,000 | N.D. | U,IH | 20 | 5 | 3.9 | N.D. | U,IH | 20 | 5 | 3.9 | NA | N.D. | U,IH | 20 | 5 | 3.9 | N.D. | U,IH | 20 | 5 | 3.9 |
| | Carbon Tetrachloride | 5 | 520 | N.D. | U | 1 | 0.5 | 0.23 | N.D. | U | 1 | 0.5 | 0.23 | NA | N.D. | U | 1 | 0.5 | 0.23 | N.D. | U | 1 | 0.5 | 0.23 |
| | Chlorobenzene | 100 | 50 | N.D. | U | 5 | 0.5 | 0.17 | N.D. | U | 5 | 0.5 | 0.17 | NA | N.D. | U | 5 | 0.5 | 0.17 | N.D. | U | 5 | 0.5 | 0.17 |
| | Chloroethane | 21,000 | 16 | N.D. | U | 10 | 5 | 2.3 | N.D. | U | 10 | 5 | 2.3 | NA | N.D. | U | 10 | 5 | 2.3 | N.D. | U | 10 | 5 | 2.3 |
| | Chloroform | 70 | 2,400 | N.D. | U | 5 | 0.5 | 0.46 | N.D. | U | 5 | 0.5 | 0.46 | NA | N.D. | U | 5 | 0.5 | 0.46 | N.D. | U | 5 | 0.5 | 0.46 |
| | Chloromethane | 1.8 | 50,000 | N.D. | U | 10 | 2 | 1.8 | N.D. | U | 10 | 2 | 1.8 | NA | N.D. | U | 10 | 2 | 1.8 | N.D. | U | 10 | 2 | 1.8 |
| | cis-1,2-Dichloroethylene | 70 | 50,000 | N.D. | U | 1 | 0.5 | 0.48 | N.D. | U | 1 | 0.5 | 0.48 | NA | N.D. | U | 1 | 0.5 | 0.48 | N.D. | U | 1 | 0.5 | 0.48 |
| | Dibromochloromethane | 0.16 | 50,000 | N.D. | U | 1 | 0.5 | 0.25 | N.D. | U | 1 | 0.5 | 0.25 | NA | N.D. | U | 1 | 0.5 | 0.25 | N.D. | U | 1 | 0.5 | 0.25 |
| | Ethylbenzene | 700 | 30 | 0.16 | J | 1 | 0.5 | 0.14 | 0.17 | J | 1 | 0.5 | 0.14 | 6.06 | N.D. | U | 1 | 0.5 | 0.14 | N.D. | U | 1 | 0.5 | 0.14 |
| | Hexachlorobutadiene | 0.86 | 6 | N.D. | U | 1 | 0.5 | 0.32 | N.D. | U | 1 | 0.5 | 0.32 | NA | N.D. | U | 1 | 0.5 | 0.32 | N.D. | U | 1 | 0.5 | 0.32 |
| | Methyl ethyl ketone (2-Butanone) | 7,100 | 8,400 | N.D. | U | 10 | 5.0 | 2.2 | N.D. | U | 10 | 5.0 | 2.2 | NA | N.D. | U | 10 | 5.0 | 2.2 | N.D. | U | 10 | 5.0 | 2.2 |
| | Methyl isobutyl ketone (4-Methyl-2-Pentanone) | 2,000 | 1300 | N.D. | U | 10 | 5.0 | 4.4 | N.D. | U | 10 | 5.0 | 4.4 | NA | N.D. | U | 10 | 5.0 | 4.4 | N.D. | U | 10 | 5.0 | 4.4 |
| | Methyl tert-butyl Ether | 12 | 5 | N.D. | U | 1 | 0.5 | 0.31 | N.D. | U | 1 | 0.5 | 0.31 | NA | N.D. | U | 1 | 0.5 | 0.31 | N.D. | U | 1 | 0.5 | 0.31 |
| | Methylene chloride | 4.8 | 9,100 | N.D. | U | 5 | 1.0 | 0.64 | N.D. | U | 5 | 1.0 | 0.64 | NA | N.D. | U | 5 | 1.0 | 0.64 | N.D. | U | 5 | 1.0 | 0.64 |
| | Styrene | 100 | 10 | N.D. | U | 1 | 0.5 | 0.17 | N.D. | U | 1 | 0.5 | 0.17 | NA | N.D. | U | 1 | 0.5 | 0.17 | N.D. | U | 1 | 0.5 | 0.17 |
| | Tetrachloroethylene | 5 | 170 | N.D. | U | 5 | 0.5 | 0.39 | N.D. | U | 5 | 0.5 | 0.39 | NA | N.D. | U | 5 | 0.5 | 0.39 | N.D. | U | 5 | 0.5 | 0.39 |
| | Toluene | 1,000 | 40 | N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 | NA | N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 |
| | trans-1,2-Dichloroethylene | 100 | 260 | N.D. | U | 1 | 0.5 | 0.37 | N.D. | U | 1 | 0.5 | 0.37 | NA | N.D. | U | 1 | 0.5 | 0.37 | N.D. | U | 1 | 0.5 | 0.37 |
| | Trichloroethylene | 5 | 310 | | | | | | | | | | | | | | | | | | | | | |

This Page Intentionally Left Blank.

SECTION 4 – SUMMARY, CONCLUSIONS, AND RECOMMENDATIONS

On January 27 and 28, 2015, ESI personnel collected groundwater samples from four monitoring wells at the RHSF (wells RHMW01, RHMW02, RHMW03, and RHMW05) and one sampling point at Red Hill Shaft (RHMW2254-01).

The groundwater sampling was conducted as part of the long-term groundwater and soil vapor monitoring program at the RHSF, under NAVFAC Contract Number N62742-12-D-1853. The sampling was conducted in accordance with the approved WP/SAP prepared by ESI. A summary of the analytical results is provided below.

- **RHMW01** – TPH-d (33 µg/L) and lead (0.0631 µg/L) were the only analytes detected. The concentrations did not exceed the DOH EALs or the SSRBL.
- **RHMW02** – TPH-d (1,100 and 1,700 µg/L), TPH-g (54 and 59 µg/L), xylenes (0.35 and 0.35 µg/L), acenaphthene (0.59 and 0.55 µg/L), ethylbenzene (0.16 and 0.17 µg/L), fluorene (0.30 and 0.22 µg/L), 1-methylnaphthalene (34 and 25 µg/L), 2-methylnaphthalene (7.6 and 2.7 µg/L), and naphthalene (90 and 63 µg/L) were detected in both the primary and duplicate samples collected. TPH-d, 1-methylnaphthalene, and naphthalene were detected at concentrations above their respective DOH EALs for both drinking water toxicity and gross contamination. However, the TPH-d concentrations did not exceed the SSRBL.
- **RHMW03** – TPH-d (39 µg/L) was the only analyte detected. The concentration did not exceed the DOH EALs or the SSRBL.
- **RHMW05** – None of the chemical constituents analyzed for were detected at a concentration at or above the LOD.
- **RHMW2254-01** – None of the chemical constituents analyzed for were detected at a concentration at or above the LOD.

Groundwater Contaminant Trends

- **RHMW01** – COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW01. TPH-d has historically been detected at concentrations above the DOH EAL for both drinking water toxicity and gross contamination. TPH-d concentrations continue to show an overall decreasing trend from a high of 1,500 µg/L in February 2005.
- **RHMW02** – COPCs detected during this round of quarterly sampling are generally consistent with the historical data for RHMW02. TPH-g, TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene have historically been detected at concentrations above the DOH EALs. During the January 2015 event, concentrations of TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene decreased from the previous event in October 2014, with the concentrations of 2-methylnaphthalene decreasing to levels below the DOH EALs. The concentrations of TPH-g remained below the DOH EALs for gross contamination and drinking water toxicity and were comparable to the concentrations

detected during the previous event. Trichloroethylene was detected once in RHMW02 in September 2005 in the primary sample at a concentration above the DOH EAL for drinking water toxicity; however, trichloroethylene was not detected in the duplicate sample, and this may have been an anomalous result.

- **RHMW03** – COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW03. TPH-d has historically been detected at concentrations above the DOH EALs; however, it has not been detected at concentrations above the DOH EALs since October 2010.
- **RHMW05** – COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW05. TPH-d has historically been detected in RHMW05 at concentrations above the DOH EALs for both drinking water toxicity and gross contamination; however, it has not been detected at concentrations above the DOH EALs since January 2010.
- **RHMW2254-01** – COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW2254-01. Although the method reporting limits for TPH-d exceeded one or both DOH EALs for drinking water toxicity and gross contamination between May 2009 and July 2010, TPH-d was last detected in RHMW2254-01 at a concentration above the DOH EAL for gross contamination in January 2008.

Conclusions and Recommendations

During the sampling event conducted on January 27 and 28, 2015, TPH-d, 1-methylnaphthalene, and naphthalene in RHMW02 were detected at concentrations exceeding the DOH EALs. The concentration of TPH-d in RHMW01 decreased from the previous event in October 2014 to a concentration below the DOH EALs. Groundwater contaminant concentrations in RHMW03, RHMW05, and RHMW2254-01 remained at low concentrations and did not change significantly from the previous event, or were not detected.

Concentrations of 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene in RHMW02 had shown an increasing trend between March and October 2014; however, concentrations of these three analytes and TPH-d all decreased since the previous event in October 2014. All other analytical results were generally consistent with historical data.

For this event, LODs and LOQs for several analytes were greater than the DOH EALs. However, with the exception of 1,2-dichloroethane, these analytes are not likely to be related to a fuel release, and these elevated LODs and LOQs should not have a significant impact on project objectives.

Based on the groundwater monitoring results and the reported release at Tank 5 in January 2014, continued groundwater monitoring at the wells inside the RHSF tunnel is recommended.

SECTION 5 – FUTURE WORK

Future work includes the second quarter 2015 groundwater monitoring which is tentatively scheduled for April 2015. A quarterly groundwater monitoring report will be prepared to document the sampling event.

This Page Intentionally Left Blank.

SECTION 6 – REFERENCES

Atlas of Hawaii, 1983, Department of Geography, University of Hawaii Press.

DLNR, 1985, Pan Evaporation: State of Hawai'i 1894-1983: Report R74, Division of Water and Land Development, August 1995.

DLNR, 1986, Rainfall Atlas of Hawaii: Report R76, Division of Water and Land Development, June 1986.

DOH, 2000, Hawaii Department of Health, Technical Guidance Manual for Underground Storage Tank Closure and Release Response, March 2000.

DOH, 2011, Evaluation of Environmental Hazards at Sites with Contaminated Soil and Groundwater, Hawaii Department of Health, Hazard Evaluation and Emergency Response, Fall 2011, Revised January 2012.

DOH, 2013, Hawaii Administrative Rules Title 11, Department of Health, Chapter 281 Underground Storage Tanks (HAR 11-281), Subchapter 7.

DON, 2007, Project Procedures Manual, U.S. Navy Installation Restoration Program, NAVFAC Pacific, Prepared for Pacific Division, Naval Facilities Engineering Command (NAVFAC Pacific), February 2007.

Earth Tech, 1999, Remedial Investigation, Phase II Technical Report, Red Hill Oily Waste Disposal Facility, Halawa, Oahu, Hawaii, Prepared for Naval Facilities Engineering Command, Pearl Harbor, Hawaii, 1999.

Environet, 2010, Work Plan, Long-Term Monitoring, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, September 2010.

ESI, 2012, Work Plan/Sampling and Analysis Plan, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, October 2012.

ESI, 2013a, Fourth Quarter 2012 - Quarterly Groundwater Monitoring Report Inside Tunnel Wells, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, January 2013.

ESI, 2013b, First Quarter 2013 - Quarterly Groundwater Monitoring Report Inside Tunnel Wells, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, April 2013.

ESI, 2013c, Second Quarter 2013 - Quarterly Groundwater Monitoring Report Inside Tunnel Wells, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, July 2013.

ESI, 2013d, Third Quarter 2013 - Quarterly Groundwater Monitoring Report Inside Tunnel Wells, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, September 2013.

ESI, 2014a, Fourth Quarter 2013 - Quarterly Groundwater Monitoring Report Inside Tunnel Wells, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, January 2014.

ESI, 2014b, Groundwater Sampling Report for Additional Sampling, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, January 2014.

ESI, 2014c, Final Groundwater Sampling Report for Tank 5 Release Response on March 5 and 6, 2014, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, March 2014.

ESI, 2014d, Final Groundwater Sampling Report for Tank 5 Release Response on March 10, 2014, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, March 2014.

ESI, 2014e, First Quarter 2014 - Quarterly Groundwater Monitoring Report Inside Tunnel Wells, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, April 2014.

ESI, 2014f, Final Groundwater Sampling Report for Tank 5 Release Response on March 25 and 26, 2014, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, April 2014.

ESI, 2014g, Final Groundwater Sampling Report for Tank 5 Release Response on April 7, 2014, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, April 2014.

ESI, 2014h, Second Quarter 2014 - Quarterly Groundwater Monitoring Report Inside Tunnel Wells, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, June 2014.

ESI, 2014i, Final Groundwater Sampling Report for Tank 5 Release Response on May 27 and 28, 2014, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, June 2014.

ESI, 2014j, Final Groundwater Sampling Report for Tank 5 Release Response on June 23 and 24, 2014, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, July 2014.

ESI, 2014k, Third Quarter 2014 - Quarterly Groundwater Monitoring Report Inside Tunnel Wells, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, September 2014.

ESI, 2015, Fourth Quarter 2014 - Quarterly Groundwater Monitoring Report Inside Tunnel Wells, Red Hill Bulk Fuel Storage Facility, Pearl Harbor, Oahu, Hawaii, January 2015.

Foote et al., 1972, Soil Survey of the Islands of Kauai, Oahu, Maui, Molokai, and Lanai, State of Hawaii.

Mink, J. F. and Lau, L. S., 1990, Aquifer Identification and Classification for Oahu: Groundwater Protection Strategy for Hawaii: Water Resources Research Center Technical Report No. 179, February 1990.

Stearns, H. T. and Vaksvik, K. N., 1935, Geology and Groundwater Resources of the Island of Oahu, Hawaii: Hawaii Div. Hydrogr. Bull.

Stearns, H. T. and Vaksvik, K. N., 1938, Records of the Drilled Wells on the Island of Oahu, Hawaii: Hawaii Div. Hydrogr. Bull. 4, 213 p.

TEC, 2007, Final Technical Report, Red Hill Bulk Fuel Storage Facility, prepared for Naval Facilities Engineering Command, Pacific, Pearl Harbor, Hawaii, August 2007

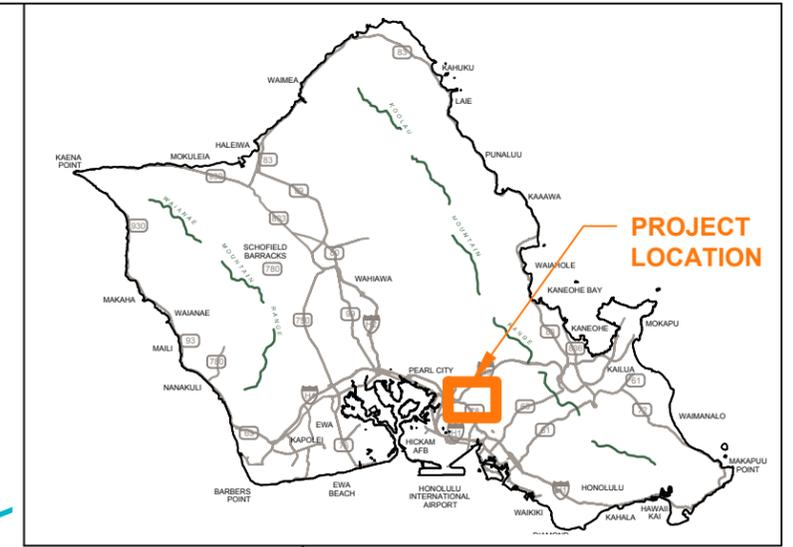
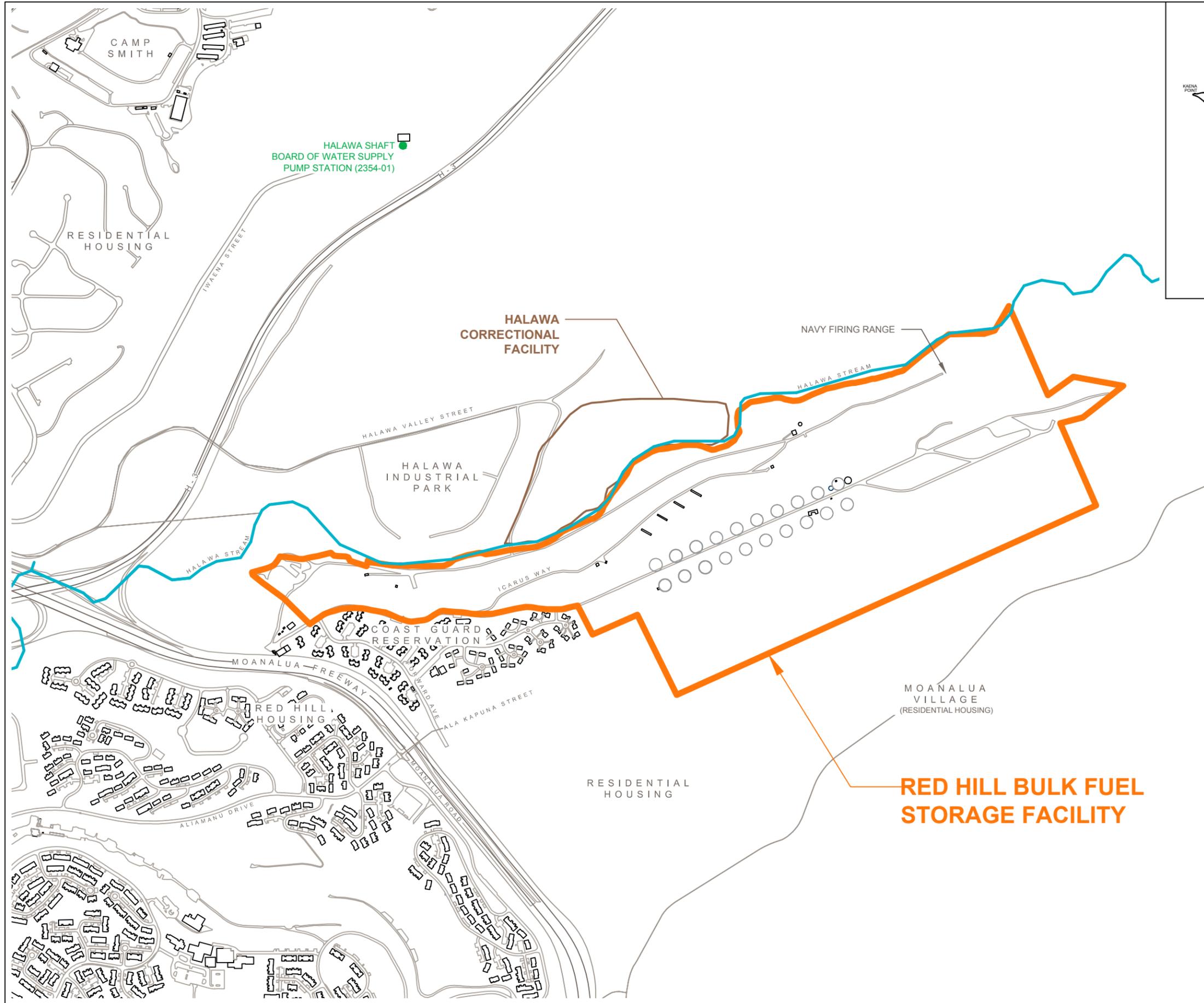
TEC, 2008, Final Groundwater Protection Plan, Red Hill Fuel Storage Facility, Prepared for Navy Region Hawaii, Pearl Harbor, Hawaii, January 2008, revised December 2009.

TEC, 2009, Quarterly Groundwater Monitoring Report, Red Hill Fuel Storage Facility, Prepared for Navy Region Hawaii, Pearl Harbor, Hawaii, September 2009.

This Page Intentionally Left Blank

FIGURES

This Page Intentionally Left Blank.



| |
|---|
| NOTES |
| The accuracy of this document is limited to the quality and scale of the source information. This document is not a legal representation of an engineered survey. |
| SOURCES |
| Pearl Harbor Base Map |
| Navy GIS files |

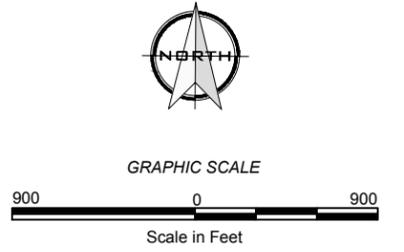
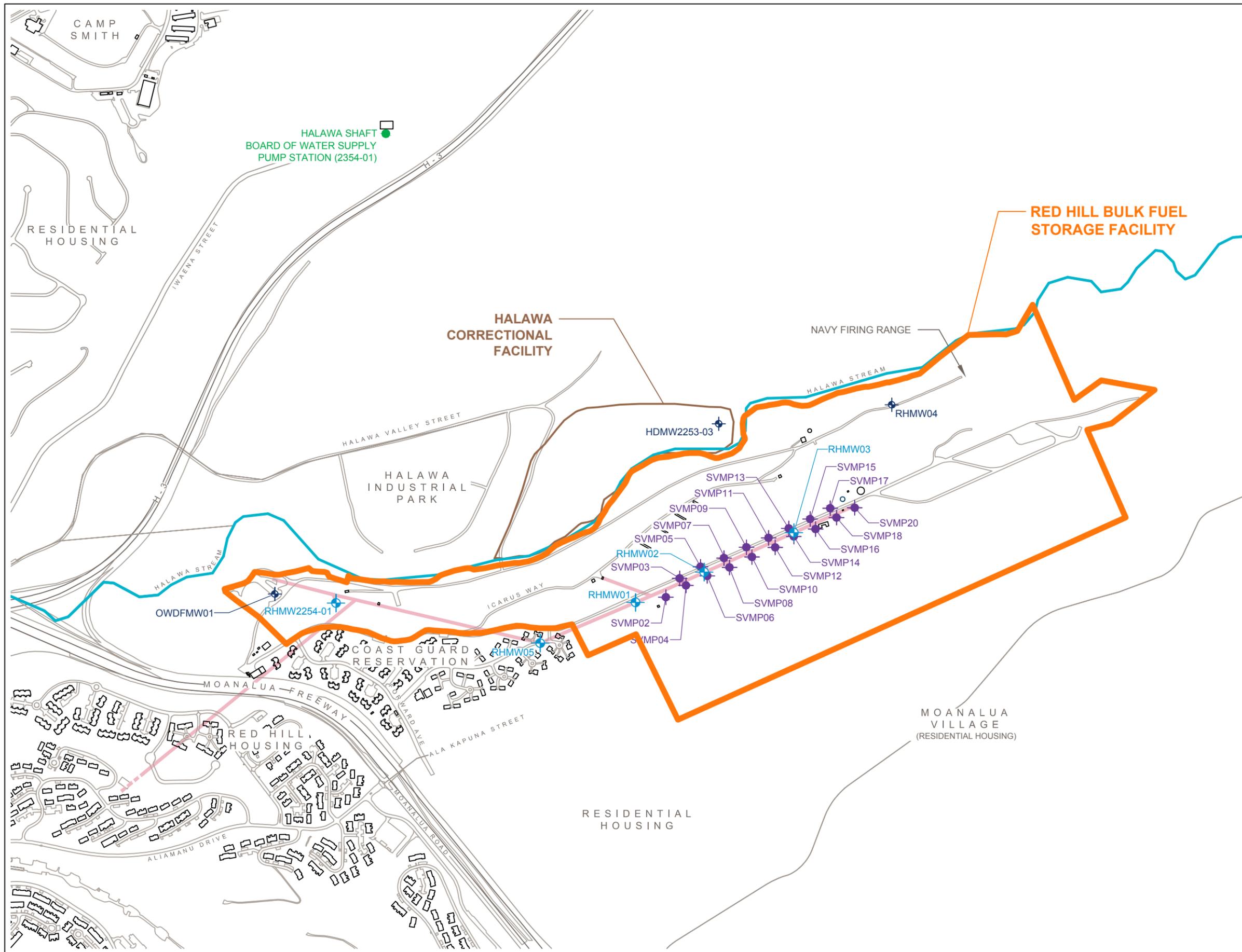


FIGURE 1
SITE LOCATION
 GROUNDWATER MONITORING
 RED HILL BULK FUEL STORAGE FACILITY
 NAVAL SUPPLY SYSTEM COMMAND (NAVSUP)
 FLEET LOGISTICS CENTER
 JBPHH, OAHU, HAWAII

This Page Intentionally Left Blank.



| LEGEND | |
|--------|--|
| | RED HILL BULK FUEL STORAGE FACILITY |
| | HALAWA CORRECTIONAL FACILITY |
| | HALAWA STREAM |
| | BUILDING |
| | ROAD |
| | ABOVEGROUND STORAGE TANK |
| | WATER TANK |
| | SOIL VAPOR MONITORING POINT |
| | GROUNDWATER MONITORING WELL LOCATED INSIDE TUNNEL |
| | GROUNDWATER MONITORING WELL LOCATED OUTSIDE TUNNEL |
| | BOARD OF WATER SUPPLY PUMP STATION |
| | TUNNEL |

NOTES

The accuracy of this document is limited to the quality and scale of the source information. This document is not a legal representation of an engineered survey.

SOURCES

Pearl Harbor Base Map
Navy GIS files

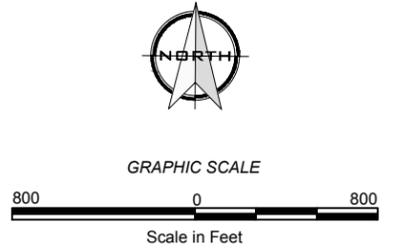


FIGURE 2
SITE LAYOUT
GROUNDWATER MONITORING
RED HILL BULK FUEL STORAGE FACILITY
NAVAL SUPPLY SYSTEM COMMAND (NAVSUP)
FLEET LOGISTICS CENTER
JBPHH, OAHU, HAWAII

This Page Intentionally Left Blank.

APPENDIX A

Groundwater Sampling Logs

This Page Intentionally Left Blank.



Groundwater Sampling Log

Well ID: RHMW01 Location: Red Hill Bulk Fuel Storage Facility Project No.: 112066

Initial Water Level: 83.63 ft Date: 1/27/2015 Time: 1205

Total Depth of Well: 97.35 ft Personnel Involved: Kirk Markle, Jeff Hattemer

Length of Saturated Zone: 13.72 ft Weather Conditions: Not applicable – well is located indoors

Volume of Water to be Removed: 2.5 L Method of Removal: Bladder Pump

Water Level After Purging: 83.63 ft Pumping Rate: 0.10 L/min

Well Purge Data:

| Time | Volume Removed | pH | Conductivity (mS/cm) | DO (mg/l) | Temperature | Salinity | Redox (ORP) (mV) |
|------|----------------|------|----------------------|-----------|-------------|----------|------------------|
| 1220 | 0.0 L | 7.74 | 0.322 | 8.96 | 24.67 | - | -29.9 |
| 1224 | 0.5 L | 7.27 | 0.318 | 4.67 | 24.31 | - | -62.2 |
| 1226 | 1.0 L | 7.07 | 0.317 | 2.56 | 24.19 | - | -72.2 |
| 1230 | 1.5 L | 6.99 | 0.318 | 2.32 | 24.10 | - | -73.5 |
| 1238 | 2.0 L | 6.90 | 0.319 | 2.29 | 24.04 | - | -75.7 |
| 1245 | 2.5 L | 6.90 | 0.320 | 2.39 | 24.27 | - | -72.8 |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |

Sample Withdrawal Method: Bladder Pump

Appearance of Sample:

Color: Clear
 Turbidity: None
 Sediment: None
 Other: None

Laboratory Analysis Parameters and Preservatives: TPH-d - 8015; TPH-g, VOCs - 8260; PAHs - 8270c sim; lead - 6020

Number and Types of Sample Containers: 6 - 40ml VOAs, 2 - 1L amber jar, 1 - 500ml amber jar, 1 - 250ml HDPE

Sample Identification Numbers: ES120X [1300]

Decontamination Procedures: Triple Rinsed

Notes: YSI did not have salinity parameter.

Sampled by: Kirk Markle, Jeff Hattemer

Sampled Delivered to: Calscience Environmental Lab Transporters: FedEx

Date: 1/27/2015 Time: 1500

Capacity of Casing (Gallons/Linear Feet)
2"-0.16 • 4"-0.65 • 8"-2.61 • 10"-4.08 • 12"-5.87



Groundwater Sampling Log

Well ID: RHMW02 Location: Red Hill Bulk Fuel Storage Facility Project No.: 112066

Initial Water Level: 86.35 ft Date: 1/28/2015 Time: 955

Total Depth of Well: 92.91 ft Personnel Involved: Justin Lam, Jeff Hattemer

Length of Saturated Zone: 6.56 ft Weather Conditions: Not applicable – well is located indoors

Volume of Water to be Removed: 5.0 L Method of Removal: Bladder Pump

Water Level After Purging: 86.81 ft Pumping Rate: 0.31 L/min

Well Purge Data:

| Time | Volume Removed | pH | Conductivity (mS/cm) | DO (mg/l) | Temperature | Salinity | Redox (ORP) (mV) |
|------|----------------|------|----------------------|-----------|-------------|----------|------------------|
| 1009 | 0.0 L | 7.17 | 0.554 | 4.47 | 24.94 | - | -61.0 |
| 1012 | 1.0 L | 7.08 | 0.552 | 1.51 | 24.12 | - | -105.4 |
| 1015 | 2.0 L | 6.98 | 0.553 | 1.40 | 23.85 | - | -113.1 |
| 1019 | 3.0 L | 6.96 | 0.557 | 1.34 | 23.77 | - | -116.7 |
| 1022 | 4.0 L | 6.96 | 0.562 | 1.30 | 23.76 | - | -119.1 |
| 1025 | 5.0 L | 6.96 | 0.563 | 1.28 | 23.76 | - | -119.8 |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |

Sample Withdrawal Method: Bladder Pump

Appearance of Sample:

Color: Clear
 Turbidity: Low
 Sediment: None
 Other: None

Laboratory Analysis Parameters and Preservatives: TPH-d - 8015; TPH-g, VOCs - 8260; PAHs - 8270c sim; lead - 6020

Number and Types of Sample Containers: 16 - 40ml VOAs, 6 - 1L amber jar, 4 - 500ml amber jar, 4 - 500ml HDPE

Sample Identification Numbers: ES126 [1035], ES127 MS/MSD [1035], ES127 (Dup) [1130]

Decontamination Procedures: Triple Rinsed

Notes: YSI did not have salinity parameter.

Sampled by: Justin Lam, Jeff Hattemer

Sampled Delivered to: Calscience Environmental Lab Transporters: FedEx

Date: 1/28/2015 Time: 1500

Capacity of Casing (Gallons/Linear Feet)
2"-0.16 • 4"-0.65 • 8"-2.61 • 10"-4.08 • 12"-5.87

This Page Intentionally Left Blank.

APPENDIX B

Field Notes

This Page Intentionally Left Blank.

Location RHSE
Project / Client NAVAC

Date 1/26/15

1350 Back @ office Drop off ice

JA
1/26/15

- tank fittings
- flexible tubing
- bucket lids
- gallon bags

Location RHSE
Project / Client NAVAC

Date 1/27/15

Purpose: GW sampling
Personnel: JJA, KM
0800 @ RHSE. Safety meeting
0830 @ Adit 3 pump house
0855 Gauge RHMW2254-01
DTW = 81.37
0903 Begin purge RHMW2254-01
0930 Collected sample ES125.
ES125 OK
1015 Set up @ RHMW05.
Leak in air line. Cut
and reattach fittings
Begin purge RHMW05
Collected sample ES124.
1205 Set up @ RHMW01
Gauge RHMW01
DTW =
Begin purge RHMW01
1300 Collected sample ES120
1410 Depart tunnel and pack
samples.
1430 Done cleaning up. Depart
Red Hill
1500 Dropped off samples @
Fed. GT. JA 1/27/15

Location RASF

Date 1/28/15

Project / Client NAVFAC

Purpose: SV, GW sampling

Personnel: JH, JL

0745 @ Red Hill, Calhoun
PID Safety meeting

0800 Entry @ Adit 3.

0850 @ RHMW03, C

0855 Gauge RHMW03, DTW = 102.63

0905 Begin purge RHMW03

0930 Collected sample ES123.

0950 @ RHMW02.

0955 Gauge RHMW02, DTW = 86.35

1009 Begin purge RHMW02

1035 Collected sample ES126,
ES126(MS/MSD), ES127 (labeled
as 1130)

1345 Out of tunnel

1415 Done cleaning up, packing
up, and dumping IDiv. Dept etc1430 Drop off samples at
Fed Ex.

1445 Back to office.

JH 1/28/15

Location RASF

Date 1/28/15

Project / Client NAVFAC

0908 SUMP20

| | | | | | |
|---|----|-----|-----|-----|-----|
| S | 72 | 112 | 109 | 112 | 101 |
| M | 75 | 80 | 54 | 80 | 72 |
| D | 80 | 60 | 89 | 89 | 80 |

0930 SUMP18

| | | | | | |
|---|------|------|------|------|------|
| S | 1032 | 1156 | 1263 | 1263 | 1179 |
| D | 1410 | 1757 | 1941 | 1941 | 1762 |

0938 SUMP17

| | | | | | |
|---|-----|-----|-----|-----|-----|
| S | 161 | 167 | 176 | 176 | 170 |
| M | 187 | 207 | 193 | 219 | 202 |
| D | 225 | 279 | 311 | 311 | 282 |

1000 SUMP16

| | | | | | |
|---|----|----|----|----|----|
| S | 60 | 54 | 69 | 67 | 63 |
| M | 14 | 20 | 20 | 23 | 19 |
| D | 46 | 49 | 60 | 60 | 54 |

1013 SUMP15

| | | | | | |
|---|-----|---------|-----------|----|----|
| S | 63 | 66 | 51 | 66 | 62 |
| M | mid | blocked | No sample | | |
| D | 43 | 80 | 80 | 80 | 71 |

1022 SUMP14

| | | | | | |
|---|----|----|----|----|----|
| S | 49 | 28 | 49 | 49 | 44 |
| M | 34 | 11 | 17 | 34 | 24 |
| D | 80 | 83 | 89 | 89 | 85 |

Location RHF

Date 1/28/15

Project / Client NAVFA C

| | | | | | |
|------|--------|-----|----|-----|----|
| 1032 | SVMP13 | | | | |
| S | 60 | 103 | 75 | 103 | 85 |
| M | 54 | 66 | 72 | 72 | 66 |
| D | 69 | 54 | 89 | 87 | 75 |

| | | | | | |
|------|--------|----|----|----|----|
| 1047 | SVMP12 | | | | |
| S | 17 | 11 | 14 | 17 | 15 |
| M | 5 | 60 | 57 | 60 | 46 |
| D | 40 | 40 | 86 | 86 | 63 |

| | | | | | |
|------|--------|-----|----|-----|----|
| 1058 | SVMP11 | | | | |
| S | | | | | |
| M | 100 | 100 | 95 | 100 | 99 |

D Deep bled. no sample.

| | | | | | |
|------|--------|----|----|----|----|
| 1110 | SVMP10 | | | | |
| S | 43 | 40 | 51 | 57 | 46 |
| D | 75 | 81 | 72 | 91 | 74 |

| | | | | | |
|------|--------|----|-----|-----|----|
| 1119 | SVMP09 | | | | |
| S | 43 | 34 | 34 | 51 | 41 |
| M | 28 | 31 | 57 | 57 | 43 |
| D | 46 | 63 | 115 | 124 | 87 |

| | | | | | |
|------|--------|-----|-----|-----|-----|
| 1130 | SVMP08 | | | | |
| S | 69 | 54 | 83 | 90 | 74 |
| M | 170 | 253 | 259 | 265 | 239 |
| D | 51 | 57 | 34 | 57 | 50 |

Location RHF

Date 1/28/15⁹⁹

Project / Client NAVFA C

| | | | | | |
|------|--------|------|------|------|----------|
| 1143 | SVMP07 | | | | |
| S | 164 | 216 | 225 | 225 | 208 |
| M | 109 | 106 | 106 | 111 | 108 |
| D | 17.3 | 21.1 | 21.3 | 21.4 | ppm 15.0 |

| | | | | | |
|------|--------|------|------|------|------|
| 1156 | SVMP06 | | | | |
| S | 1821 | 1996 | 2163 | 2224 | 2051 |
| M | 464 | 484 | 481 | 499 | 482 |

| | | | | | |
|-------|--------|------|-------|-----|----------|
| 12075 | SVMP05 | | | | |
| S | 81.5 | 93.6 | 100.3 | 116 | ppm 97.9 |
| M | 125 | 186 | 189 | 190 | ppm 172 |
| D | 193 | 216 | 213 | 217 | ppm 208 |

| | | | | | |
|------|--------|------|------|------|----------|
| 1225 | SVMP04 | | | | |
| S | 13.0 | 14.6 | 15.6 | 15.9 | ppm 14.8 |
| M | 11.9 | 13.4 | 14.0 | 14.4 | ppm 13.4 |
| D | 12.5 | 14.9 | 15.7 | 15.9 | ppm 14.8 |

| | | | | | |
|------|--------|------|------|------|----------|
| 1236 | SVMP03 | | | | |
| S | 6401 | 7953 | 8423 | 8871 | 8087 |
| M | 9673 | 10.3 | 10.6 | 10.6 | ppm 10.3 |
| D | 7311 | 7859 | 8565 | 8565 | 8075 |

| | | | | | |
|------|--------|------|------|------|------|
| 1249 | SVMP02 | | | | |
| S | 3467 | 3796 | 3914 | 4053 | 3808 |
| M | 1895 | 2135 | 2285 | 2285 | 2150 |
| D | 2189 | 2611 | 2660 | 2660 | 2530 |

SW 1/28/15

This Page Intentionally Left Blank.

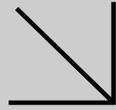
APPENDIX C

Laboratory Reports

This Page Intentionally Left Blank.



Calscience



WORK ORDER NUMBER: 15-01-1715

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For

Client: Environmental Science International, Inc.

Client Project Name: Red Hill LTM 112066

Attention: Jeff Hattemer
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Approved for release on 02/04/2015 by:
Terri Chang
Project Manager

ResultLink ▶

Email your PM ▶



Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.

Contents

Client Project Name: Red Hill LTM 112066
 Work Order Number: 15-01-1715

| | | |
|---|--|----|
| 1 | Work Order Narrative. | 3 |
| 2 | Client Sample Data. | 4 |
| | 2.1 EPA 8015B (M) TPH Diesel (Aqueous). | 4 |
| | 2.2 EPA 6020 ICP/MS Metals (Aqueous). | 5 |
| | 2.3 EPA 8270C SIM PAHs (Aqueous). | 6 |
| | 2.4 GC/MS GRO/EPA 8260B Volatile Organics (Aqueous). | 10 |
| 3 | Quality Control Sample Data. | 20 |
| | 3.1 MS/MSD. | 20 |
| | 3.2 PDS/PDSD. | 24 |
| | 3.3 LCS/LCSD. | 25 |
| 4 | Sample Analysis Summary. | 30 |
| 5 | Glossary of Terms and Qualifiers. | 31 |
| 6 | Chain-of-Custody/Sample Receipt Form. | 32 |

Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 01/28/15. They were assigned to Work Order 15-01-1715.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

New York NELAP air certification does not certify for all reported methods and analytes, reference the accredited items here: http://www.calscience.com/PDF/New_York.pdf

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.

The client requested the sample ID changed to ES120X on 01/29/2015 to differentiate the Q1 2015 ID numbers from the ones used for Q4 2014.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.



Calscience

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 3510C
Method: EPA 8015B (M)
Units: ug/L

Project: Red Hill LTM 112066

Page 1 of 1

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES125 | 15-01-1715-2-G | 01/27/15 09:30 | Aqueous | GC 45 | 01/29/15 | 01/30/15 10:49 | 150129B23A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
- TPH as Diesel is quantified in the carbon range C10-C28.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|---------------|--------|----|-----|-----|------|------------|
| TPH as Diesel | <12 | 11 | 12 | 25 | 1.00 | U |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|--------------|----------|----------------|------------|
| n-Octacosane | 77 | 51-141 | |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES124 | 15-01-1715-4-H | 01/27/15 11:15 | Aqueous | GC 45 | 01/29/15 | 01/30/15 11:06 | 150129B23A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
- TPH as Diesel is quantified in the carbon range C10-C28.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|---------------|--------|----|-----|-----|------|------------|
| TPH as Diesel | <13 | 12 | 13 | 26 | 1.00 | U |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|--------------|----------|----------------|------------|
| n-Octacosane | 74 | 51-141 | |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES120X | 15-01-1715-5-H | 01/27/15 13:00 | Aqueous | GC 45 | 01/29/15 | 01/30/15 11:24 | 150129B23A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
- TPH as Diesel is quantified in the carbon range C10-C28.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|---------------|--------|----|-----|-----|------|------------|
| TPH as Diesel | 33 | 11 | 12 | 25 | 1.00 | HD |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|--------------|----------|----------------|------------|
| n-Octacosane | 78 | 51-141 | |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-516-237 | N/A | Aqueous | GC 45 | 01/29/15 | 01/30/15 05:03 | 150129B23A |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|---------------|--------|----|-----|-----|------|------------|
| TPH as Diesel | <12 | 11 | 12 | 25 | 1.00 | U |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|--------------|----------|----------------|------------|
| n-Octacosane | 71 | 51-141 | |



Calscience

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: ug/L

Project: Red Hill LTM 112066

Page 1 of 1

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES124 | 15-01-1715-4-G | 01/27/15 11:15 | Aqueous | ICP/MS 03 | 01/29/15 | 01/29/15 22:50 | 150129L08F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|-----------|--------|--------|-------|-------|------|------------|
| Lead | <0.200 | 0.0898 | 0.200 | 0.500 | 1.00 | U |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES120X | 15-01-1715-5-G | 01/27/15 13:00 | Aqueous | ICP/MS 03 | 01/29/15 | 01/29/15 22:54 | 150129L08F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|-----------|--------|--------|-------|-------|------|------------|
| Lead | 0.631 | 0.0898 | 0.200 | 0.500 | 1.00 | |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-497-116 | N/A | Aqueous | ICP/MS 03 | 01/29/15 | 01/29/15 22:28 | 150129L08F |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|-----------|--------|--------|-------|-------|------|------------|
| Lead | <0.200 | 0.0898 | 0.200 | 0.500 | 1.00 | U |

Return to Contents

Analytical Report

Environmental Science International, Inc.
 354 Uluniu Street, Suite 304
 Kailua, HI 96734-2500

Date Received: 01/28/15
 Work Order: 15-01-1715
 Preparation: EPA 3510C
 Method: EPA 8270C SIM PAHs
 Units: ug/L

Project: Red Hill LTM 112066

Page 1 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES125 | 15-01-1715-2-I | 01/27/15 09:30 | Aqueous | GC/MS AAA | 01/30/15 | 02/03/15 00:28 | 150130L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|---------------------------|--------|-------|-------|------|------|------------|
| Naphthalene | <0.050 | 0.034 | 0.050 | 0.20 | 1.00 | U |
| 2-Methylnaphthalene | <0.050 | 0.046 | 0.050 | 0.20 | 1.00 | U |
| 1-Methylnaphthalene | <0.10 | 0.051 | 0.10 | 0.20 | 1.00 | U |
| Acenaphthylene | <0.050 | 0.044 | 0.050 | 0.20 | 1.00 | U |
| Acenaphthene | <0.050 | 0.027 | 0.050 | 0.20 | 1.00 | U |
| Fluorene | <0.050 | 0.042 | 0.050 | 0.20 | 1.00 | U |
| Phenanthrene | <0.050 | 0.027 | 0.050 | 0.20 | 1.00 | U |
| Anthracene | <0.050 | 0.029 | 0.050 | 0.20 | 1.00 | U |
| Fluoranthene | <0.050 | 0.047 | 0.050 | 0.20 | 1.00 | U |
| Pyrene | <0.050 | 0.020 | 0.050 | 0.20 | 1.00 | U |
| Benzo (a) Anthracene | <0.050 | 0.033 | 0.050 | 0.20 | 1.00 | U |
| Chrysene | <0.050 | 0.025 | 0.050 | 0.20 | 1.00 | U |
| Benzo (k) Fluoranthene | <0.050 | 0.031 | 0.050 | 0.20 | 1.00 | U |
| Benzo (b) Fluoranthene | <0.050 | 0.018 | 0.050 | 0.20 | 1.00 | U |
| Benzo (a) Pyrene | <0.050 | 0.022 | 0.050 | 0.20 | 1.00 | U |
| Indeno (1,2,3-c,d) Pyrene | <0.050 | 0.021 | 0.050 | 0.20 | 1.00 | U |
| Dibenz (a,h) Anthracene | <0.050 | 0.047 | 0.050 | 0.20 | 1.00 | U |
| Benzo (g,h,i) Perylene | <0.10 | 0.082 | 0.10 | 0.20 | 1.00 | U |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------|----------|----------------|------------|
| Nitrobenzene-d5 | 60 | 28-139 | |
| 2-Fluorobiphenyl | 56 | 33-144 | |
| p-Terphenyl-d14 | 63 | 23-160 | |

Analytical Report

Environmental Science International, Inc.
 354 Uluniu Street, Suite 304
 Kailua, HI 96734-2500

Date Received: 01/28/15
 Work Order: 15-01-1715
 Preparation: EPA 3510C
 Method: EPA 8270C SIM PAHs
 Units: ug/L

Project: Red Hill LTM 112066

Page 2 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES124 | 15-01-1715-4-I | 01/27/15 11:15 | Aqueous | GC/MS AAA | 01/30/15 | 02/03/15 00:48 | 150130L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|---------------------------|--------|-------|-------|------|------|------------|
| Naphthalene | <0.048 | 0.033 | 0.048 | 0.19 | 1.00 | U |
| 2-Methylnaphthalene | <0.048 | 0.045 | 0.048 | 0.19 | 1.00 | U |
| 1-Methylnaphthalene | <0.096 | 0.050 | 0.096 | 0.19 | 1.00 | U |
| Acenaphthylene | <0.048 | 0.043 | 0.048 | 0.19 | 1.00 | U |
| Acenaphthene | <0.048 | 0.026 | 0.048 | 0.19 | 1.00 | U |
| Fluorene | <0.048 | 0.041 | 0.048 | 0.19 | 1.00 | U |
| Phenanthrene | <0.048 | 0.026 | 0.048 | 0.19 | 1.00 | U |
| Anthracene | <0.048 | 0.028 | 0.048 | 0.19 | 1.00 | U |
| Fluoranthene | <0.048 | 0.045 | 0.048 | 0.19 | 1.00 | U |
| Pyrene | <0.048 | 0.020 | 0.048 | 0.19 | 1.00 | U |
| Benzo (a) Anthracene | <0.048 | 0.031 | 0.048 | 0.19 | 1.00 | U |
| Chrysene | <0.048 | 0.024 | 0.048 | 0.19 | 1.00 | U |
| Benzo (k) Fluoranthene | <0.048 | 0.030 | 0.048 | 0.19 | 1.00 | U |
| Benzo (b) Fluoranthene | <0.048 | 0.017 | 0.048 | 0.19 | 1.00 | U |
| Benzo (a) Pyrene | <0.048 | 0.021 | 0.048 | 0.19 | 1.00 | U |
| Indeno (1,2,3-c,d) Pyrene | <0.048 | 0.020 | 0.048 | 0.19 | 1.00 | U |
| Dibenz (a,h) Anthracene | <0.048 | 0.046 | 0.048 | 0.19 | 1.00 | U |
| Benzo (g,h,i) Perylene | <0.096 | 0.079 | 0.096 | 0.19 | 1.00 | U |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------|----------|----------------|------------|
| Nitrobenzene-d5 | 61 | 28-139 | |
| 2-Fluorobiphenyl | 61 | 33-144 | |
| p-Terphenyl-d14 | 67 | 23-160 | |



Calscience

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 3510C
Method: EPA 8270C SIM PAHs
Units: ug/L

Project: Red Hill LTM 112066

Page 3 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES120X | 15-01-1715-5-J | 01/27/15 13:00 | Aqueous | GC/MS AAA | 01/30/15 | 02/03/15 01:08 | 150130L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|---------------------------|--------|-------|-------|------|------|------------|
| Naphthalene | <0.054 | 0.037 | 0.054 | 0.22 | 1.00 | U |
| 2-Methylnaphthalene | <0.054 | 0.050 | 0.054 | 0.22 | 1.00 | U |
| 1-Methylnaphthalene | <0.11 | 0.056 | 0.11 | 0.22 | 1.00 | U |
| Acenaphthylene | <0.054 | 0.048 | 0.054 | 0.22 | 1.00 | U |
| Acenaphthene | <0.054 | 0.029 | 0.054 | 0.22 | 1.00 | U |
| Fluorene | <0.054 | 0.046 | 0.054 | 0.22 | 1.00 | U |
| Phenanthrene | <0.054 | 0.029 | 0.054 | 0.22 | 1.00 | U |
| Anthracene | <0.054 | 0.031 | 0.054 | 0.22 | 1.00 | U |
| Fluoranthene | <0.054 | 0.051 | 0.054 | 0.22 | 1.00 | U |
| Pyrene | <0.054 | 0.022 | 0.054 | 0.22 | 1.00 | U |
| Benzo (a) Anthracene | <0.054 | 0.035 | 0.054 | 0.22 | 1.00 | U |
| Chrysene | <0.054 | 0.027 | 0.054 | 0.22 | 1.00 | U |
| Benzo (k) Fluoranthene | <0.054 | 0.034 | 0.054 | 0.22 | 1.00 | U |
| Benzo (b) Fluoranthene | <0.054 | 0.019 | 0.054 | 0.22 | 1.00 | U |
| Benzo (a) Pyrene | <0.054 | 0.024 | 0.054 | 0.22 | 1.00 | U |
| Indeno (1,2,3-c,d) Pyrene | <0.054 | 0.023 | 0.054 | 0.22 | 1.00 | U |
| Dibenz (a,h) Anthracene | <0.054 | 0.052 | 0.054 | 0.22 | 1.00 | U |
| Benzo (g,h,i) Perylene | <0.11 | 0.089 | 0.11 | 0.22 | 1.00 | U |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------|----------|----------------|------------|
| Nitrobenzene-d5 | 51 | 28-139 | |
| 2-Fluorobiphenyl | 54 | 33-144 | |
| p-Terphenyl-d14 | 60 | 23-160 | |

Return to Contents

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 3510C
Method: EPA 8270C SIM PAHs
Units: ug/L

Project: Red Hill LTM 112066

Page 4 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-148-80 | N/A | Aqueous | GC/MS AAA | 01/30/15 | 02/02/15 21:28 | 150130L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|---------------------------|--------|-------|-------|------|------|------------|
| Naphthalene | <0.050 | 0.034 | 0.050 | 0.20 | 1.00 | U |
| 2-Methylnaphthalene | <0.050 | 0.046 | 0.050 | 0.20 | 1.00 | U |
| 1-Methylnaphthalene | <0.10 | 0.052 | 0.10 | 0.20 | 1.00 | U |
| Acenaphthylene | <0.050 | 0.045 | 0.050 | 0.20 | 1.00 | U |
| Acenaphthene | <0.050 | 0.027 | 0.050 | 0.20 | 1.00 | U |
| Fluorene | <0.050 | 0.043 | 0.050 | 0.20 | 1.00 | U |
| Phenanthrene | <0.050 | 0.027 | 0.050 | 0.20 | 1.00 | U |
| Anthracene | <0.050 | 0.029 | 0.050 | 0.20 | 1.00 | U |
| Fluoranthene | <0.050 | 0.047 | 0.050 | 0.20 | 1.00 | U |
| Pyrene | <0.050 | 0.020 | 0.050 | 0.20 | 1.00 | U |
| Benzo (a) Anthracene | <0.050 | 0.033 | 0.050 | 0.20 | 1.00 | U |
| Chrysene | <0.050 | 0.025 | 0.050 | 0.20 | 1.00 | U |
| Benzo (k) Fluoranthene | <0.050 | 0.031 | 0.050 | 0.20 | 1.00 | U |
| Benzo (b) Fluoranthene | <0.050 | 0.018 | 0.050 | 0.20 | 1.00 | U |
| Benzo (a) Pyrene | <0.050 | 0.022 | 0.050 | 0.20 | 1.00 | U |
| Indeno (1,2,3-c,d) Pyrene | <0.050 | 0.021 | 0.050 | 0.20 | 1.00 | U |
| Dibenz (a,h) Anthracene | <0.050 | 0.048 | 0.050 | 0.20 | 1.00 | U |
| Benzo (g,h,i) Perylene | <0.10 | 0.082 | 0.10 | 0.20 | 1.00 | U |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------|----------|----------------|------------|
| Nitrobenzene-d5 | 69 | 28-139 | |
| 2-Fluorobiphenyl | 67 | 33-144 | |
| p-Terphenyl-d14 | 69 | 23-160 | |



Calscience

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B
Units: ug/L

Project: Red Hill LTM 112066

Page 1 of 10

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ESTRIP | 15-01-1715-1-A | 01/27/15 09:00 | Aqueous | GC/MS OO | 01/29/15 | 01/29/15 18:29 | 150129L018 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|-----------------------------|--------|------|------|-----|------|------------|
| Acetone | <10 | 6.0 | 10 | 20 | 1.00 | U |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1.00 | U,IH |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1.00 | U |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1.00 | U |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1.00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1.00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1.00 | U |

Return to Contents



Calscience

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B
Units: ug/L

Project: Red Hill LTM 112066

Page 2 of 10

| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|-----------------|-----------------------|-------------------|------------|-----------|-------------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | <30 | 26 | 30 | 50 | 1.00 | U |
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> | | | |
| Dibromofluoromethane | 101 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 105 | 80-134 | | | | |
| Toluene-d8 | 97 | 80-120 | | | | |
| Toluene-d8-TPPH | 98 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 92 | 80-120 | | | | |


 Return to Contents

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B
Units: ug/L

Project: Red Hill LTM 112066

Page 3 of 10

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES125 | 15-01-1715-2-A | 01/27/15 09:30 | Aqueous | GC/MS OO | 01/29/15 | 01/29/15 21:09 | 150129L018 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|-----------------------------|--------|------|------|-----|------|------------|
| Acetone | <10 | 6.0 | 10 | 20 | 1.00 | U |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1.00 | U,IH |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1.00 | U |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1.00 | U |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1.00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1.00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1.00 | U |



Calscience

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B
Units: ug/L

Project: Red Hill LTM 112066

Page 4 of 10

| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|-----------------|-----------------------|-------------------|------------|-----------|-------------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | <30 | 26 | 30 | 50 | 1.00 | U |
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> | | | |
| Dibromofluoromethane | 101 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 105 | 80-134 | | | | |
| Toluene-d8 | 96 | 80-120 | | | | |
| Toluene-d8-TPPH | 98 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 92 | 80-120 | | | | |


 Return to Contents



Calscience

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B
Units: ug/L

Project: Red Hill LTM 112066

Page 5 of 10

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES124 | 15-01-1715-4-A | 01/27/15 11:15 | Aqueous | GC/MS OO | 01/29/15 | 01/29/15 21:36 | 150129L018 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|-----------------------------|--------|------|------|-----|------|------------|
| Acetone | <10 | 6.0 | 10 | 20 | 1.00 | U |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1.00 | U,IH |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1.00 | U |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1.00 | U |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1.00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1.00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1.00 | U |

Return to Contents



Calscience

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B
Units: ug/L

Project: Red Hill LTM 112066

Page 6 of 10

| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|-----------------|-----------------------|-------------------|------------|-----------|-------------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | <30 | 26 | 30 | 50 | 1.00 | U |
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> | | | |
| Dibromofluoromethane | 100 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 105 | 80-134 | | | | |
| Toluene-d8 | 96 | 80-120 | | | | |
| Toluene-d8-TPPH | 97 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 91 | 80-120 | | | | |



Return to Contents



Calscience

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B
Units: ug/L

Project: Red Hill LTM 112066

Page 7 of 10

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES120X | 15-01-1715-5-A | 01/27/15 13:00 | Aqueous | GC/MS OO | 01/29/15 | 01/29/15 22:03 | 150129L018 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|-----------------------------|--------|------|------|-----|------|------------|
| Acetone | <10 | 6.0 | 10 | 20 | 1.00 | U |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1.00 | U,IH |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1.00 | U |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1.00 | U |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1.00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1.00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1.00 | U |

Return to Contents

Analytical Report

Environmental Science International, Inc.
 354 Uluniu Street, Suite 304
 Kailua, HI 96734-2500

Date Received: 01/28/15
 Work Order: 15-01-1715
 Preparation: EPA 5030C
 Method: GC/MS / EPA 8260B
 Units: ug/L

Project: Red Hill LTM 112066

Page 8 of 10

| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|-----------------|-----------------------|-------------------|------------|-----------|-------------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | <30 | 26 | 30 | 50 | 1.00 | U |
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> | | | |
| Dibromofluoromethane | 102 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 106 | 80-134 | | | | |
| Toluene-d8 | 97 | 80-120 | | | | |
| Toluene-d8-TPPH | 98 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 92 | 80-120 | | | | |



Calscience

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B
Units: ug/L

Project: Red Hill LTM 112066

Page 9 of 10

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|----------------------|---------------------|----------------|-----------------|-----------------|---------------------------|-------------------|
| Method Blank | 099-13-057-75 | N/A | Aqueous | GC/MS OO | 01/29/15 | 01/29/15 17:31 | 150129L018 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|------------|-----------|-------------------|
| Acetone | <10 | 6.0 | 10 | 20 | 1.00 | U |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1.00 | U |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1.00 | U |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1.00 | U |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1.00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1.00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1.00 | U |

Return to Contents

Analytical Report

Environmental Science International, Inc.
 354 Uluniu Street, Suite 304
 Kailua, HI 96734-2500

Date Received: 01/28/15
 Work Order: 15-01-1715
 Preparation: EPA 5030C
 Method: GC/MS / EPA 8260B
 Units: ug/L

Project: Red Hill LTM 112066

Page 10 of 10

| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|-----------------|-----------------------|-------------------|------------|-----------|-------------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | <30 | 26 | 30 | 50 | 1.00 | U |
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> | | | |
| Dibromofluoromethane | 99 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 101 | 80-134 | | | | |
| Toluene-d8 | 96 | 80-120 | | | | |
| Toluene-d8-TPPH | 98 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 93 | 80-120 | | | | |



Calscience

Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 3005A Filt.
Method: EPA 6020

Project: Red Hill LTM 112066

Page 1 of 4

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| ES124 | Sample | Aqueous | ICP/MS 03 | 01/29/15 | 01/29/15 22:50 | 150129S08 |
| ES124 | Matrix Spike | Aqueous | ICP/MS 03 | 01/29/15 | 01/29/15 22:34 | 150129S08 |
| ES124 | Matrix Spike Duplicate | Aqueous | ICP/MS 03 | 01/29/15 | 01/29/15 22:37 | 150129S08 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Lead | ND | 100.0 | 109.8 | 110 | 107.5 | 107 | 80-120 | 2 | 0-20 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 3510C
Method: EPA 8270C SIM PAHs

Project: Red Hill LTM 112066

Page 2 of 4

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-01-1609-2 | Sample | Aqueous | GC/MS AAA | 01/30/15 | 02/02/15 22:08 | 150130S19 |
| 15-01-1609-2 | Matrix Spike | Aqueous | GC/MS AAA | 01/30/15 | 02/02/15 20:48 | 150130S19 |
| 15-01-1609-2 | Matrix Spike Duplicate | Aqueous | GC/MS AAA | 01/30/15 | 02/03/15 13:38 | 150130S19 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|---------------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Naphthalene | ND | 2.000 | 1.752 | 88 | 1.824 | 91 | 21-133 | 4 | 0-25 | |
| 2-Methylnaphthalene | ND | 2.000 | 1.713 | 86 | 1.802 | 90 | 21-140 | 5 | 0-25 | |
| 1-Methylnaphthalene | ND | 2.000 | 1.745 | 87 | 1.792 | 90 | 20-140 | 3 | 0-25 | |
| Acenaphthylene | ND | 2.000 | 1.756 | 88 | 1.756 | 88 | 33-145 | 0 | 0-25 | |
| Acenaphthene | ND | 2.000 | 1.836 | 92 | 1.863 | 93 | 49-121 | 1 | 0-25 | |
| Fluorene | ND | 2.000 | 1.830 | 92 | 1.810 | 90 | 59-121 | 1 | 0-25 | |
| Phenanthrene | ND | 2.000 | 1.877 | 94 | 1.874 | 94 | 54-120 | 0 | 0-25 | |
| Anthracene | ND | 2.000 | 1.539 | 77 | 1.632 | 82 | 27-133 | 6 | 0-25 | |
| Fluoranthene | ND | 2.000 | 1.853 | 93 | 1.820 | 91 | 26-137 | 2 | 0-25 | |
| Pyrene | ND | 2.000 | 2.018 | 101 | 1.897 | 95 | 18-168 | 6 | 0-25 | |
| Benzo (a) Anthracene | ND | 2.000 | 1.876 | 94 | 1.774 | 89 | 33-143 | 6 | 0-25 | |
| Chrysene | ND | 2.000 | 1.929 | 96 | 1.876 | 94 | 17-168 | 3 | 0-25 | |
| Benzo (k) Fluoranthene | ND | 2.000 | 1.951 | 98 | 1.869 | 93 | 24-159 | 4 | 0-25 | |
| Benzo (b) Fluoranthene | ND | 2.000 | 1.583 | 79 | 1.508 | 75 | 24-159 | 5 | 0-25 | |
| Benzo (a) Pyrene | ND | 2.000 | 1.673 | 84 | 1.621 | 81 | 17-163 | 3 | 0-25 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 2.000 | 1.726 | 86 | 1.722 | 86 | 10-171 | 0 | 0-25 | |
| Dibenz (a,h) Anthracene | ND | 2.000 | 1.811 | 91 | 1.796 | 90 | 10-219 | 1 | 0-25 | |
| Benzo (g,h,i) Perylene | ND | 2.000 | 1.627 | 81 | 1.716 | 86 | 10-227 | 5 | 0-25 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B

Project: Red Hill LTM 112066

Page 3 of 4

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| 15-01-1609-2 | Sample | Aqueous | GC/MS OO | 01/29/15 | 01/29/15 18:56 | 150129S039 |
| 15-01-1609-2 | Matrix Spike | Aqueous | GC/MS OO | 01/29/15 | 01/29/15 19:22 | 150129S039 |
| 15-01-1609-2 | Matrix Spike Duplicate | Aqueous | GC/MS OO | 01/29/15 | 01/29/15 19:49 | 150129S039 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Acetone | ND | 50.00 | 101.2 | 202 | 106.5 | 213 | 40-140 | 5 | 0-20 | 3 |
| Benzene | ND | 50.00 | 44.16 | 88 | 44.64 | 89 | 80-120 | 1 | 0-20 | |
| Bromodichloromethane | ND | 50.00 | 44.82 | 90 | 45.91 | 92 | 75-120 | 2 | 0-20 | |
| Bromoform | ND | 50.00 | 44.29 | 89 | 46.61 | 93 | 70-130 | 5 | 0-20 | |
| Bromomethane | ND | 50.00 | 60.69 | 121 | 49.43 | 99 | 30-145 | 20 | 0-20 | |
| 2-Butanone | ND | 50.00 | 61.40 | 123 | 65.63 | 131 | 30-150 | 7 | 0-20 | |
| Carbon Tetrachloride | ND | 50.00 | 39.91 | 80 | 41.28 | 83 | 65-140 | 3 | 0-20 | |
| Chlorobenzene | ND | 50.00 | 47.73 | 95 | 48.32 | 97 | 80-120 | 1 | 0-20 | |
| Chloroethane | ND | 50.00 | 41.79 | 84 | 41.02 | 82 | 60-135 | 2 | 0-20 | |
| Chloroform | ND | 50.00 | 46.18 | 92 | 46.34 | 93 | 65-135 | 0 | 0-20 | |
| Chloromethane | ND | 50.00 | 39.45 | 79 | 38.78 | 78 | 40-125 | 2 | 0-20 | |
| Dibromochloromethane | ND | 50.00 | 48.15 | 96 | 50.21 | 100 | 60-135 | 4 | 0-20 | |
| 1,2-Dibromo-3-Chloropropane | ND | 50.00 | 36.92 | 74 | 38.82 | 78 | 50-130 | 5 | 0-20 | |
| 1,2-Dibromoethane | ND | 50.00 | 50.57 | 101 | 52.07 | 104 | 80-120 | 3 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 47.95 | 96 | 48.23 | 96 | 70-120 | 1 | 0-20 | |
| 1,3-Dichlorobenzene | ND | 50.00 | 47.14 | 94 | 47.15 | 94 | 75-125 | 0 | 0-20 | |
| 1,4-Dichlorobenzene | ND | 50.00 | 47.60 | 95 | 47.53 | 95 | 75-125 | 0 | 0-20 | |
| 1,1-Dichloroethane | ND | 50.00 | 42.07 | 84 | 42.56 | 85 | 70-135 | 1 | 0-20 | |
| 1,2-Dichloroethane | ND | 50.00 | 51.51 | 103 | 52.61 | 105 | 70-130 | 2 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 42.04 | 84 | 42.35 | 85 | 70-130 | 1 | 0-20 | |
| c-1,2-Dichloroethene | ND | 50.00 | 44.23 | 88 | 45.20 | 90 | 70-125 | 2 | 0-20 | |
| t-1,2-Dichloroethene | ND | 50.00 | 40.55 | 81 | 41.70 | 83 | 60-140 | 3 | 0-20 | |
| 1,2-Dichloropropane | ND | 50.00 | 43.82 | 88 | 44.82 | 90 | 75-125 | 2 | 0-20 | |
| c-1,3-Dichloropropene | ND | 50.00 | 43.03 | 86 | 44.24 | 88 | 70-130 | 3 | 0-20 | |
| t-1,3-Dichloropropene | ND | 50.00 | 41.92 | 84 | 43.75 | 87 | 55-140 | 4 | 0-20 | |
| Ethylbenzene | ND | 50.00 | 45.53 | 91 | 46.40 | 93 | 75-125 | 2 | 0-20 | |
| Methylene Chloride | ND | 50.00 | 41.82 | 84 | 42.37 | 85 | 55-140 | 1 | 0-20 | |
| 4-Methyl-2-Pentanone | ND | 50.00 | 44.55 | 89 | 47.13 | 94 | 60-135 | 6 | 0-20 | |
| Styrene | ND | 50.00 | 44.83 | 90 | 45.51 | 91 | 65-135 | 1 | 0-20 | |
| 1,1,1,2-Tetrachloroethane | ND | 50.00 | 45.01 | 90 | 46.20 | 92 | 80-130 | 3 | 0-20 | |
| 1,1,2,2-Tetrachloroethane | ND | 50.00 | 0.2428 | 0 | 1.304 | 3 | 65-130 | 137 | 0-20 | 3,4 |
| Tetrachloroethene | ND | 50.00 | 68.54 | 137 | 69.48 | 139 | 45-150 | 1 | 0-20 | |
| Toluene | ND | 50.00 | 44.70 | 89 | 45.03 | 90 | 75-120 | 1 | 0-20 | |
| 1,2,4-Trichlorobenzene | ND | 50.00 | 47.30 | 95 | 46.60 | 93 | 65-135 | 1 | 0-20 | |
| 1,1,1-Trichloroethane | ND | 50.00 | 42.95 | 86 | 43.94 | 88 | 65-130 | 2 | 0-20 | |

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B

Project: Red Hill LTM 112066

Page 4 of 4

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>MS Conc.</u> | <u>MS %Rec.</u> | <u>MSD Conc.</u> | <u>MSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------|---------------------|--------------------|-----------------|-----------------|------------------|------------------|-----------------|------------|---------------|-------------------|
| Hexachloro-1,3-Butadiene | ND | 50.00 | 46.50 | 93 | 46.05 | 92 | 50-140 | 1 | 0-20 | |
| 1,1,2-Trichloroethane | ND | 50.00 | 46.30 | 93 | 48.01 | 96 | 75-125 | 4 | 0-20 | |
| Trichloroethene | ND | 50.00 | 77.69 | 155 | 78.02 | 156 | 70-125 | 0 | 0-20 | 3 |
| 1,2,3-Trichloropropane | ND | 50.00 | 38.71 | 77 | 39.92 | 80 | 75-125 | 3 | 0-20 | |
| Vinyl Chloride | ND | 50.00 | 43.44 | 87 | 42.43 | 85 | 50-145 | 2 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 90.64 | 91 | 90.75 | 91 | 75-130 | 0 | 0-20 | |
| o-Xylene | ND | 50.00 | 45.07 | 90 | 45.59 | 91 | 80-120 | 1 | 0-20 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 43.30 | 87 | 45.26 | 91 | 65-125 | 4 | 0-20 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - PDS

Environmental Science International, Inc.
 354 Uluniu Street, Suite 304
 Kailua, HI 96734-2500

Date Received: 01/28/15
 Work Order: 15-01-1715
 Preparation: EPA 3005A Filt.
 Method: EPA 6020

Project: Red Hill LTM 112066

Page 1 of 1

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | PDS/PDSD Batch Number |
|---------------------------|--------|---------|------------|----------------|----------------|-----------------------|
| ES124 | Sample | Aqueous | ICP/MS 03 | 01/29/15 00:00 | 01/29/15 22:50 | 150129S08 |
| ES124 | PDS | Aqueous | ICP/MS 03 | 01/29/15 00:00 | 01/29/15 22:40 | 150129S08 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>PDS Conc.</u> | <u>PDS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|---------------------|--------------------|------------------|------------------|-----------------|-------------------|
| Lead | ND | 100.0 | 111.4 | 111 | 75-125 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 3510C
Method: EPA 8015B (M)

Project: Red Hill LTM 112066

Page 1 of 5

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-516-237 | LCS | Aqueous | GC 45 | 01/29/15 | 01/30/15 05:20 | 150129B23A | | | |
| 099-15-516-237 | LCSD | Aqueous | GC 45 | 01/29/15 | 01/30/15 05:38 | 150129B23A | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| TPH as Diesel | 2000 | 1741 | 87 | 1692 | 85 | 60-132 | 3 | 0-11 | |

RPD: Relative Percent Difference. CL: Control Limits

Quality Control - LCS

Environmental Science International, Inc.
 354 Uluniu Street, Suite 304
 Kailua, HI 96734-2500

Date Received: 01/28/15
 Work Order: 15-01-1715
 Preparation: EPA 3005A Filt.
 Method: EPA 6020

Project: Red Hill LTM 112066

Page 2 of 5

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|--------------------|------------------------|------------------|-----------------------|-------------------|
| 099-14-497-116 | LCS | Aqueous | ICP/MS 03 | 01/29/15 | 01/29/15 22:31 | 150129L08F |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Lead | | 100.0 | 104.5 | 105 | 80-120 | |



Calscience

Quality Control - LCS

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 3510C
Method: EPA 8270C SIM PAHs

Project: Red Hill LTM 112066

Page 3 of 5

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|--------------------|------------------------|------------------|-----------------------|-------------------|
| 099-15-148-80 | LCS | Aqueous | GC/MS AAA | 01/30/15 | 02/03/15 03:08 | 150130L19 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Naphthalene | | 2.000 | 1.510 | 75 | 21-133 | |
| 2-Methylnaphthalene | | 2.000 | 1.537 | 77 | 21-140 | |
| 1-Methylnaphthalene | | 2.000 | 1.528 | 76 | 20-140 | |
| Acenaphthylene | | 2.000 | 1.458 | 73 | 33-145 | |
| Acenaphthene | | 2.000 | 1.547 | 77 | 55-121 | |
| Fluorene | | 2.000 | 1.454 | 73 | 59-121 | |
| Phenanthrene | | 2.000 | 1.536 | 77 | 54-120 | |
| Anthracene | | 2.000 | 1.262 | 63 | 27-133 | |
| Fluoranthene | | 2.000 | 1.595 | 80 | 26-137 | |
| Pyrene | | 2.000 | 1.784 | 89 | 45-129 | |
| Benzo (a) Anthracene | | 2.000 | 1.585 | 79 | 33-143 | |
| Chrysene | | 2.000 | 1.722 | 86 | 17-168 | |
| Benzo (k) Fluoranthene | | 2.000 | 1.970 | 98 | 24-159 | |
| Benzo (b) Fluoranthene | | 2.000 | 1.389 | 69 | 24-159 | |
| Benzo (a) Pyrene | | 2.000 | 1.473 | 74 | 17-163 | |
| Indeno (1,2,3-c,d) Pyrene | | 2.000 | 1.508 | 75 | 25-175 | |
| Dibenz (a,h) Anthracene | | 2.000 | 1.577 | 79 | 25-175 | |
| Benzo (g,h,i) Perylene | | 2.000 | 1.351 | 68 | 25-157 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B

Project: Red Hill LTM 112066

Page 4 of 5

| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | |
|-----------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|--------|------------|
| 099-13-057-75 | LCS | Aqueous | | GC/MS OO | 01/29/15 | 01/29/15 16:06 | 150129L018 | | |
| 099-13-057-75 | LCSD | Aqueous | | GC/MS OO | 01/29/15 | 01/29/15 16:33 | 150129L018 | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Acetone | 50.00 | 87.61 | 175 | N/A | N/A | 40-140 | N/A | 0-20 | X |
| Benzene | 50.00 | 46.83 | 94 | N/A | N/A | 80-120 | N/A | 0-20 | |
| Bromodichloromethane | 50.00 | 47.24 | 94 | N/A | N/A | 75-120 | N/A | 0-20 | |
| Bromoform | 50.00 | 46.74 | 93 | N/A | N/A | 70-130 | N/A | 0-20 | |
| Bromomethane | 50.00 | 53.74 | 107 | N/A | N/A | 30-145 | N/A | 0-20 | |
| 2-Butanone | 50.00 | 74.86 | 150 | N/A | N/A | 30-150 | N/A | 0-20 | |
| Carbon Tetrachloride | 50.00 | 45.16 | 90 | N/A | N/A | 65-140 | N/A | 0-20 | |
| Chlorobenzene | 50.00 | 49.91 | 100 | N/A | N/A | 80-120 | N/A | 0-20 | |
| Chloroethane | 50.00 | 46.95 | 94 | N/A | N/A | 60-135 | N/A | 0-20 | |
| Chloroform | 50.00 | 48.63 | 97 | N/A | N/A | 65-135 | N/A | 0-20 | |
| Chloromethane | 50.00 | 42.46 | 85 | N/A | N/A | 40-125 | N/A | 0-20 | |
| Dibromochloromethane | 50.00 | 50.11 | 100 | N/A | N/A | 60-135 | N/A | 0-20 | |
| 1,2-Dibromo-3-Chloropropane | 50.00 | 39.79 | 80 | N/A | N/A | 50-130 | N/A | 0-20 | |
| 1,2-Dibromoethane | 50.00 | 51.49 | 103 | N/A | N/A | 80-120 | N/A | 0-20 | |
| 1,2-Dichlorobenzene | 50.00 | 49.75 | 100 | N/A | N/A | 70-120 | N/A | 0-20 | |
| 1,3-Dichlorobenzene | 50.00 | 49.91 | 100 | N/A | N/A | 75-125 | N/A | 0-20 | |
| 1,4-Dichlorobenzene | 50.00 | 49.48 | 99 | N/A | N/A | 75-125 | N/A | 0-20 | |
| 1,1-Dichloroethane | 50.00 | 46.12 | 92 | N/A | N/A | 70-135 | N/A | 0-20 | |
| 1,2-Dichloroethane | 50.00 | 52.47 | 105 | N/A | N/A | 70-130 | N/A | 0-20 | |
| 1,1-Dichloroethene | 50.00 | 44.94 | 90 | N/A | N/A | 70-130 | N/A | 0-20 | |
| c-1,2-Dichloroethene | 50.00 | 47.74 | 95 | N/A | N/A | 70-125 | N/A | 0-20 | |
| t-1,2-Dichloroethene | 50.00 | 44.53 | 89 | N/A | N/A | 60-140 | N/A | 0-20 | |
| 1,2-Dichloropropane | 50.00 | 46.70 | 93 | N/A | N/A | 75-125 | N/A | 0-20 | |
| c-1,3-Dichloropropene | 50.00 | 47.08 | 94 | N/A | N/A | 70-130 | N/A | 0-20 | |
| t-1,3-Dichloropropene | 50.00 | 47.60 | 95 | N/A | N/A | 55-140 | N/A | 0-20 | |
| Ethylbenzene | 50.00 | 48.75 | 97 | N/A | N/A | 75-125 | N/A | 0-20 | |
| Methylene Chloride | 50.00 | 44.38 | 89 | N/A | N/A | 55-140 | N/A | 0-20 | |
| 4-Methyl-2-Pentanone | 50.00 | 45.89 | 92 | N/A | N/A | 60-135 | N/A | 0-20 | |
| Styrene | 50.00 | 47.01 | 94 | N/A | N/A | 65-135 | N/A | 0-20 | |
| 1,1,1,2-Tetrachloroethane | 50.00 | 48.20 | 96 | N/A | N/A | 80-130 | N/A | 0-20 | |
| 1,1,2,2-Tetrachloroethane | 50.00 | 40.07 | 80 | N/A | N/A | 65-130 | N/A | 0-20 | |
| Tetrachloroethene | 50.00 | 57.95 | 116 | N/A | N/A | 45-150 | N/A | 0-20 | |
| Toluene | 50.00 | 46.98 | 94 | N/A | N/A | 75-120 | N/A | 0-20 | |
| 1,2,4-Trichlorobenzene | 50.00 | 49.94 | 100 | N/A | N/A | 65-135 | N/A | 0-20 | |
| 1,1,1-Trichloroethane | 50.00 | 48.36 | 97 | N/A | N/A | 65-130 | N/A | 0-20 | |
| Hexachloro-1,3-Butadiene | 50.00 | 49.00 | 98 | N/A | N/A | 50-140 | N/A | 0-20 | |

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B

Project: Red Hill LTM 112066

Page 5 of 5

| <u>Parameter</u> | <u>Spike Added</u> | <u>LCS Conc.</u> | <u>LCS %Rec.</u> | <u>LCSD Conc.</u> | <u>LCSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------|--------------------|------------------|------------------|-------------------|-------------------|-----------------|------------|---------------|-------------------|
| 1,1,2-Trichloroethane | 50.00 | 48.87 | 98 | N/A | N/A | 75-125 | N/A | 0-20 | |
| Trichloroethene | 50.00 | 50.10 | 100 | N/A | N/A | 70-125 | N/A | 0-20 | |
| 1,2,3-Trichloropropane | 50.00 | 40.04 | 80 | N/A | N/A | 75-125 | N/A | 0-20 | |
| Vinyl Chloride | 50.00 | 47.19 | 94 | N/A | N/A | 50-145 | N/A | 0-20 | |
| p/m-Xylene | 100.0 | 95.89 | 96 | N/A | N/A | 75-130 | N/A | 0-20 | |
| o-Xylene | 50.00 | 47.74 | 95 | N/A | N/A | 80-120 | N/A | 0-20 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 47.26 | 95 | N/A | N/A | 65-125 | N/A | 0-20 | |
| Gasoline Range Organics | 1000 | 1077 | 108 | 1121 | 112 | 80-120 | 4 | 0-20 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits

Sample Analysis Summary Report

Work Order: 15-01-1715

Page 1 of 1

| <u>Method</u> | <u>Extraction</u> | <u>Chemist ID</u> | <u>Instrument</u> | <u>Analytical Location</u> |
|--------------------|-------------------|-------------------|-------------------|----------------------------|
| EPA 6020 | EPA 3005A Filtr. | 776 | ICP/MS 03 | 1 |
| EPA 8015B (M) | EPA 3510C | 682 | GC 45 | 1 |
| EPA 8270C SIM PAHs | EPA 3510C | 907 | GC/MS AAA | 1 |
| GC/MS / EPA 8260B | EPA 5030C | 849 | GC/MS OO | 2 |


Return to Contents

Location 1: 7440 Lincoln Way, Garden Grove, CA 92841

Location 2: 7445 Lampson Avenue, Garden Grove, CA 92841

| <u>Qualifiers</u> | <u>Definition</u> |
|-------------------|---|
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| B | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| DL | The Detection Limit (DL) is the smallest analyte concentration that can be demonstrated to be different from zero or a blank concentration at the 99% level of confidence. |
| E | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| ICH | Initial calibration verification recovery is above the control limit for this analyte. |
| ICJ | Initial calibration verification recovery is below the control limit for this analyte. |
| IH | Calibration verification recovery is above the control limit for this analyte. |
| IJ | Calibration verification recovery is below the control limit for this analyte. |
| J | Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| LOD | The Limit of Detection (LOD) is the smallest amount or concentration of a substance that must be present in a sample in order to be detected at 99% confidence level. |
| LOQ | The Limit of Quantitation (LOQ) is the lowest concentration of a substance that produces a quantitative result within specified limits of precision and bias. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| SG | The sample extract was subjected to Silica Gel treatment prior to analysis. |
| U | Undetected at Detection Limit (DL) and is reported as less than the Limit of Detection (LOD). |
| X | % Recovery and/or RPD out-of-range. |
| Z | Analyte presence was not confirmed by second column or GC/MS analysis. |
| | Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis. |
| | Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time. |
| | A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations. |



Calscience

7440 Lincoln Way, Garden Grove, CA 92841-1427 • (714) 895-5494
For courier service / sample drop off information, contact us28_sales@eurofins.com or call us.

LABORATORY CLIENT: ESI

ADDRESS: 354 Uluniv St., Suite 304

CITY: Kailua STATE: HI ZIP: 96734

TEL: 808-261-0740 E-MAIL: jhatter@escience.com, d@escience.com

TURNAROUND TIME (Rush surcharges may apply to any VAI not 'STANDARD'):
 SAME DAY 24 HR 48 HR 72 HR 5 DAYS STANDARD

EDD: COELT EDF OTHER

SPECIAL INSTRUCTIONS:

WO NO. / LAB USE ONLY
15-01-1715

CLIENT PROJECT NAME / NO.: Red Hill LTM 112066

PROJECT CONTACT: Jeff Hattermer

GLOBAL ID:
com

LOG CODE:

SAMPLER(S): (PRINT)

P.O. NO.:

LAB CONTACT OR QUOTE NO.:

REQUESTED ANALYSES

Please check box or fill in blank as needed.

| LAB USE ONLY | SAMPLE ID | SAMPLING DATE | SAMPLING TIME | MATRIX | NO. OF CONT. | Unpreserved | Preserved | Field Filtered | TPH (g) <input type="checkbox"/> C60 (8260) | TPH (d) <input type="checkbox"/> BPE (8015) | TPH <input type="checkbox"/> C6-C36 <input type="checkbox"/> C6-C44 | TPH | BTEX / MTBE <input type="checkbox"/> 8260 <input type="checkbox"/> | VOCs (8260) | Oxygenates (8260) | Prep (5035) <input type="checkbox"/> En Core <input type="checkbox"/> Terra Core | SVOCs (8270) | Pesticides (8081) | PCBs (8082) | PAHs <input type="checkbox"/> 8270 <input checked="" type="checkbox"/> 8270 SIM | T22 Metals <input type="checkbox"/> 6010/747X <input type="checkbox"/> 6020/747X | Cr(VI) <input type="checkbox"/> 7196 <input type="checkbox"/> 7199 <input type="checkbox"/> 218.6 | Lead (6012) | Lead (200.8) | |
|--------------|-----------|---------------|---------------|--------|--------------|-------------------------------------|--------------------------|--------------------------|---|---|---|--------------------------|--|-------------------------------------|--------------------------|--|--------------------------|--------------------------|--------------------------|---|--|---|--------------------------|--------------------------|--------------------------|
| 1 | ESTRIP | 1/27/15 | 0900 | water | 3 | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| 2 | ES 125 | 1/27/15 | 0930 | water | 9 | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| 3 | ES 125 UF | 1/27/15 | 0930 | water | 1 | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| 4 | ES 124 | 1/27/15 | 1115 | water | 10 | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| 5 | ES 120 | 1/27/15 | 1300 | water | 10 | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

| | | |
|---|---------------|------------|
| Received by: (Signature/Affiliation) Fed Ex | Date: 1/27/15 | Time: 1500 |
| Received by: (Signature/Affiliation) [Signature] | Date: 1/28/15 | Time: 1020 |
| Received by: (Signature/Affiliation) | Date: | Time: |

Relinquished by: (Signature) [Signature] SASH Hattermer

Relinquished by: (Signature)

Relinquished by: (Signature)

ORIGIN ID:HNLA (714) 895-5494
CALSCIENCE ENVIRONMENTAL LAB
7440 LINCOLN WAY
GARDEN GROVE, CA 928411427
UNITED STATES US

SHIP DATE: 27JAN15
ACTWGT: 55.6 LB
CAD: /POS1525
DIMS: 23x15x14 IN
BILL SENDER

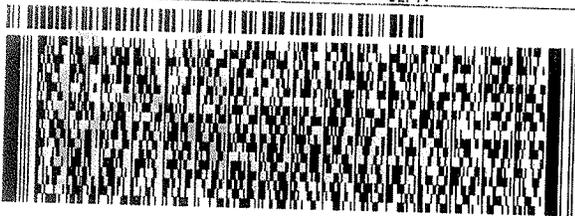
17LS

TO **SAMPLE CONTROL**
CALSCIENCE LABORTORIES
7440 LINCOLN WAY

GARDEN GROVE CA 92841

(714) 895-5494 REF: DEPT:
INU:
PO:

SPR1073 01/27/15 53711/REF15/FEAR

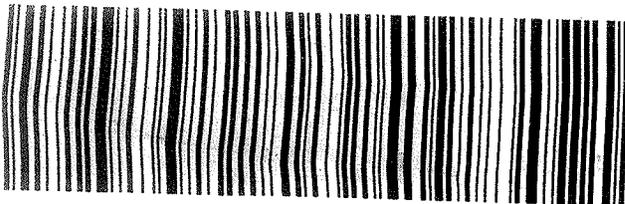


WED - 28 JAN AA
STANDARD OVERNIGHT

TRK# 8059 2709 6788
0200

WZ APVA

92841
CA-US SNA



Open End of FedEx Pouch Here

Calscience

WORK ORDER #: 15-01-1715

SAMPLE RECEIPT FORM

Cooler 1 of 1

CLIENT: ESI

DATE: 01/28/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 2.7 °C + 0.2 °C (CF) = 2.9 °C Blank Sample

Sample(s) outside temperature criteria (PM/APM contacted by: _____)

Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: Air Filter Checked by: LS

CUSTODY SEALS INTACT:

Cooler _____ No (Not Intact) Not Present N/A Checked by: LS

Sample _____ No (Not Intact) Not Present Checked by: LS

| SAMPLE CONDITION: | Yes | No | N/A |
|---|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-Of-Custody (COC) document(s) received with samples..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Collection date/time, matrix, and/or # of containers logged in based on sample labels. <input type="checkbox"/> No analysis requested. <input type="checkbox"/> Not relinquished. <input type="checkbox"/> No date/time relinquished. | | | |
| Sampler's name indicated on COC..... | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and good condition..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers and sufficient volume for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Analyses received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples received within 15-minute holding time | | | |
| <input type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfides <input type="checkbox"/> Dissolved Oxygen..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| Proper preservation noted on COC or sample container..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Unpreserved vials received for Volatiles analysis | | | |
| Volatile analysis container(s) free of headspace..... | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Tedlar bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

Solid: 4ozCGJ 8ozCGJ 16ozCGJ Sleeve (____) EnCores® TerraCores® _____

Aqueous: VOA VOA⁶h VOAna₂ 125AGB 125AGBh 125AGBp 1AGB 1AGBna₂ 1AGBs

500AGB 500AGJ 500AGJs 250AGB 250CGB 250CGBs 1PB 1PBna 500PB

250PB 250PBn 125PB 125PBz³⁽⁻¹⁾na 100PJ 100PJna₂ 250PBn² _____ _____

Air: Tedlar® Canister Other: _____ Trip Blank Lot#: 140908B Labeled/Checked by: LS

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope Reviewed by: LS

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure z^{na}: ZnAc₂+NaOH f: Filtered Scanned by: LS

Return to Contents

Terri Chang

From: Domonkos Feher [DFeher@esciencei.com]
Sent: Thursday, January 29, 2015 11:41 AM
To: Terri Chang
Cc: Jeff Hattemer; Ann Dang; Scott Simmons
Subject: 15-01-1715 and 15-01-1609

Terri,

The following Sample ID's have already been used for the Q4 2014 monitoring event:

Sample receipt 15-01-1715:

ES120

Sample receipt 15-01-1609:

ES121 (ES121 MS/MSD)

ES122

Please append an "X" to the Sample IDs in the final report, to differentiate the Q1 2015 ID numbers from the ones used for Q4 2014. The new sample IDs should read:

ES120X

ES121X (ES121X MS/MSD)

ES122X

Please include this email in the COC sections of the final laboratory reports as a record for the sample ID change. Mention the change in the case narratives.

Thank you,

Domonkos Fehér, Ph.D.

Project Chemist

Environmental Science International, Inc.

354 Uluniu Street, Suite 304, Kailua, Hawaii 96734

Cell: (808) 232-1261, Office: (808) 261-0740 ext. 118;

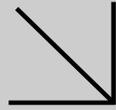
Email: dfeyer@esciencei.com

Notify us [here](#) to report this email as spam.



Environmental
Calscience

Supplemental Report 1



WORK ORDER NUMBER: 15-01-1715

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For

Client: Environmental Science International, Inc.

Client Project Name: Red Hill LTM 112066

Attention: Jeff Hattemer
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Approved for release on 02/04/2015 by:
Terri Chang
Project Manager

ResultLink ▶

Email your PM ▶



Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.



Contents

Client Project Name: Red Hill LTM 112066
Work Order Number: 15-01-1715

| | | |
|---|--|---|
| 1 | Work Order Narrative. | 3 |
| 2 | Client Sample Data. | 4 |
| | 2.1 EPA 200.8 ICP/MS Metals (Aqueous). | 4 |
| 3 | Quality Control Sample Data. | 5 |
| | 3.1 MS/MSD. | 5 |
| | 3.2 LCS/LCSD. | 6 |
| 4 | Sample Analysis Summary. | 7 |
| 5 | Glossary of Terms and Qualifiers. | 8 |
| 6 | Chain-of-Custody/Sample Receipt Form. | 9 |

Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 01/28/15. They were assigned to Work Order 15-01-1715.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

New York NELAP air certification does not certify for all reported methods and analytes, reference the accredited items here: http://www.calscience.com/PDF/New_York.pdf

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

Analytical Report

Environmental Science International, Inc.
 354 Uluniu Street, Suite 304
 Kailua, HI 96734-2500

Date Received: 01/28/15
 Work Order: 15-01-1715
 Preparation: N/A
 Method: EPA 200.8
 Units: ug/L

Project: Red Hill LTM 112066

Page 1 of 1

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES125UF | 15-01-1715-3-A | 01/27/15 09:30 | Aqueous | ICP/MS 03 | 01/29/15 | 01/30/15 00:35 | 150129L07 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|---------|------|--------|------|------------|
| Lead | <0.0898 | 1.00 | 0.0898 | 1.00 | U |

| Method Blank | 099-16-094-697 | N/A | Aqueous | ICP/MS 03 | 01/29/15 | 01/29/15 23:29 | 150129L07 |
|--------------|----------------|-----|---------|-----------|----------|-------------------|-----------|
|--------------|----------------|-----|---------|-----------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | RL | MDL | DF | Qualifiers |
|-----------|---------|------|--------|------|------------|
| Lead | <0.0898 | 1.00 | 0.0898 | 1.00 | U |



Calscience

Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: N/A
Method: EPA 200.8

Project: Red Hill LTM 112066

Page 1 of 1

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| ES125UF | Sample | Aqueous | ICP/MS 03 | 01/29/15 | 01/30/15 00:35 | 150129S07 |
| ES125UF | Matrix Spike | Aqueous | ICP/MS 03 | 01/29/15 | 01/30/15 00:18 | 150129S07 |
| ES125UF | Matrix Spike Duplicate | Aqueous | ICP/MS 03 | 01/29/15 | 01/30/15 00:22 | 150129S07 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Lead | ND | 100.0 | 111.0 | 111 | 110.1 | 110 | 80-120 | 1 | 0-20 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits

Quality Control - LCS

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/28/15
Work Order: 15-01-1715
Preparation: N/A
Method: EPA 200.8

Project: Red Hill LTM 112066

Page 1 of 1

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|--------------------|------------------------|------------------|-----------------------|-------------------|
| 099-16-094-697 | LCS | Aqueous | ICP/MS 03 | 01/29/15 | 01/29/15 23:59 | 150129L07 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Lead | | 100.0 | 103.9 | 104 | 80-120 | |



Calscience

Sample Analysis Summary Report

Work Order: 15-01-1715

Page 1 of 1

| <u>Method</u> | <u>Extraction</u> | <u>Chemist ID</u> | <u>Instrument</u> | <u>Analytical Location</u> |
|---------------|-------------------|-------------------|-------------------|----------------------------|
| EPA 200.8 | N/A | 776 | ICP/MS 03 | 1 |


Return to Contents

Location 1: 7440 Lincoln Way, Garden Grove, CA 92841

| <u>Qualifiers</u> | <u>Definition</u> |
|-------------------|---|
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| B | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| DL | The Detection Limit (DL) is the smallest analyte concentration that can be demonstrated to be different from zero or a blank concentration at the 99% level of confidence. |
| E | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| ICH | Initial calibration verification recovery is above the control limit for this analyte. |
| ICJ | Initial calibration verification recovery is below the control limit for this analyte. |
| IH | Calibration verification recovery is above the control limit for this analyte. |
| IJ | Calibration verification recovery is below the control limit for this analyte. |
| J | Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| LOD | The Limit of Detection (LOD) is the smallest amount or concentration of a substance that must be present in a sample in order to be detected at 99% confidence level. |
| LOQ | The Limit of Quantitation (LOQ) is the lowest concentration of a substance that produces a quantitative result within specified limits of precision and bias. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| SG | The sample extract was subjected to Silica Gel treatment prior to analysis. |
| U | Undetected at Detection Limit (DL) and is reported as less than the Limit of Detection (LOD). |
| X | % Recovery and/or RPD out-of-range. |
| Z | Analyte presence was not confirmed by second column or GC/MS analysis. |
| | Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis. |
| | Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time. |
| | A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations. |



Calscience

7440 Lincoln Way, Garden Grove, CA 92841-1427 • (714) 895-5494
For courier service / sample drop off information, contact us28_sales@eurofins.com or call us.

LABORATORY CLIENT: ESI

ADDRESS: 354 Uluniv St, Suite 304

CITY: Kailua STATE: HI ZIP: 96734

TEL: 808-261-0740 E-MAIL: jhatter@escience.com, d@escience.com

TURNAROUND TIME (Rush surcharges may apply to any VAI not 'STANDARD'):
 SAME DAY 24 HR 48 HR 72 HR 5 DAYS STANDARD

EDD: COELT EDF OTHER

SPECIAL INSTRUCTIONS:

WO NO. / LAB USE ONLY
15-01-1715

CLIENT PROJECT NAME / NO.:

Red Hill LTM 112066

PROJECT CONTACT:

Jeff Hattermer

GLOBAL ID:

com

LOG CODE:

SAMPLER(S): (PRINT)

P.O. NO.:

LAB CONTACT OR QUOTE NO.:

REQUESTED ANALYSES

Please check box or fill in blank as needed.

| LAB USE ONLY | SAMPLE ID | SAMPLING DATE | SAMPLING TIME | MATRIX | NO. OF CONT. | Unpreserved | Preserved | Field Filtered | TPH (g) <input type="checkbox"/> GPO (8260) | TPH (d) <input type="checkbox"/> DPE (8015) | TPH <input type="checkbox"/> C6-C36 <input type="checkbox"/> C6-C44 | TPH | BTEX / MTBE <input type="checkbox"/> 8260 <input type="checkbox"/> | VOCs (8260) | Oxygenates (8260) | Prep (5035) <input type="checkbox"/> En Core <input type="checkbox"/> Terra Core | SVOCs (8270) | Pesticides (8081) | PCBs (8082) | PAHs <input type="checkbox"/> 8270 <input checked="" type="checkbox"/> 8270 SIM | T22 Metals <input type="checkbox"/> 6010/747X <input type="checkbox"/> 6020/747X | Cr(VI) <input type="checkbox"/> 7196 <input type="checkbox"/> 7199 <input type="checkbox"/> 218.6 | Lead (6012) | Lead (200.8) | |
|--------------|-----------|---------------|---------------|--------|--------------|-------------------------------------|--------------------------|--------------------------|---|---|---|--------------------------|--|-------------------------------------|--------------------------|--|--------------------------|--------------------------|--------------------------|---|--|---|--------------------------|--------------------------|--------------------------|
| 1 | ES121P | 1/27/15 | 0900 | water | 3 | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| 2 | ES125 | 1/27/15 | 0930 | water | 9 | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| 3 | ES125 UF | 1/27/15 | 0930 | water | 1 | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| 4 | ES124 | 1/27/15 | 1115 | water | 10 | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| 5 | ES120 | 1/27/15 | 1300 | water | 10 | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

| | | |
|--|---------------|------------|
| Received by: (Signature/Affiliation) Fed Ex | Date: 1/27/15 | Time: 1500 |
| Received by: (Signature/Affiliation) | Date: 1/28/15 | Time: 1020 |
| Received by: (Signature/Affiliation) | Date: | Time: |

Relinquished by: (Signature) *Jeff Hattermer*

Relinquished by: (Signature)

Relinquished by: (Signature)

ORIGIN ID:HNLA (714) 895-5494
CALSCIENCE ENVIRONMENTAL LAB
7440 LINCOLN WAY

SHIP DATE: 27JAN15
ACTWGT: 55.6 LB
CAD: /POS1525
DIMS: 23x15x14 IN
BILL SENDER

GARDEN GROVE, CA 928411427
UNITED STATES US

TO **SAMPLE CONTROL**
CALSCIENCE LABORTORIES
7440 LINCOLN WAY

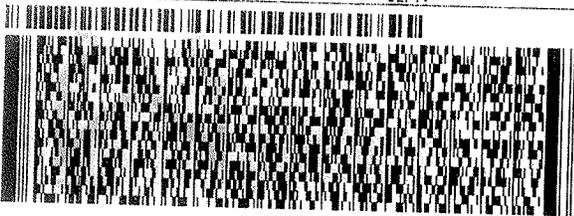
GARDEN GROVE CA 92841

(714) 895-5494
TNU:
PO:

REF:

DEPT:

SPR1073 01/27/15 53711/REF15/FEAR



FedEx
Express



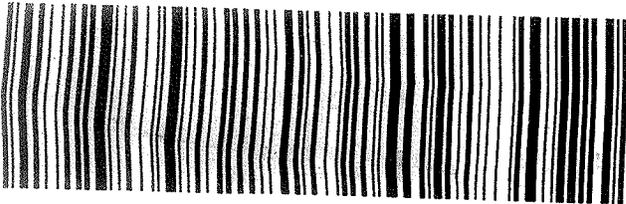
J16101501140110

TRK# 8059 2709 6788
0200

WED - 28 JAN AA
STANDARD OVERNIGHT

WZ APVA

92841
CA-US SNA



Open End of FedEx Pouch Here

17LS

SAMPLE RECEIPT FORM

Cooler 1 of 1

CLIENT: ESI

DATE: 01/28/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 2.7 °C + 0.2 °C (CF) = 2.9 °C Blank Sample

Sample(s) outside temperature criteria (PM/APM contacted by: _____)

Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: Air Filter Checked by: LS

CUSTODY SEALS INTACT:

Cooler _____ No (Not Intact) Not Present N/A Checked by: LS

Sample _____ No (Not Intact) Not Present Checked by: LS

| SAMPLE CONDITION: | Yes | No | N/A |
|---|-------------------------------------|-------------------------------------|-------------------------------------|
| Chain-Of-Custody (COC) document(s) received with samples..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Collection date/time, matrix, and/or # of containers logged in based on sample labels. <input type="checkbox"/> No analysis requested. <input type="checkbox"/> Not relinquished. <input type="checkbox"/> No date/time relinquished. | | | |
| Sampler's name indicated on COC..... | <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and good condition..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers and sufficient volume for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Analyses received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples received within 15-minute holding time | | | |
| <input type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfides <input type="checkbox"/> Dissolved Oxygen..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| Proper preservation noted on COC or sample container..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Unpreserved vials received for Volatiles analysis | | | |
| Volatile analysis container(s) free of headspace..... | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Tedlar bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

Solid: 4ozCGJ 8ozCGJ 16ozCGJ Sleeve (____) EnCores® TerraCores® _____

Aqueous: VOA VOA³⁽⁻¹⁾h VOAna₂ 125AGB 125AGBh 125AGBp 1AGB 1AGBna₂ 1AGBs

500AGB 500AGJ 500AGJs 250AGB 250CGB 250CGBs 1PB 1PBna 500PB

250PB 250PBn 125PB 125PBz^{na} 100PJ 100PJna₂ 250PBn^f _____ _____

Air: Tedlar® Canister Other: _____ Trip Blank Lot#: 140908B Labeled/Checked by: LS

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope Reviewed by: LS

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure z^{na}: ZnAc₂+NaOH f: Filtered Scanned by: LS

Return to Contents

RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER: 15-01-1715
INSTRUMENT: GC 45
EXTRACTION: EPA 3510C
D/T EXTRACTED: 2015-01-29 00:00

ANALYZED BY: 682
D/T ANALYZED: 2015-01-30 11:24
REVIEWED BY:
D/T REVIEWED:

DATA FILE: W:\GC_45\GC 45 DATA\2015\150129\15012974.D\15012974

5 **CLIENT SAMPLE NUMBER: ES120X**

| | |
|---------------------------------|---|
| LCS/MB BATCH: 150129B23A | SAMPLE VOLUME / WEIGHT: DEFAULT: 500.00 ml / ACTUAL: 500.00 ml |
| MS/MSD BATCH: | FINAL VOLUME / WEIGHT: DEFAULT: 5.00 ml / ACTUAL: 2.50 ml |
| UNITS: ug/L | ADJUSTMENT RATIO TO PF: 0.50 |

COMMENT: Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

| <u>COMPOUND</u> | <u>INI. CONC</u> | <u>DF</u> | <u>CONC</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>QUAL</u> |
|-----------------|------------------|-----------|-------------|-----------|------------|------------|-------------|
| TPH as Diesel | 6540 | 1.00 | 32.7 | 11 | 12 | 25 | b |

=====
 Area Percent Report
 =====

Data File Name : W:\GC_45\GC 45 DATA\2015\150129\15012974.D
 Page Number : 1
 Operator : Vial Number : Vial 74
 Instrument : GC 45 Injection Number : 1
 Sample Name : 15-01-1715-5 Sequence Line : 74
 Instrument Method: C:\CHEM32\1\METHODS\ ->
 Acquired on : 30 Jan 15 11:24 am
 Report Created on: 02 Feb 15 10:24 am Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Sig. 1 in W:\GC_45\GC 45 DATA\2015\150129\ ->

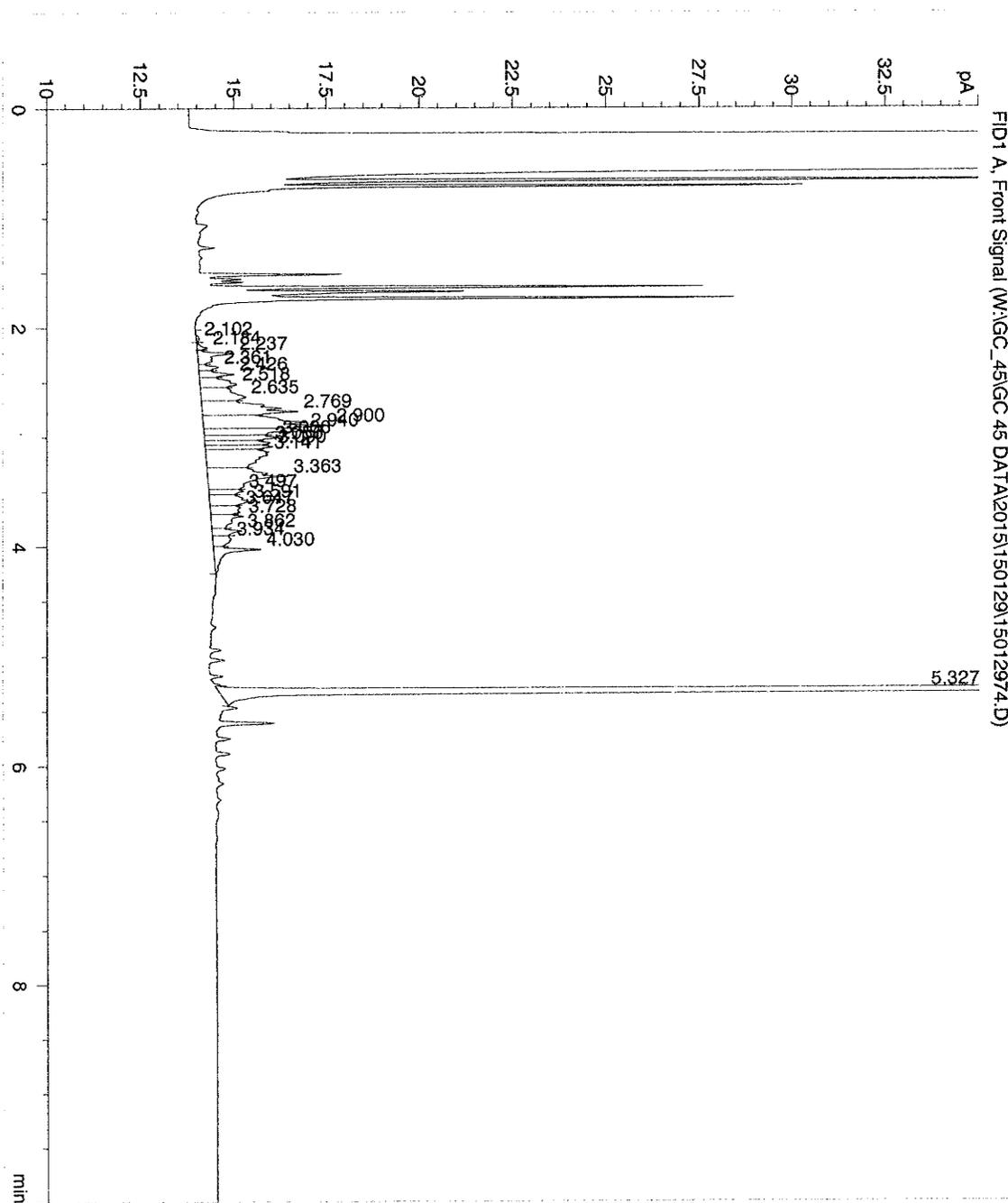
| Pk | Ret Time | Area | Height | Peak | Width | Response % |
|----|----------|--------|--------|-------|--------|------------|
| 1 | 2.102 | 0.19 | 0 BV | 0.046 | 0.028 | |
| 2 | 2.184 | 0.54 | 0 VV | 0.027 | 0.078 | |
| 3 | 2.237 | 3.13 | 1 VV | 0.043 | 0.455 | |
| 4 | 2.361 | 1.20 | 1 VV | 0.033 | 0.174 | |
| 5 | 2.426 | 2.25 | 1 VV | 0.034 | 0.326 | |
| 6 | 2.518 | 4.19 | 1 VV | 0.055 | 0.609 | |
| 7 | 2.635 | 7.01 | 1 VV | 0.079 | 1.018 | |
| 8 | 2.769 | 13.33 | 3 VV | 0.065 | 1.937 | |
| 9 | 2.900 | 17.13 | 3 VV | 0.063 | 2.488 | |
| 10 | 2.940 | 7.71 | 3 VV | 0.038 | 1.120 | |
| 11 | 3.006 | 4.98 | 2 VV | 0.036 | 0.724 | |
| 12 | 3.056 | 3.97 | 2 VV | 0.032 | 0.577 | |
| 13 | 3.090 | 3.99 | 2 VV | 0.030 | 0.580 | |
| 14 | 3.141 | 14.00 | 2 VV | 0.101 | 2.034 | |
| 15 | 3.363 | 15.19 | 2 VV | 0.086 | 2.207 | |
| 16 | 3.497 | 2.40 | 1 VV | 0.036 | 0.349 | |
| 17 | 3.591 | 5.12 | 1 VV | 0.066 | 0.744 | |
| 18 | 3.647 | 3.56 | 1 VV | 0.055 | 0.517 | |
| 19 | 3.728 | 4.11 | 1 VV | 0.059 | 0.598 | |
| 20 | 3.862 | 2.21 | 1 VV | 0.038 | 0.322 | |
| 21 | 3.934 | 2.36 | 0 VV | 0.059 | 0.343 | |
| 22 | 4.030 | 3.98 | 1 VB | 0.042 | 0.578 | |
| 23 | 5.327 | 565.69 | 292 VV | 0.032 | 82.193 | |

Total area = 688.24

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2015\150129\15012974.D
Page Number : 2
Operator : Vial Number : Vial 74
Instrument : GC 45 Injection Number : 1
Sample Name : 15-01-1715-5 Sequence Line : 74
Instrument Method: C:\CHEM32\1\METHODS\ ->
Acquired on : 30 Jan 15 11:24 am
Report Created on: 02 Feb 15 10:24 am Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies



**RAW DATA SHEET
FOR METHOD: EPA 8015B (M)**

WORK ORDER: 099-15-516
INSTRUMENT: GC 45
EXTRACTION: EPA 3510C
D/T EXTRACTED: 2015-01-29 00:00

ANALYZED BY: 682
D/T ANALYZED: 2015-01-30 05:03
REVIEWED BY:
D/T REVIEWED:

DATA FILE: W:\GC 45\GC 45 DATA\2015\150129\15012952.D\15012952

MB **CLIENT SAMPLE NUMBER: Method Blank**

LCS/MB BATCH: 150129B23A **SAMPLE VOLUME / WEIGHT:** DEFAULT: 500.00 ml / ACTUAL: 500.00 ml
MS/MSD BATCH: **FINAL VOLUME / WEIGHT:** DEFAULT: 5.00 ml / ACTUAL: 2.50 ml
UNITS: ug/L **ADJUSTMENT RATIO TO PF:** 0.50

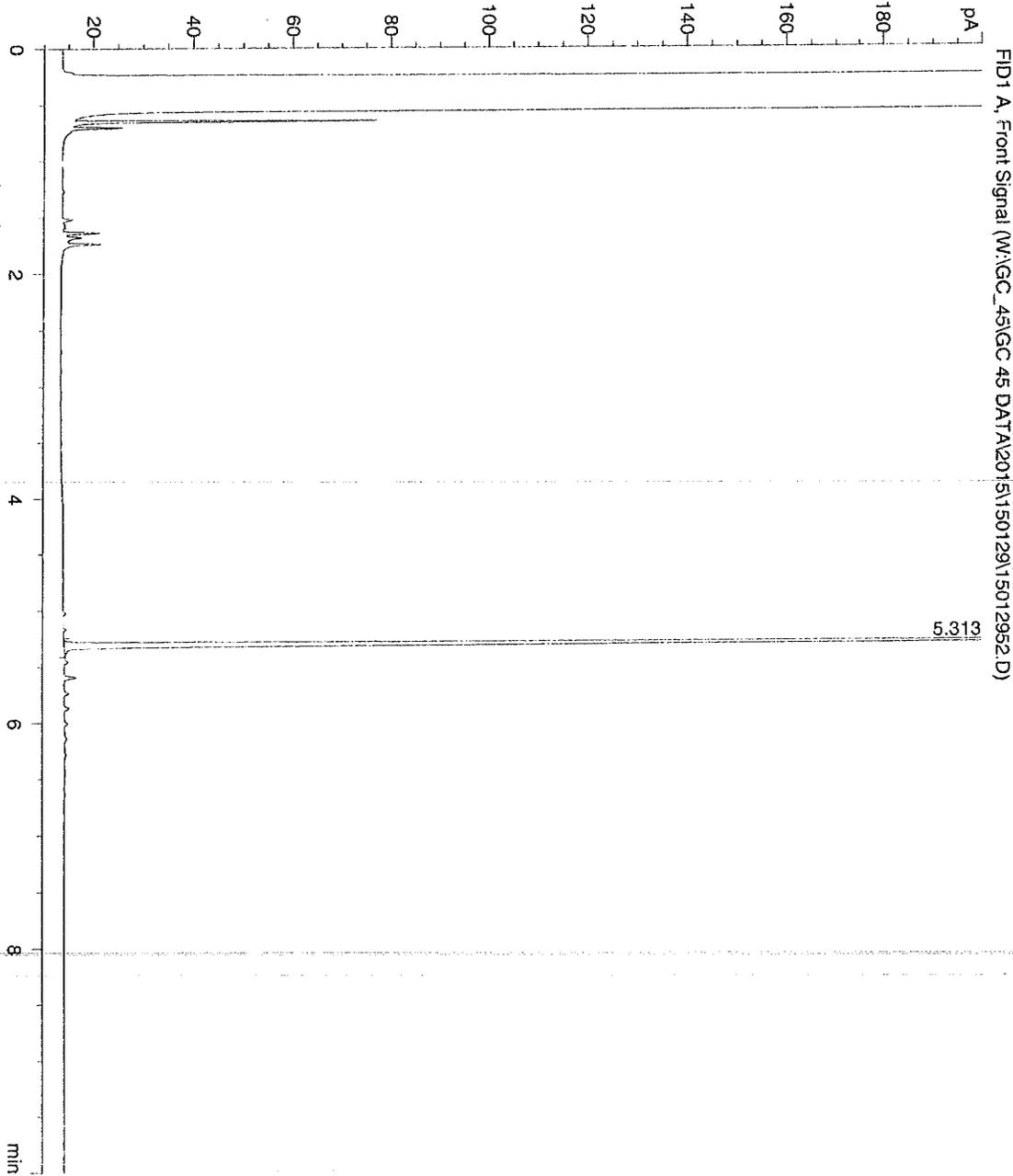
COMMENT:

| <u>COMPOUND</u> | <u>INI. CONC</u> | <u>DF</u> | <u>CONC</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>QUAL</u> |
|-----------------|------------------|-----------|-------------|-----------|------------|------------|-------------|
| TPH as Diesel | 0.000 | 1.00 | ND | 11 | 12 | 25 | |

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2015\150129\15012952.D
Page Number : 2
Operator :
Instrument : GC 45
Sample Name : MB 15012923/24
Vial Number : Vial 52
Injection Number : 1
Sequence Line : 52
Instrument Method: C:\CHEM32\1\METHODS\
Acquired on : 30 Jan 15 05:03 am
Report Created on: 30 Jan 15 07:35 pm
Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies



CONTINUING CALIBRATION VERIFICATION QUALITY CONTROL SHEET

FOR METHOD: EPA 8015B (M)

CCV WORK ORDER: 099-15-515-361-6279

BATCH ID:
INITIAL: 150119I007
CCV: 150129A127
INSTRUMENT: GC 45

ANALYZED BY: 682

D/T ANALYZED:
INITIAL: 2015-01-19 18:09
CCV: 2015-01-30 04:28
REVIEWED BY:
D/T REVIEWED:

DATA FILE: W:\GC_45\GC 45 DATA\2015\150129\15012950.D\15012950

| <u>COMPOUND NAME</u> | <u>COMP TYPE</u> | <u>CALIB MODEL</u> | <u>MIN RF</u> | <u>AVG RE</u> | <u>CCV RF</u> | <u>AMOUNT</u> | <u>CCV CONC</u> | <u>CCV %D</u> | <u>CCV %D CL</u> | <u>STATUS</u> |
|----------------------|------------------|--------------------|---------------|---------------|---------------|---------------|-----------------|---------------|------------------|---------------|
| TPH as Diesel | C | Avg Resp | 0.00 | 0.019 | 0.021 | | | -11 | 0-15 | PASS |

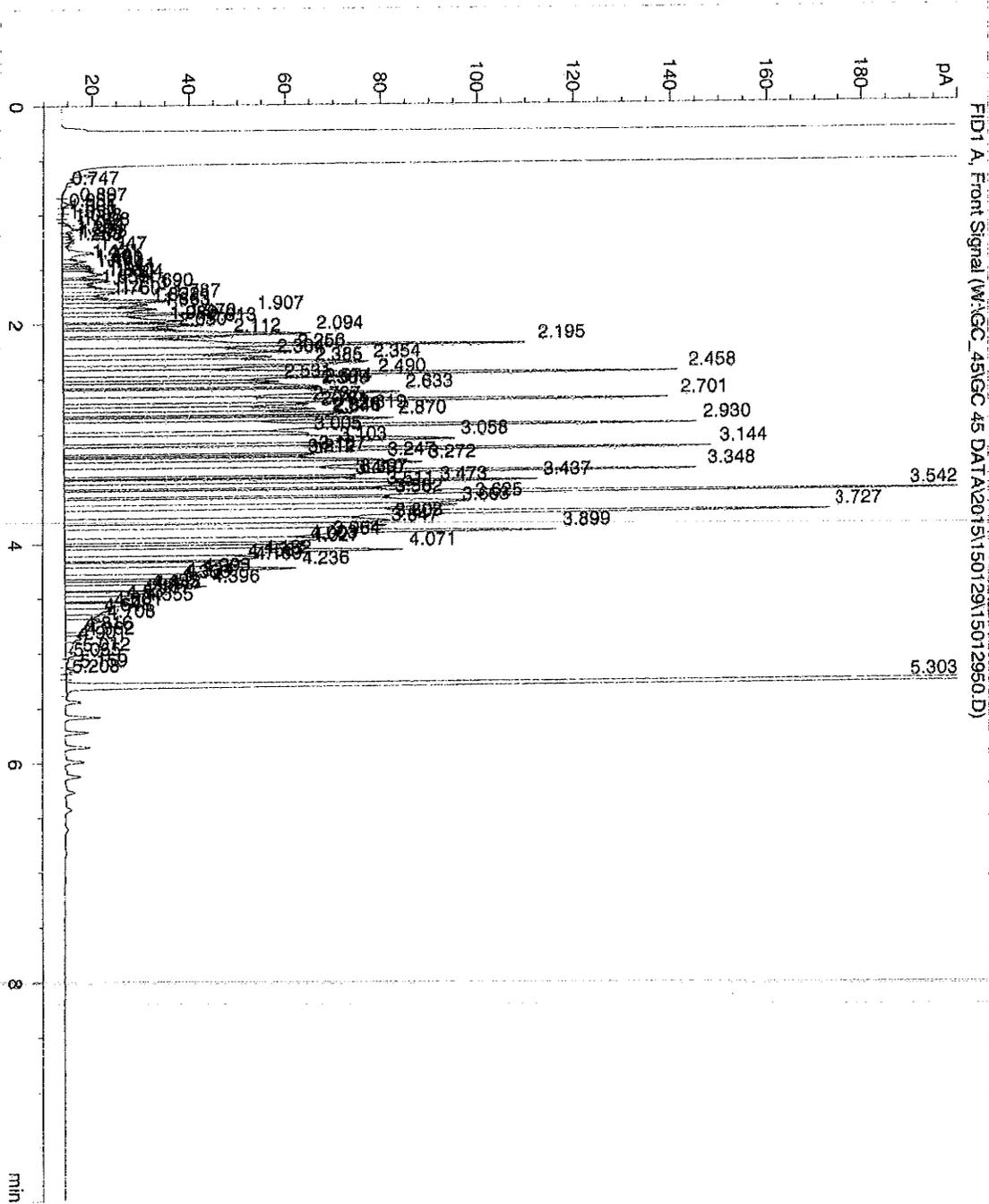
MIN RF: Method Specified Minimum Response Factor

Area Percent Report

```

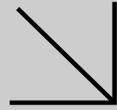
Data File Name   : W:\GC_45\GC 45 DATA\2015\150129\15012950.D
Page Number      : 4
Operator         :
Instrument       : GC 45
Sample Name      : D400 C28 50 L012815D
Vial Number     : Vial 50
Injection Number : 1
Sequence Line   : 50
Instrument Method : C:\CHEM32\1\METHODS\
Report Created on : 30 Jan 15 04:28 am
Analysis Method  : 8015B.MTH
    
```

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies





Calscience



WORK ORDER NUMBER: 15-01-1810

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For

Client: Environmental Science International, Inc.

Client Project Name: Red Hill LTM 112066

Attention: Jeff Hattemer
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Approved for release on 02/05/2015 by:
Terri Chang
Project Manager

ResultLink ▶

Email your PM ▶



Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.

Contents

Client Project Name: Red Hill LTM 112066
 Work Order Number: 15-01-1810

| | | |
|---|--|----|
| 1 | Work Order Narrative. | 3 |
| 2 | Client Sample Data. | 4 |
| | 2.1 EPA 8015B (M) TPH Diesel (Aqueous). | 4 |
| | 2.2 EPA 6020 ICP/MS Metals (Aqueous). | 5 |
| | 2.3 EPA 8270C SIM PAHs (Aqueous). | 6 |
| | 2.4 GC/MS GRO/EPA 8260B Volatile Organics (Aqueous). | 10 |
| 3 | Quality Control Sample Data. | 20 |
| | 3.1 MS/MSD. | 20 |
| | 3.2 PDS/PDSD. | 25 |
| | 3.3 LCS/LCSD. | 26 |
| 4 | Sample Analysis Summary. | 31 |
| 5 | Glossary of Terms and Qualifiers. | 32 |
| 6 | Chain-of-Custody/Sample Receipt Form. | 33 |

Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 01/29/15. They were assigned to Work Order 15-01-1810.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of ≤ 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

New York NELAP air certification does not certify for all reported methods and analytes, reference the accredited items here: http://www.calscience.com/PDF/New_York.pdf

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.



Calscience

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/29/15
Work Order: 15-01-1810
Preparation: EPA 3510C
Method: EPA 8015B (M)
Units: ug/L

Project: Red Hill LTM 112066

Page 1 of 1

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES126 | 15-01-1810-2-K | 01/28/15 10:35 | Aqueous | GC 45 | 02/02/15 | 02/03/15 03:03 | 150202B09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
- TPH as Diesel is quantified in the carbon range C10-C28.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|---------------|--------|-----|-----|-----|------|------------|
| TPH as Diesel | 1100 | 2.9 | 10 | 25 | 1.00 | HD |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|--------------|----------|----------------|------------|
| n-Octacosane | 70 | 51-141 | |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES127 | 15-01-1810-3-I | 01/28/15 11:30 | Aqueous | GC 45 | 02/02/15 | 02/03/15 03:24 | 150202B09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
- TPH as Diesel is quantified in the carbon range C10-C28.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|---------------|--------|-----|-----|-----|------|------------|
| TPH as Diesel | 1700 | 2.9 | 10 | 25 | 1.00 | HD |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|--------------|----------|----------------|------------|
| n-Octacosane | 93 | 51-141 | |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES123 | 15-01-1810-4-I | 01/28/15 09:30 | Aqueous | GC 45 | 02/02/15 | 02/03/15 03:44 | 150202B09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.
- TPH as Diesel is quantified in the carbon range C10-C28.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|---------------|--------|-----|-----|-----|------|------------|
| TPH as Diesel | 39 | 2.9 | 10 | 25 | 1.00 | HD |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|--------------|----------|----------------|------------|
| n-Octacosane | 64 | 51-141 | |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-15-516-238 | N/A | Aqueous | GC 45 | 02/02/15 | 02/03/15 00:43 | 150202B09 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|---------------|--------|-----|-----|-----|------|------------|
| TPH as Diesel | <10 | 2.9 | 10 | 25 | 1.00 | U |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|--------------|----------|----------------|------------|
| n-Octacosane | 83 | 51-141 | |



Calscience

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/29/15
Work Order: 15-01-1810
Preparation: EPA 3005A Filt.
Method: EPA 6020
Units: ug/L

Project: Red Hill LTM 112066

Page 1 of 1

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES126 | 15-01-1810-2-R | 01/28/15 10:35 | Aqueous | ICP/MS 04 | 02/03/15 | 02/03/15 20:52 | 150203LA3D |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|-----------|--------|--------|-------|-------|------|------------|
| Lead | <0.200 | 0.0898 | 0.200 | 0.500 | 1.00 | U |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES127 | 15-01-1810-3-J | 01/28/15 11:30 | Aqueous | ICP/MS 04 | 02/03/15 | 02/03/15 20:54 | 150203LA3D |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|-----------|--------|--------|-------|-------|------|------------|
| Lead | <0.200 | 0.0898 | 0.200 | 0.500 | 1.00 | U |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES123 | 15-01-1810-4-G | 01/28/15 09:30 | Aqueous | ICP/MS 04 | 02/03/15 | 02/03/15 20:56 | 150203LA3D |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|-----------|--------|--------|-------|-------|------|------------|
| Lead | <0.200 | 0.0898 | 0.200 | 0.500 | 1.00 | U |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| Method Blank | 099-14-497-118 | N/A | Aqueous | ICP/MS 04 | 02/03/15 | 02/03/15 20:40 | 150203LA3D |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|-----------|--------|--------|-------|-------|------|------------|
| Lead | <0.200 | 0.0898 | 0.200 | 0.500 | 1.00 | U |

Return to Contents

Analytical Report

Environmental Science International, Inc.
 354 Uluniu Street, Suite 304
 Kailua, HI 96734-2500

Date Received: 01/29/15
 Work Order: 15-01-1810
 Preparation: EPA 3510C
 Method: EPA 8270C SIM PAHs
 Units: ug/L

Project: Red Hill LTM 112066

Page 1 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES126 | 15-01-1810-2-N | 01/28/15 10:35 | Aqueous | GC/MS AAA | 01/30/15 | 02/02/15 22:48 | 150130L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|---------------------------|--------|-------|-------|------|------|------------|
| 2-Methylnaphthalene | 7.6 | 0.046 | 0.050 | 0.20 | 1.00 | |
| Acenaphthylene | <0.050 | 0.044 | 0.050 | 0.20 | 1.00 | U |
| Acenaphthene | 0.59 | 0.027 | 0.050 | 0.20 | 1.00 | |
| Fluorene | 0.30 | 0.042 | 0.050 | 0.20 | 1.00 | |
| Phenanthrene | <0.050 | 0.027 | 0.050 | 0.20 | 1.00 | U |
| Anthracene | <0.050 | 0.029 | 0.050 | 0.20 | 1.00 | U |
| Fluoranthene | <0.050 | 0.046 | 0.050 | 0.20 | 1.00 | U |
| Pyrene | <0.050 | 0.020 | 0.050 | 0.20 | 1.00 | U |
| Benzo (a) Anthracene | <0.050 | 0.032 | 0.050 | 0.20 | 1.00 | U |
| Chrysene | <0.050 | 0.025 | 0.050 | 0.20 | 1.00 | U |
| Benzo (k) Fluoranthene | <0.050 | 0.031 | 0.050 | 0.20 | 1.00 | U |
| Benzo (b) Fluoranthene | <0.050 | 0.017 | 0.050 | 0.20 | 1.00 | U |
| Benzo (a) Pyrene | <0.050 | 0.022 | 0.050 | 0.20 | 1.00 | U |
| Indeno (1,2,3-c,d) Pyrene | <0.050 | 0.021 | 0.050 | 0.20 | 1.00 | U |
| Dibenz (a,h) Anthracene | <0.050 | 0.047 | 0.050 | 0.20 | 1.00 | U |
| Benzo (g,h,i) Perylene | <0.099 | 0.081 | 0.099 | 0.20 | 1.00 | U |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------|----------|----------------|------------|
| Nitrobenzene-d5 | 90 | 28-139 | |
| 2-Fluorobiphenyl | 68 | 33-144 | |
| p-Terphenyl-d14 | 66 | 23-160 | |

| ES126 | 15-01-1810-2-N | 01/28/15 10:35 | Aqueous | GC/MS AAA | 01/30/15 | 02/03/15 16:18 | 150130L19 |
|-------|----------------|-------------------|---------|-----------|----------|-------------------|-----------|
|-------|----------------|-------------------|---------|-----------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|---------------------|--------|------|------|-----|------|------------|
| Naphthalene | 90 | 0.34 | 0.50 | 2.0 | 10.0 | |
| 1-Methylnaphthalene | 34 | 0.51 | 0.99 | 2.0 | 10.0 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------|----------|----------------|------------|
| Nitrobenzene-d5 | 0 | 28-139 | 1,2,6 |
| 2-Fluorobiphenyl | 64 | 33-144 | |
| p-Terphenyl-d14 | 69 | 23-160 | |

Analytical Report

Environmental Science International, Inc.
 354 Uluniu Street, Suite 304
 Kailua, HI 96734-2500

Date Received: 01/29/15
 Work Order: 15-01-1810
 Preparation: EPA 3510C
 Method: EPA 8270C SIM PAHs
 Units: ug/L

Project: Red Hill LTM 112066

Page 2 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES127 | 15-01-1810-3-H | 01/28/15 11:30 | Aqueous | GC/MS AAA | 01/30/15 | 02/02/15 23:08 | 150130L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|---------------------------|--------|-------|-------|------|------|------------|
| 2-Methylnaphthalene | 2.7 | 0.046 | 0.049 | 0.20 | 1.00 | |
| Acenaphthylene | <0.049 | 0.044 | 0.049 | 0.20 | 1.00 | U |
| Acenaphthene | 0.55 | 0.026 | 0.049 | 0.20 | 1.00 | |
| Fluorene | 0.22 | 0.042 | 0.049 | 0.20 | 1.00 | |
| Phenanthrene | <0.049 | 0.027 | 0.049 | 0.20 | 1.00 | U |
| Anthracene | <0.049 | 0.028 | 0.049 | 0.20 | 1.00 | U |
| Fluoranthene | <0.049 | 0.046 | 0.049 | 0.20 | 1.00 | U |
| Pyrene | <0.049 | 0.020 | 0.049 | 0.20 | 1.00 | U |
| Benzo (a) Anthracene | <0.049 | 0.032 | 0.049 | 0.20 | 1.00 | U |
| Chrysene | <0.049 | 0.024 | 0.049 | 0.20 | 1.00 | U |
| Benzo (k) Fluoranthene | <0.049 | 0.030 | 0.049 | 0.20 | 1.00 | U |
| Benzo (b) Fluoranthene | <0.049 | 0.017 | 0.049 | 0.20 | 1.00 | U |
| Benzo (a) Pyrene | <0.049 | 0.022 | 0.049 | 0.20 | 1.00 | U |
| Indeno (1,2,3-c,d) Pyrene | <0.049 | 0.021 | 0.049 | 0.20 | 1.00 | U |
| Dibenz (a,h) Anthracene | <0.049 | 0.047 | 0.049 | 0.20 | 1.00 | U |
| Benzo (g,h,i) Perylene | <0.098 | 0.080 | 0.098 | 0.20 | 1.00 | U |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------|----------|----------------|------------|
| Nitrobenzene-d5 | 61 | 28-139 | |
| 2-Fluorobiphenyl | 58 | 33-144 | |
| p-Terphenyl-d14 | 61 | 23-160 | |

| ES127 | 15-01-1810-3-H | 01/28/15 11:30 | Aqueous | GC/MS AAA | 01/30/15 | 02/03/15 16:38 | 150130L19 |
|-------|----------------|-------------------|---------|-----------|----------|-------------------|-----------|
|-------|----------------|-------------------|---------|-----------|----------|-------------------|-----------|

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|---------------------|--------|------|------|-----|------|------------|
| Naphthalene | 63 | 0.33 | 0.49 | 2.0 | 10.0 | |
| 1-Methylnaphthalene | 25 | 0.51 | 0.98 | 2.0 | 10.0 | |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------|----------|----------------|------------|
| Nitrobenzene-d5 | 0 | 28-139 | 1,2,6 |
| 2-Fluorobiphenyl | 50 | 33-144 | |
| p-Terphenyl-d14 | 56 | 23-160 | |



Calscience

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/29/15
Work Order: 15-01-1810
Preparation: EPA 3510C
Method: EPA 8270C SIM PAHs
Units: ug/L

Project: Red Hill LTM 112066

Page 3 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES123 | 15-01-1810-4-J | 01/28/15 09:30 | Aqueous | GC/MS AAA | 01/30/15 | 02/02/15 23:28 | 150130L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|---------------------------|--------|-------|-------|------|------|------------|
| Naphthalene | <0.048 | 0.033 | 0.048 | 0.19 | 1.00 | U |
| 2-Methylnaphthalene | <0.048 | 0.045 | 0.048 | 0.19 | 1.00 | U |
| 1-Methylnaphthalene | <0.097 | 0.050 | 0.097 | 0.19 | 1.00 | U |
| Acenaphthylene | <0.048 | 0.043 | 0.048 | 0.19 | 1.00 | U |
| Acenaphthene | <0.048 | 0.026 | 0.048 | 0.19 | 1.00 | U |
| Fluorene | <0.048 | 0.041 | 0.048 | 0.19 | 1.00 | U |
| Phenanthrene | <0.048 | 0.026 | 0.048 | 0.19 | 1.00 | U |
| Anthracene | <0.048 | 0.028 | 0.048 | 0.19 | 1.00 | U |
| Fluoranthene | <0.048 | 0.045 | 0.048 | 0.19 | 1.00 | U |
| Pyrene | <0.048 | 0.020 | 0.048 | 0.19 | 1.00 | U |
| Benzo (a) Anthracene | <0.048 | 0.032 | 0.048 | 0.19 | 1.00 | U |
| Chrysene | <0.048 | 0.024 | 0.048 | 0.19 | 1.00 | U |
| Benzo (k) Fluoranthene | <0.048 | 0.030 | 0.048 | 0.19 | 1.00 | U |
| Benzo (b) Fluoranthene | <0.048 | 0.017 | 0.048 | 0.19 | 1.00 | U |
| Benzo (a) Pyrene | <0.048 | 0.022 | 0.048 | 0.19 | 1.00 | U |
| Indeno (1,2,3-c,d) Pyrene | <0.048 | 0.020 | 0.048 | 0.19 | 1.00 | U |
| Dibenz (a,h) Anthracene | <0.048 | 0.046 | 0.048 | 0.19 | 1.00 | U |
| Benzo (g,h,i) Perylene | <0.097 | 0.079 | 0.097 | 0.19 | 1.00 | U |

| Surrogate | Rec. (%) | Control Limits | Qualifiers |
|------------------|----------|----------------|------------|
| Nitrobenzene-d5 | 50 | 28-139 | |
| 2-Fluorobiphenyl | 55 | 33-144 | |
| p-Terphenyl-d14 | 67 | 23-160 | |

Return to Contents

Analytical Report

Environmental Science International, Inc.
 354 Uluniu Street, Suite 304
 Kailua, HI 96734-2500

Date Received: 01/29/15
 Work Order: 15-01-1810
 Preparation: EPA 3510C
 Method: EPA 8270C SIM PAHs
 Units: ug/L

Project: Red Hill LTM 112066

Page 4 of 4

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|----------------------|---------------------|----------------|------------------|-----------------|---------------------------|------------------|
| Method Blank | 099-15-148-80 | N/A | Aqueous | GC/MS AAA | 01/30/15 | 02/02/15 21:28 | 150130L19 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>DF</u> | <u>Qualifiers</u> |
|---------------------------|---------------|-----------|------------|------------|-----------|-------------------|
| Naphthalene | <0.050 | 0.034 | 0.050 | 0.20 | 1.00 | U |
| 2-Methylnaphthalene | <0.050 | 0.046 | 0.050 | 0.20 | 1.00 | U |
| 1-Methylnaphthalene | <0.10 | 0.052 | 0.10 | 0.20 | 1.00 | U |
| Acenaphthylene | <0.050 | 0.045 | 0.050 | 0.20 | 1.00 | U |
| Acenaphthene | <0.050 | 0.027 | 0.050 | 0.20 | 1.00 | U |
| Fluorene | <0.050 | 0.043 | 0.050 | 0.20 | 1.00 | U |
| Phenanthrene | <0.050 | 0.027 | 0.050 | 0.20 | 1.00 | U |
| Anthracene | <0.050 | 0.029 | 0.050 | 0.20 | 1.00 | U |
| Fluoranthene | <0.050 | 0.047 | 0.050 | 0.20 | 1.00 | U |
| Pyrene | <0.050 | 0.020 | 0.050 | 0.20 | 1.00 | U |
| Benzo (a) Anthracene | <0.050 | 0.033 | 0.050 | 0.20 | 1.00 | U |
| Chrysene | <0.050 | 0.025 | 0.050 | 0.20 | 1.00 | U |
| Benzo (k) Fluoranthene | <0.050 | 0.031 | 0.050 | 0.20 | 1.00 | U |
| Benzo (b) Fluoranthene | <0.050 | 0.018 | 0.050 | 0.20 | 1.00 | U |
| Benzo (a) Pyrene | <0.050 | 0.022 | 0.050 | 0.20 | 1.00 | U |
| Indeno (1,2,3-c,d) Pyrene | <0.050 | 0.021 | 0.050 | 0.20 | 1.00 | U |
| Dibenz (a,h) Anthracene | <0.050 | 0.048 | 0.050 | 0.20 | 1.00 | U |
| Benzo (g,h,i) Perylene | <0.10 | 0.082 | 0.10 | 0.20 | 1.00 | U |

| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> |
|------------------|-----------------|-----------------------|-------------------|
| Nitrobenzene-d5 | 69 | 28-139 | |
| 2-Fluorobiphenyl | 67 | 33-144 | |
| p-Terphenyl-d14 | 69 | 23-160 | |

Analytical Report

Environmental Science International, Inc.
 354 Uluniu Street, Suite 304
 Kailua, HI 96734-2500

Date Received: 01/29/15
 Work Order: 15-01-1810
 Preparation: EPA 5030C
 Method: GC/MS / EPA 8260B
 Units: ug/L

Project: Red Hill LTM 112066

Page 1 of 10

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ESTRIP | 15-01-1810-1-A | 01/28/15 09:00 | Aqueous | GC/MS OO | 01/31/15 | 01/31/15 14:11 | 150131L001 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|-----------------------------|--------|------|------|-----|------|------------|
| Acetone | <10 | 6.0 | 10 | 20 | 1.00 | U,IJ |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1.00 | U,IH |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1.00 | U |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1.00 | U |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1.00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1.00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1.00 | U |

Analytical Report

Environmental Science International, Inc.
 354 Uluniu Street, Suite 304
 Kailua, HI 96734-2500

Date Received: 01/29/15
 Work Order: 15-01-1810
 Preparation: EPA 5030C
 Method: GC/MS / EPA 8260B
 Units: ug/L

Project: Red Hill LTM 112066

Page 2 of 10

| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|-----------------|-----------------------|-------------------|------------|-----------|-------------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | <30 | 26 | 30 | 50 | 1.00 | U |
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> | | | |
| Dibromofluoromethane | 104 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 105 | 80-134 | | | | |
| Toluene-d8 | 97 | 80-120 | | | | |
| Toluene-d8-TPPH | 98 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 90 | 80-120 | | | | |

Analytical Report

Environmental Science International, Inc.
 354 Uluniu Street, Suite 304
 Kailua, HI 96734-2500

Date Received: 01/29/15
 Work Order: 15-01-1810
 Preparation: EPA 5030C
 Method: GC/MS / EPA 8260B
 Units: ug/L

Project: Red Hill LTM 112066

Page 3 of 10

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES126 | 15-01-1810-2-A | 01/28/15 10:35 | Aqueous | GC/MS OO | 01/31/15 | 01/31/15 14:38 | 150131L001 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|-----------------------------|--------|------|------|-----|------|------------|
| Acetone | <10 | 6.0 | 10 | 20 | 1.00 | U,IJ |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1.00 | U,IH |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1.00 | U |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1.00 | U |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| Ethylbenzene | 0.16 | 0.14 | 0.50 | 1.0 | 1.00 | J |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1.00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1.00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1.00 | U |

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/29/15
Work Order: 15-01-1810
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B
Units: ug/L

Project: Red Hill LTM 112066

Page 4 of 10

| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|-----------------|-----------------------|-------------------|------------|-----------|-------------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | 0.35 | 0.23 | 0.50 | 1.0 | 1.00 | J |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | 54 | 26 | 30 | 50 | 1.00 | |
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> | | | |
| Dibromofluoromethane | 104 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 107 | 80-134 | | | | |
| Toluene-d8 | 97 | 80-120 | | | | |
| Toluene-d8-TPPH | 98 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 93 | 80-120 | | | | |

Analytical Report

Environmental Science International, Inc.
 354 Uluniu Street, Suite 304
 Kailua, HI 96734-2500

Date Received: 01/29/15
 Work Order: 15-01-1810
 Preparation: EPA 5030C
 Method: GC/MS / EPA 8260B
 Units: ug/L

Project: Red Hill LTM 112066

Page 5 of 10

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES127 | 15-01-1810-3-A | 01/28/15 11:30 | Aqueous | GC/MS OO | 01/31/15 | 01/31/15 16:52 | 150131L001 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|-----------------------------|--------|------|------|-----|------|------------|
| Acetone | <10 | 6.0 | 10 | 20 | 1.00 | U,IJ |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1.00 | U,IH |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1.00 | U |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1.00 | U |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| Ethylbenzene | 0.17 | 0.14 | 0.50 | 1.0 | 1.00 | J |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1.00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1.00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1.00 | U |



Calscience

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/29/15
Work Order: 15-01-1810
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B
Units: ug/L

Project: Red Hill LTM 112066

Page 6 of 10

| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|-----------------|-----------------------|-------------------|------------|-----------|-------------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | 0.35 | 0.23 | 0.50 | 1.0 | 1.00 | J |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | 59 | 26 | 30 | 50 | 1.00 | |
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> | | | |
| Dibromofluoromethane | 102 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 106 | 80-134 | | | | |
| Toluene-d8 | 96 | 80-120 | | | | |
| Toluene-d8-TPPH | 97 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 94 | 80-120 | | | | |


 Return to Contents



Calscience

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/29/15
Work Order: 15-01-1810
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B
Units: ug/L

Project: Red Hill LTM 112066

Page 7 of 10

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|-------------------|---------------------|---------|------------|---------------|--------------------|-------------|
| ES123 | 15-01-1810-4-A | 01/28/15 09:30 | Aqueous | GC/MS OO | 01/31/15 | 01/31/15 17:19 | 150131L001 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| Parameter | Result | DL | LOD | LOQ | DF | Qualifiers |
|-----------------------------|--------|------|------|-----|------|------------|
| Acetone | <10 | 6.0 | 10 | 20 | 1.00 | U,IJ |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1.00 | U,IH |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1.00 | U |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1.00 | U |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1.00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1.00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1.00 | U |

Return to Contents

Analytical Report

Environmental Science International, Inc.
 354 Uluniu Street, Suite 304
 Kailua, HI 96734-2500

Date Received: 01/29/15
 Work Order: 15-01-1810
 Preparation: EPA 5030C
 Method: GC/MS / EPA 8260B
 Units: ug/L

Project: Red Hill LTM 112066

Page 8 of 10

| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|-----------------|-----------------------|-------------------|------------|-----------|-------------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | <30 | 26 | 30 | 50 | 1.00 | U |
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> | | | |
| Dibromofluoromethane | 100 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 103 | 80-134 | | | | |
| Toluene-d8 | 96 | 80-120 | | | | |
| Toluene-d8-TPPH | 97 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 93 | 80-120 | | | | |



Calscience

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/29/15
Work Order: 15-01-1810
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B
Units: ug/L

Project: Red Hill LTM 112066

Page 9 of 10

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|----------------------|---------------------|----------------|-----------------|-----------------|---------------------------|-------------------|
| Method Blank | 099-13-057-76 | N/A | Aqueous | GC/MS OO | 01/31/15 | 01/31/15 13:10 | 150131L001 |

Comment(s): - Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag.

| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|-----------|------------|------------|-----------|-------------------|
| Acetone | <10 | 6.0 | 10 | 20 | 1.00 | U |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1.00 | U |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1.00 | U |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1.00 | U |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1.00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1.00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1.00 | U |

Return to Contents



Calscience

Analytical Report

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/29/15
Work Order: 15-01-1810
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B
Units: ug/L

Project: Red Hill LTM 112066

Page 10 of 10

| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|-----------------|-----------------------|-------------------|------------|-----------|-------------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | <30 | 26 | 30 | 50 | 1.00 | U |
| <u>Surrogate</u> | <u>Rec. (%)</u> | <u>Control Limits</u> | <u>Qualifiers</u> | | | |
| Dibromofluoromethane | 102 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 103 | 80-134 | | | | |
| Toluene-d8 | 97 | 80-120 | | | | |
| Toluene-d8-TPPH | 99 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 92 | 80-120 | | | | |


 Return to Contents



Calscience

Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/29/15
Work Order: 15-01-1810
Preparation: EPA 3510C
Method: EPA 8015B (M)

Project: Red Hill LTM 112066

Page 1 of 5

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|--------------------|-----------------|-----------------|------------------|---------------------|-----------------|------------|---------------|-------------------|
| ES126 | Sample | Aqueous | GC 45 | 02/02/15 | 02/03/15 03:03 | 150202S09 | | | | |
| ES126 | Matrix Spike | Aqueous | GC 45 | 02/02/15 | 02/03/15 02:23 | 150202S09 | | | | |
| ES126 | Matrix Spike Duplicate | Aqueous | GC 45 | 02/02/15 | 02/03/15 02:43 | 150202S09 | | | | |
| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>MS Conc.</u> | <u>MS %Rec.</u> | <u>MSD Conc.</u> | <u>MSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| TPH as Diesel | 1101 | 2000 | 2835 | 87 | 2958 | 93 | 55-133 | 4 | 0-30 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

| | | |
|---|----------------|-----------------|
| Environmental Science International, Inc. | Date Received: | 01/29/15 |
| 354 Uluniu Street, Suite 304 | Work Order: | 15-01-1810 |
| Kailua, HI 96734-2500 | Preparation: | EPA 3005A Filt. |
| | Method: | EPA 6020 |
| Project: Red Hill LTM 112066 | | Page 2 of 5 |

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| ES126 | Sample | Aqueous | ICP/MS 04 | 02/03/15 | 02/03/15 20:52 | 150203SA3 |
| ES126 | Matrix Spike | Aqueous | ICP/MS 04 | 02/03/15 | 02/03/15 20:44 | 150203SA3 |
| ES126 | Matrix Spike Duplicate | Aqueous | ICP/MS 04 | 02/03/15 | 02/03/15 20:46 | 150203SA3 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Lead | ND | 100.0 | 99.87 | 100 | 103.6 | 104 | 80-120 | 4 | 0-20 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/29/15
Work Order: 15-01-1810
Preparation: EPA 3510C
Method: EPA 8270C SIM PAHS

Project: Red Hill LTM 112066

Page 3 of 5

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number | | | | |
|---------------------------|------------------------|-------------|------------|---------------|----------------|---------------------|----------|-----|--------|------------|
| ES126 | Sample | Aqueous | GC/MS AAA | 01/30/15 | 02/03/15 16:18 | 150130S19A | | | | |
| ES126 | Matrix Spike | Aqueous | GC/MS AAA | 01/30/15 | 02/02/15 20:07 | 150130S19A | | | | |
| ES126 | Matrix Spike Duplicate | Aqueous | GC/MS AAA | 01/30/15 | 02/02/15 20:28 | 150130S19A | | | | |
| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Naphthalene | 89.96 | 2.000 | 83.41 | 0 | 72.55 | 0 | 21-133 | 14 | 0-25 | 3 |
| 2-Methylnaphthalene | 7.590 | 2.000 | 4.948 | 0 | 5.156 | 0 | 21-140 | 4 | 0-25 | 3 |
| 1-Methylnaphthalene | 34.15 | 2.000 | 28.85 | 0 | 25.83 | 0 | 20-140 | 11 | 0-25 | 3 |
| Acenaphthylene | ND | 2.000 | 1.563 | 78 | 1.498 | 75 | 33-145 | 4 | 0-25 | |
| Acenaphthene | 0.5935 | 2.000 | 2.018 | 71 | 1.843 | 62 | 49-121 | 9 | 0-25 | |
| Fluorene | 0.2969 | 2.000 | 1.704 | 70 | 1.568 | 64 | 59-121 | 8 | 0-25 | |
| Phenanthrene | ND | 2.000 | 1.611 | 81 | 1.556 | 78 | 54-120 | 3 | 0-25 | |
| Anthracene | ND | 2.000 | 1.320 | 66 | 1.240 | 62 | 27-133 | 6 | 0-25 | |
| Fluoranthene | ND | 2.000 | 1.451 | 73 | 1.374 | 69 | 26-137 | 5 | 0-25 | |
| Pyrene | ND | 2.000 | 1.468 | 73 | 1.444 | 72 | 18-168 | 2 | 0-25 | |
| Benzo (a) Anthracene | ND | 2.000 | 1.409 | 70 | 1.328 | 66 | 33-143 | 6 | 0-25 | |
| Chrysene | ND | 2.000 | 1.471 | 74 | 1.401 | 70 | 17-168 | 5 | 0-25 | |
| Benzo (k) Fluoranthene | ND | 2.000 | 1.556 | 78 | 1.356 | 68 | 24-159 | 14 | 0-25 | |
| Benzo (b) Fluoranthene | ND | 2.000 | 1.200 | 60 | 1.070 | 54 | 24-159 | 11 | 0-25 | |
| Benzo (a) Pyrene | ND | 2.000 | 1.293 | 65 | 1.207 | 60 | 17-163 | 7 | 0-25 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 2.000 | 1.330 | 66 | 1.261 | 63 | 10-171 | 5 | 0-25 | |
| Dibenz (a,h) Anthracene | ND | 2.000 | 1.381 | 69 | 1.271 | 64 | 10-219 | 8 | 0-25 | |
| Benzo (g,h,i) Perylene | ND | 2.000 | 1.264 | 63 | 1.348 | 67 | 10-227 | 6 | 0-25 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/29/15
Work Order: 15-01-1810
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B

Project: Red Hill LTM 112066

Page 4 of 5

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Batch Number |
|---------------------------|------------------------|---------|------------|---------------|----------------|---------------------|
| ES126 | Sample | Aqueous | GC/MS OO | 01/31/15 | 01/31/15 14:38 | 150131S001 |
| ES126 | Matrix Spike | Aqueous | GC/MS OO | 01/31/15 | 01/31/15 15:05 | 150131S001 |
| ES126 | Matrix Spike Duplicate | Aqueous | GC/MS OO | 01/31/15 | 01/31/15 15:31 | 150131S001 |

| Parameter | Sample Conc. | Spike Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------------|--------------|-------------|----------|----------|-----------|-----------|----------|-----|--------|------------|
| Acetone | ND | 50.00 | 65.98 | 132 | 63.42 | 127 | 40-140 | 4 | 0-20 | |
| Benzene | ND | 50.00 | 43.74 | 87 | 43.62 | 87 | 80-120 | 0 | 0-20 | |
| Bromodichloromethane | ND | 50.00 | 44.53 | 89 | 44.71 | 89 | 75-120 | 0 | 0-20 | |
| Bromoform | ND | 50.00 | 43.49 | 87 | 44.63 | 89 | 70-130 | 3 | 0-20 | |
| Bromomethane | ND | 50.00 | 59.65 | 119 | 53.85 | 108 | 30-145 | 10 | 0-20 | |
| 2-Butanone | ND | 50.00 | 53.01 | 106 | 51.20 | 102 | 30-150 | 3 | 0-20 | |
| Carbon Tetrachloride | ND | 50.00 | 40.25 | 80 | 40.78 | 82 | 65-140 | 1 | 0-20 | |
| Chlorobenzene | ND | 50.00 | 46.77 | 94 | 46.50 | 93 | 80-120 | 1 | 0-20 | |
| Chloroethane | ND | 50.00 | 39.98 | 80 | 41.50 | 83 | 60-135 | 4 | 0-20 | |
| Chloroform | ND | 50.00 | 45.37 | 91 | 45.16 | 90 | 65-135 | 0 | 0-20 | |
| Chloromethane | ND | 50.00 | 37.73 | 75 | 39.18 | 78 | 40-125 | 4 | 0-20 | |
| Dibromochloromethane | ND | 50.00 | 48.10 | 96 | 47.60 | 95 | 60-135 | 1 | 0-20 | |
| 1,2-Dibromo-3-Chloropropane | ND | 50.00 | 40.53 | 81 | 39.34 | 79 | 50-130 | 3 | 0-20 | |
| 1,2-Dibromoethane | ND | 50.00 | 49.12 | 98 | 48.98 | 98 | 80-120 | 0 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 47.05 | 94 | 46.53 | 93 | 70-120 | 1 | 0-20 | |
| 1,3-Dichlorobenzene | ND | 50.00 | 46.08 | 92 | 45.62 | 91 | 75-125 | 1 | 0-20 | |
| 1,4-Dichlorobenzene | ND | 50.00 | 46.22 | 92 | 45.85 | 92 | 75-125 | 1 | 0-20 | |
| 1,1-Dichloroethane | ND | 50.00 | 42.14 | 84 | 42.19 | 84 | 70-135 | 0 | 0-20 | |
| 1,2-Dichloroethane | ND | 50.00 | 50.52 | 101 | 49.45 | 99 | 70-130 | 2 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 39.86 | 80 | 40.36 | 81 | 70-130 | 1 | 0-20 | |
| c-1,2-Dichloroethene | ND | 50.00 | 44.04 | 88 | 43.78 | 88 | 70-125 | 1 | 0-20 | |
| t-1,2-Dichloroethene | ND | 50.00 | 39.54 | 79 | 39.99 | 80 | 60-140 | 1 | 0-20 | |
| 1,2-Dichloropropane | ND | 50.00 | 43.81 | 88 | 43.85 | 88 | 75-125 | 0 | 0-20 | |
| c-1,3-Dichloropropene | ND | 50.00 | 43.26 | 87 | 43.00 | 86 | 70-130 | 1 | 0-20 | |
| t-1,3-Dichloropropene | ND | 50.00 | 40.98 | 82 | 41.61 | 83 | 55-140 | 2 | 0-20 | |
| Ethylbenzene | ND | 50.00 | 45.15 | 90 | 44.90 | 90 | 75-125 | 1 | 0-20 | |
| Methylene Chloride | ND | 50.00 | 41.05 | 82 | 41.07 | 82 | 55-140 | 0 | 0-20 | |
| 4-Methyl-2-Pentanone | ND | 50.00 | 43.39 | 87 | 42.42 | 85 | 60-135 | 2 | 0-20 | |
| Styrene | ND | 50.00 | 44.39 | 89 | 43.61 | 87 | 65-135 | 2 | 0-20 | |
| 1,1,1,2-Tetrachloroethane | ND | 50.00 | 44.95 | 90 | 45.50 | 91 | 80-130 | 1 | 0-20 | |
| 1,1,2,2-Tetrachloroethane | ND | 50.00 | 41.23 | 82 | 41.67 | 83 | 65-130 | 1 | 0-20 | |
| Tetrachloroethene | ND | 50.00 | 55.26 | 111 | 54.04 | 108 | 45-150 | 2 | 0-20 | |
| Toluene | ND | 50.00 | 44.30 | 89 | 43.48 | 87 | 75-120 | 2 | 0-20 | |
| 1,2,4-Trichlorobenzene | ND | 50.00 | 47.52 | 95 | 46.31 | 93 | 65-135 | 3 | 0-20 | |
| 1,1,1-Trichloroethane | ND | 50.00 | 42.47 | 85 | 42.76 | 86 | 65-130 | 1 | 0-20 | |

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/29/15
Work Order: 15-01-1810
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B

Project: Red Hill LTM 112066

Page 5 of 5

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>MS Conc.</u> | <u>MS %Rec.</u> | <u>MSD Conc.</u> | <u>MSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------|---------------------|--------------------|-----------------|-----------------|------------------|------------------|-----------------|------------|---------------|-------------------|
| Hexachloro-1,3-Butadiene | ND | 50.00 | 46.89 | 94 | 46.45 | 93 | 50-140 | 1 | 0-20 | |
| 1,1,2-Trichloroethane | ND | 50.00 | 46.89 | 94 | 47.41 | 95 | 75-125 | 1 | 0-20 | |
| Trichloroethene | ND | 50.00 | 44.46 | 89 | 43.96 | 88 | 70-125 | 1 | 0-20 | |
| 1,2,3-Trichloropropane | ND | 50.00 | 38.07 | 76 | 37.07 | 74 | 75-125 | 3 | 0-20 | 3 |
| Vinyl Chloride | ND | 50.00 | 40.78 | 82 | 41.96 | 84 | 50-145 | 3 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 89.32 | 89 | 88.22 | 88 | 75-130 | 1 | 0-20 | |
| o-Xylene | ND | 50.00 | 44.60 | 89 | 44.47 | 89 | 80-120 | 0 | 0-20 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 41.69 | 83 | 42.71 | 85 | 65-125 | 2 | 0-20 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - PDS

Environmental Science International, Inc.
 354 Uluniu Street, Suite 304
 Kailua, HI 96734-2500

Date Received: 01/29/15
 Work Order: 15-01-1810
 Preparation: EPA 3005A Filt.
 Method: EPA 6020

Project: Red Hill LTM 112066

Page 1 of 1

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | PDS/PDSD Batch Number |
|---------------------------|--------|---------|------------|----------------|----------------|-----------------------|
| ES126 | Sample | Aqueous | ICP/MS 04 | 02/03/15 00:00 | 02/03/15 20:52 | 150203SA3 |
| ES126 | PDS | Aqueous | ICP/MS 04 | 02/03/15 00:00 | 02/03/15 20:48 | 150203SA3 |

| <u>Parameter</u> | <u>Sample Conc.</u> | <u>Spike Added</u> | <u>PDS Conc.</u> | <u>PDS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
|------------------|---------------------|--------------------|------------------|------------------|-----------------|-------------------|
| Lead | ND | 100.0 | 98.15 | 98 | 75-125 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/29/15
Work Order: 15-01-1810
Preparation: EPA 3510C
Method: EPA 8015B (M)

Project: Red Hill LTM 112066

Page 1 of 5

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | | |
|---------------------------|-------------|-----------|------------|---------------|----------------|-----------------------|-----|--------|------------|
| 099-15-516-238 | LCS | Aqueous | GC 45 | 02/02/15 | 02/03/15 01:03 | 150202B09 | | | |
| 099-15-516-238 | LCSD | Aqueous | GC 45 | 02/02/15 | 02/03/15 01:23 | 150202B09 | | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| TPH as Diesel | 2000 | 1897 | 95 | 1893 | 95 | 60-132 | 0 | 0-11 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits

Quality Control - LCS

Environmental Science International, Inc.
 354 Uluniu Street, Suite 304
 Kailua, HI 96734-2500

Date Received: 01/29/15
 Work Order: 15-01-1810
 Preparation: EPA 3005A Filt.
 Method: EPA 6020

Project: Red Hill LTM 112066

Page 2 of 5

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|--------------------|------------------------|------------------|-----------------------|-------------------|
| 099-14-497-118 | LCS | Aqueous | ICP/MS 04 | 02/03/15 | 02/03/15 20:42 | 150203LA3D |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Lead | | 100.0 | 96.03 | 96 | 80-120 | |



Calscience

Quality Control - LCS

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/29/15
Work Order: 15-01-1810
Preparation: EPA 3510C
Method: EPA 8270C SIM PAHs

Project: Red Hill LTM 112066

Page 3 of 5

| Quality Control Sample ID | Type | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number |
|---------------------------|------------|--------------------|------------------------|------------------|-----------------------|-------------------|
| 099-15-148-80 | LCS | Aqueous | GC/MS AAA | 01/30/15 | 02/03/15 03:08 | 150130L19 |
| <u>Parameter</u> | | <u>Spike Added</u> | <u>Conc. Recovered</u> | <u>LCS %Rec.</u> | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Naphthalene | | 2.000 | 1.510 | 75 | 21-133 | |
| 2-Methylnaphthalene | | 2.000 | 1.537 | 77 | 21-140 | |
| 1-Methylnaphthalene | | 2.000 | 1.528 | 76 | 20-140 | |
| Acenaphthylene | | 2.000 | 1.458 | 73 | 33-145 | |
| Acenaphthene | | 2.000 | 1.547 | 77 | 55-121 | |
| Fluorene | | 2.000 | 1.454 | 73 | 59-121 | |
| Phenanthrene | | 2.000 | 1.536 | 77 | 54-120 | |
| Anthracene | | 2.000 | 1.262 | 63 | 27-133 | |
| Fluoranthene | | 2.000 | 1.595 | 80 | 26-137 | |
| Pyrene | | 2.000 | 1.784 | 89 | 45-129 | |
| Benzo (a) Anthracene | | 2.000 | 1.585 | 79 | 33-143 | |
| Chrysene | | 2.000 | 1.722 | 86 | 17-168 | |
| Benzo (k) Fluoranthene | | 2.000 | 1.970 | 98 | 24-159 | |
| Benzo (b) Fluoranthene | | 2.000 | 1.389 | 69 | 24-159 | |
| Benzo (a) Pyrene | | 2.000 | 1.473 | 74 | 17-163 | |
| Indeno (1,2,3-c,d) Pyrene | | 2.000 | 1.508 | 75 | 25-175 | |
| Dibenz (a,h) Anthracene | | 2.000 | 1.577 | 79 | 25-175 | |
| Benzo (g,h,i) Perylene | | 2.000 | 1.351 | 68 | 25-157 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/29/15
Work Order: 15-01-1810
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B

Project: Red Hill LTM 112066

Page 4 of 5

| Quality Control Sample ID | Type | Matrix | | Instrument | Date Prepared | Date Analyzed | LCS/LCSD Batch Number | | |
|-----------------------------|-------------|-----------|-----------|------------|---------------|----------------|-----------------------|--------|------------|
| 099-13-057-76 | LCS | Aqueous | | GC/MS OO | 01/31/15 | 01/31/15 11:51 | 150131L001 | | |
| 099-13-057-76 | LCSD | Aqueous | | GC/MS OO | 01/31/15 | 01/31/15 12:11 | 150131L001 | | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Acetone | 50.00 | 82.13 | 164 | N/A | N/A | 40-140 | N/A | 0-20 | X |
| Benzene | 50.00 | 47.93 | 96 | N/A | N/A | 80-120 | N/A | 0-20 | |
| Bromodichloromethane | 50.00 | 48.08 | 96 | N/A | N/A | 75-120 | N/A | 0-20 | |
| Bromoform | 50.00 | 46.64 | 93 | N/A | N/A | 70-130 | N/A | 0-20 | |
| Bromomethane | 50.00 | 53.01 | 106 | N/A | N/A | 30-145 | N/A | 0-20 | |
| 2-Butanone | 50.00 | 62.76 | 126 | N/A | N/A | 30-150 | N/A | 0-20 | |
| Carbon Tetrachloride | 50.00 | 45.54 | 91 | N/A | N/A | 65-140 | N/A | 0-20 | |
| Chlorobenzene | 50.00 | 51.78 | 104 | N/A | N/A | 80-120 | N/A | 0-20 | |
| Chloroethane | 50.00 | 44.69 | 89 | N/A | N/A | 60-135 | N/A | 0-20 | |
| Chloroform | 50.00 | 48.16 | 96 | N/A | N/A | 65-135 | N/A | 0-20 | |
| Chloromethane | 50.00 | 42.87 | 86 | N/A | N/A | 40-125 | N/A | 0-20 | |
| Dibromochloromethane | 50.00 | 51.24 | 102 | N/A | N/A | 60-135 | N/A | 0-20 | |
| 1,2-Dibromo-3-Chloropropane | 50.00 | 38.87 | 78 | N/A | N/A | 50-130 | N/A | 0-20 | |
| 1,2-Dibromoethane | 50.00 | 50.90 | 102 | N/A | N/A | 80-120 | N/A | 0-20 | |
| 1,2-Dichlorobenzene | 50.00 | 51.42 | 103 | N/A | N/A | 70-120 | N/A | 0-20 | |
| 1,3-Dichlorobenzene | 50.00 | 51.65 | 103 | N/A | N/A | 75-125 | N/A | 0-20 | |
| 1,4-Dichlorobenzene | 50.00 | 51.48 | 103 | N/A | N/A | 75-125 | N/A | 0-20 | |
| 1,1-Dichloroethane | 50.00 | 44.96 | 90 | N/A | N/A | 70-135 | N/A | 0-20 | |
| 1,2-Dichloroethane | 50.00 | 53.05 | 106 | N/A | N/A | 70-130 | N/A | 0-20 | |
| 1,1-Dichloroethene | 50.00 | 44.07 | 88 | N/A | N/A | 70-130 | N/A | 0-20 | |
| c-1,2-Dichloroethene | 50.00 | 47.25 | 95 | N/A | N/A | 70-125 | N/A | 0-20 | |
| t-1,2-Dichloroethene | 50.00 | 43.09 | 86 | N/A | N/A | 60-140 | N/A | 0-20 | |
| 1,2-Dichloropropane | 50.00 | 46.41 | 93 | N/A | N/A | 75-125 | N/A | 0-20 | |
| c-1,3-Dichloropropene | 50.00 | 46.07 | 92 | N/A | N/A | 70-130 | N/A | 0-20 | |
| t-1,3-Dichloropropene | 50.00 | 45.03 | 90 | N/A | N/A | 55-140 | N/A | 0-20 | |
| Ethylbenzene | 50.00 | 49.99 | 100 | N/A | N/A | 75-125 | N/A | 0-20 | |
| Methylene Chloride | 50.00 | 44.21 | 88 | N/A | N/A | 55-140 | N/A | 0-20 | |
| 4-Methyl-2-Pentanone | 50.00 | 43.41 | 87 | N/A | N/A | 60-135 | N/A | 0-20 | |
| Styrene | 50.00 | 48.69 | 97 | N/A | N/A | 65-135 | N/A | 0-20 | |
| 1,1,1,2-Tetrachloroethane | 50.00 | 48.98 | 98 | N/A | N/A | 80-130 | N/A | 0-20 | |
| 1,1,2,2-Tetrachloroethane | 50.00 | 41.18 | 82 | N/A | N/A | 65-130 | N/A | 0-20 | |
| Tetrachloroethene | 50.00 | 61.86 | 124 | N/A | N/A | 45-150 | N/A | 0-20 | |
| Toluene | 50.00 | 48.40 | 97 | N/A | N/A | 75-120 | N/A | 0-20 | |
| 1,2,4-Trichlorobenzene | 50.00 | 50.89 | 102 | N/A | N/A | 65-135 | N/A | 0-20 | |
| 1,1,1-Trichloroethane | 50.00 | 46.72 | 93 | N/A | N/A | 65-130 | N/A | 0-20 | |
| Hexachloro-1,3-Butadiene | 50.00 | 52.10 | 104 | N/A | N/A | 50-140 | N/A | 0-20 | |

RPD: Relative Percent Difference. CL: Control Limits



Calscience

Quality Control - LCS/LCSD

Environmental Science International, Inc.
354 Uluniu Street, Suite 304
Kailua, HI 96734-2500

Date Received: 01/29/15
Work Order: 15-01-1810
Preparation: EPA 5030C
Method: GC/MS / EPA 8260B

Project: Red Hill LTM 112066

Page 5 of 5

| <u>Parameter</u> | <u>Spike Added</u> | <u>LCS Conc.</u> | <u>LCS %Rec.</u> | <u>LCSD Conc.</u> | <u>LCSD %Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
|-----------------------------|--------------------|------------------|------------------|-------------------|-------------------|-----------------|------------|---------------|-------------------|
| 1,1,2-Trichloroethane | 50.00 | 48.92 | 98 | N/A | N/A | 75-125 | N/A | 0-20 | |
| Trichloroethene | 50.00 | 49.55 | 99 | N/A | N/A | 70-125 | N/A | 0-20 | |
| 1,2,3-Trichloropropane | 50.00 | 39.93 | 80 | N/A | N/A | 75-125 | N/A | 0-20 | |
| Vinyl Chloride | 50.00 | 46.46 | 93 | N/A | N/A | 50-145 | N/A | 0-20 | |
| p/m-Xylene | 100.0 | 99.55 | 100 | N/A | N/A | 75-130 | N/A | 0-20 | |
| o-Xylene | 50.00 | 49.14 | 98 | N/A | N/A | 80-120 | N/A | 0-20 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 41.61 | 83 | N/A | N/A | 65-125 | N/A | 0-20 | |
| Gasoline Range Organics | 1000 | 1097 | 110 | 1105 | 111 | 80-120 | 1 | 0-20 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits

Sample Analysis Summary Report

Work Order: 15-01-1810

Page 1 of 1

| <u>Method</u> | <u>Extraction</u> | <u>Chemist ID</u> | <u>Instrument</u> | <u>Analytical Location</u> |
|--------------------|-------------------|-------------------|-------------------|----------------------------|
| EPA 6020 | EPA 3005A Filtr. | 598 | ICP/MS 04 | 1 |
| EPA 8015B (M) | EPA 3510C | 960 | GC 45 | 1 |
| EPA 8270C SIM PAHs | EPA 3510C | 907 | GC/MS AAA | 1 |
| GC/MS / EPA 8260B | EPA 5030C | 849 | GC/MS OO | 2 |

| <u>Qualifiers</u> | <u>Definition</u> |
|-------------------|---|
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| B | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| DL | The Detection Limit (DL) is the smallest analyte concentration that can be demonstrated to be different from zero or a blank concentration at the 99% level of confidence. |
| E | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| ICH | Initial calibration verification recovery is above the control limit for this analyte. |
| ICJ | Initial calibration verification recovery is below the control limit for this analyte. |
| IH | Calibration verification recovery is above the control limit for this analyte. |
| IJ | Calibration verification recovery is below the control limit for this analyte. |
| J | Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| LOD | The Limit of Detection (LOD) is the smallest amount or concentration of a substance that must be present in a sample in order to be detected at 99% confidence level. |
| LOQ | The Limit of Quantitation (LOQ) is the lowest concentration of a substance that produces a quantitative result within specified limits of precision and bias. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| SG | The sample extract was subjected to Silica Gel treatment prior to analysis. |
| U | Undetected at Detection Limit (DL) and is reported as less than the Limit of Detection (LOD). |
| X | % Recovery and/or RPD out-of-range. |
| Z | Analyte presence was not confirmed by second column or GC/MS analysis. |
| | Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis. |
| | Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time. |
| | A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations. |

ORIGIN ID:HNLA

SHIP DATE: 28JAN15
ACTWGT: 52.4 LB
CAD: /POS1525
DIMS: 24x15x13 IN
BILL RECIPIENT

UNITED STATES US

TO **SMPLE CONTROL
CALSCIENCE LABORTORIES
7440 LINCOLN WAY**

GARDEN GROVE CA 92841

(714) 896-5494

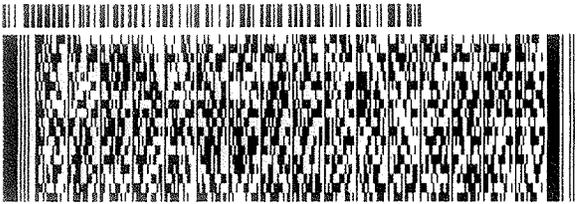
REF:

INV:

DEPT:

658056 01/28 54711/RF1E/FE48

1510



**FedEx
Express**



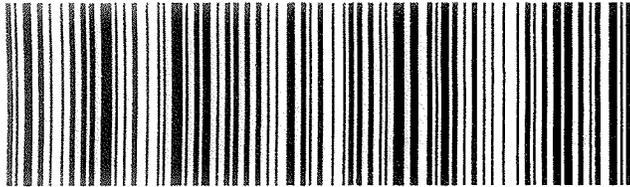
J151015011401uv

1 of 2
TRK# 8059 2709 6814
0200
MASTER

**THU - 29 JAN AA
STANDARD OVERNIGHT**

WZ APVA

92841
CA-US SNA



ORIGIN ID:HNLA

SHIP DATE: 28JAN15
ACTWGT: 51.3 LB
CAD: /POS1525
DIMS: 24x15x13 IN
BILL RECIPIENT

11/20/2014 09:46:00 # 1004
#50055 01/29 53711/RF15/FF4R

UNITED STATES US

TO SMPLE CONTROL
CALSCIENCE LABORTORIES
7440 LINCOLN WAY

GARDEN GROVE CA 92841

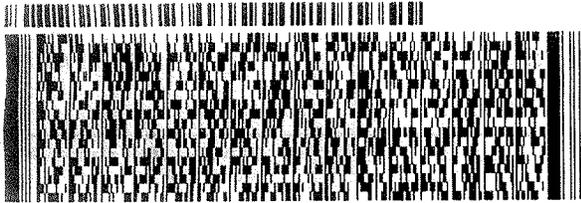
(714) 895-6494

REF:

DEPT:

INV:

PO:



FedEx
Express



J1510150114010V

2 of 2

MPS# 7801 7093 8910
0681

Mstr# 8059 2709 6814

0200

THU - 29 JAN AA
STANDARD OVERNIGHT

WZ APVA

92841
CA-US SNA



SAMPLE RECEIPT FORM

Cooler 1 of 1

CLIENT: ESI

DATE: 01/29/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 2.7 °C + 0.2 °C (CF) = 2.9 °C Blank Sample

Sample(s) outside temperature criteria (PM/APM contacted by: _____)

Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: Air Filter Checked by: LS

CUSTODY SEALS INTACT:

Cooler _____ No (Not Intact) Not Present N/A Checked by: LS

Sample _____ No (Not Intact) Not Present Checked by: 917

| SAMPLE CONDITION: | Yes | No | N/A |
|---|-------------------------------------|--------------------------|-------------------------------------|
| Chain-Of-Custody (COC) document(s) received with samples..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| COC document(s) received complete..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input type="checkbox"/> Collection date/time, matrix, and/or # of containers logged in based on sample labels. | | | |
| <input type="checkbox"/> No analysis requested. <input type="checkbox"/> Not relinquished. <input type="checkbox"/> No date/time relinquished. | | | |
| Sampler's name indicated on COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container label(s) consistent with COC..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Sample container(s) intact and good condition..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Proper containers and sufficient volume for analyses requested..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Analyses received within holding time..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Aqueous samples received within 15-minute holding time | | | |
| <input type="checkbox"/> pH <input type="checkbox"/> Residual Chlorine <input type="checkbox"/> Dissolved Sulfides <input type="checkbox"/> Dissolved Oxygen..... | | | |
| Proper preservation noted on COC or sample container..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| <input checked="" type="checkbox"/> Unpreserved vials received for Volatiles analysis | | | |
| Volatile analysis container(s) free of headspace..... | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| Tedlar bag(s) free of condensation..... | <input type="checkbox"/> | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

CONTAINER TYPE:

Solid: 4ozCGJ 8ozCGJ 16ozCGJ Sleeve (____) EnCores® TerraCores® _____

Aqueous: VOA VOA⁽⁻¹⁾ VOAna₂ 125AGB 125AGBh 125AGBp 1AGB 1AGBna₂ 1AGBs

500AGB 500AGJ 500AGJs 250AGB 250CGB 250CGBs 1PB 1PBna 500PB

250PB 250PBn 125PB 125PBz_{na} 100PJ 100PJna₂ _____ _____ _____

Air: Tedlar® Canister **Other:** _____ **Trip Blank Lot#:** 140900B **Labeled/Checked by:** 917/→

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope **Reviewed by:** 681/681

Preservative: h: HCL n: HNO₃ na₂: Na₂S₂O₃ na: NaOH p: H₃PO₄ s: H₂SO₄ u: Ultra-pure z_{na}: ZnAc₂+NaOH f: Filtered **Scanned by:** 681

Return to Contents



Calscience

WORK ORDER #: 15-01-1810

SAMPLE RECEIPT FORM

Cooler 1 of 1

CLIENT: EST

DATE: 01/30/15

TEMPERATURE: Thermometer ID: SC4 (Criteria: 0.0 °C – 6.0 °C, not frozen except sediment/tissue)

Temperature 2.9 °C + 0.2 °C (CF) = 3.1 °C [X] Blank [] Sample

[] Sample(s) outside temperature criteria (PM/APM contacted by: _____)

[] Sample(s) outside temperature criteria but received on ice/chilled on same day of sampling.

[] Received at ambient temperature, placed on ice for transport by Courier.

Ambient Temperature: [] Air [] Filter

Checked by: 15

CUSTODY SEALS INTACT:

[X] Cooler [] _____ [] No (Not Intact) [] Not Present [] N/A Checked by: 15

[X] Sample [] _____ [] No (Not Intact) [X] Not Present Checked by: 977

SAMPLE CONDITION:

Table with 4 columns: Description, Yes, No, N/A. Rows include Chain-Of-Custody (COC) document(s) received with samples, COC document(s) received complete, Sampler's name indicated on COC, Sample container label(s) consistent with COC, Sample container(s) intact and good condition, Proper containers and sufficient volume for analyses requested, Analyses received within holding time, Aqueous samples received within 15-minute holding time, Proper preservation noted on COC or sample container, Volatile analysis container(s) free of headspace, Tedlar bag(s) free of condensation.

CONTAINER TYPE:

Solid: [] 4ozCGJ [] 8ozCGJ [] 16ozCGJ [] Sleeve (____) [] EnCores® [] TerraCores® [] _____

Aqueous: [] VOA [] VOA_h [] VOA_{na2} [] 125AGB [] 125AGB_h [] 125AGB_p [X] 1AGB [] 1AGB_{na2} [] 1AGB_s

[] 500AGB [X] 500AGJ [] 500AGJ_s [] 250AGB [] 250CGB [] 250CGB_s [] 1PB [] 1PB_{na} [] 500PB

[] 250PB [] 250PB_n [] 125PB [] 125PB_{znna} [] 100PJ [] 100PJ_{na2} [] _____ [] _____ [] _____

Air: [] Tedlar® [] Canister Other: [] _____ Trip Blank Lot#: _____ Labeled/Checked by: 977

Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope Reviewed by: 681

Preservative: h: HCL n: HNO3 na2:Na2S2O3 na: NaOH p: H3PO4 s: H2SO4 u: Ultra-pure znna: ZnAc2+NaOH f: Filtered Scanned by: 681

Return to Contents

**RAW DATA SHEET
FOR METHOD: EPA 8015B (M)**

WORK ORDER: 15-01-1810
INSTRUMENT: GC 45
EXTRACTION: EPA 3510C ✓
D/T EXTRACTED: 2015-02-02 00:00 ✓

ANALYZED BY: 960
D/T ANALYZED: 2015-02-03 03:03
REVIEWED BY:
D/T REVIEWED:

DATA FILE: W:\GC 45\GC 45 DATA\2015\150202\15020252.D\15020252 ✓

2 **CLIENT SAMPLE NUMBER: ES126**

LCS/MB BATCH: 150202B09 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 500.00 ml / ACTUAL: 500.00 ml ✓
MS/MSD BATCH: 150202S09 **FINAL VOLUME / WEIGHT:** DEFAULT: 5.00 ml / ACTUAL: 2.50 ml
UNITS: ug/L **ADJUSTMENT RATIO TO PF:** 0.50

COMMENT: Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

| <u>COMPOUND</u> | <u>INI. CONC</u> | <u>DF</u> | <u>CONC</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>QUAL</u> |
|-----------------|------------------|-----------|-------------|-----------|------------|------------|-------------|
| TPH as Diesel | 220000 | 1.00 ✓ | 1100 | 2.9 | 10 | 25 | b |

=====
Area Percent Report
=====

Data File Name : W:\GC_45\GC 45 DATA\2015\150202\15020252.D
Page Number : 1
Operator : Vial Number : Vial 52
Instrument : GC 45 Injection Number : 1
Sample Name : 15-01-1810-2 Sequence Line : 52
Instrument Method: C:\CHEM32\1\METHODS\ ->
Acquired on : 03 Feb 15 03:03 am
Report Created on: 03 Feb 15 03:36 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Sig. 1 in W:\GC_45\GC 45 DATA\2015\150202\ ->

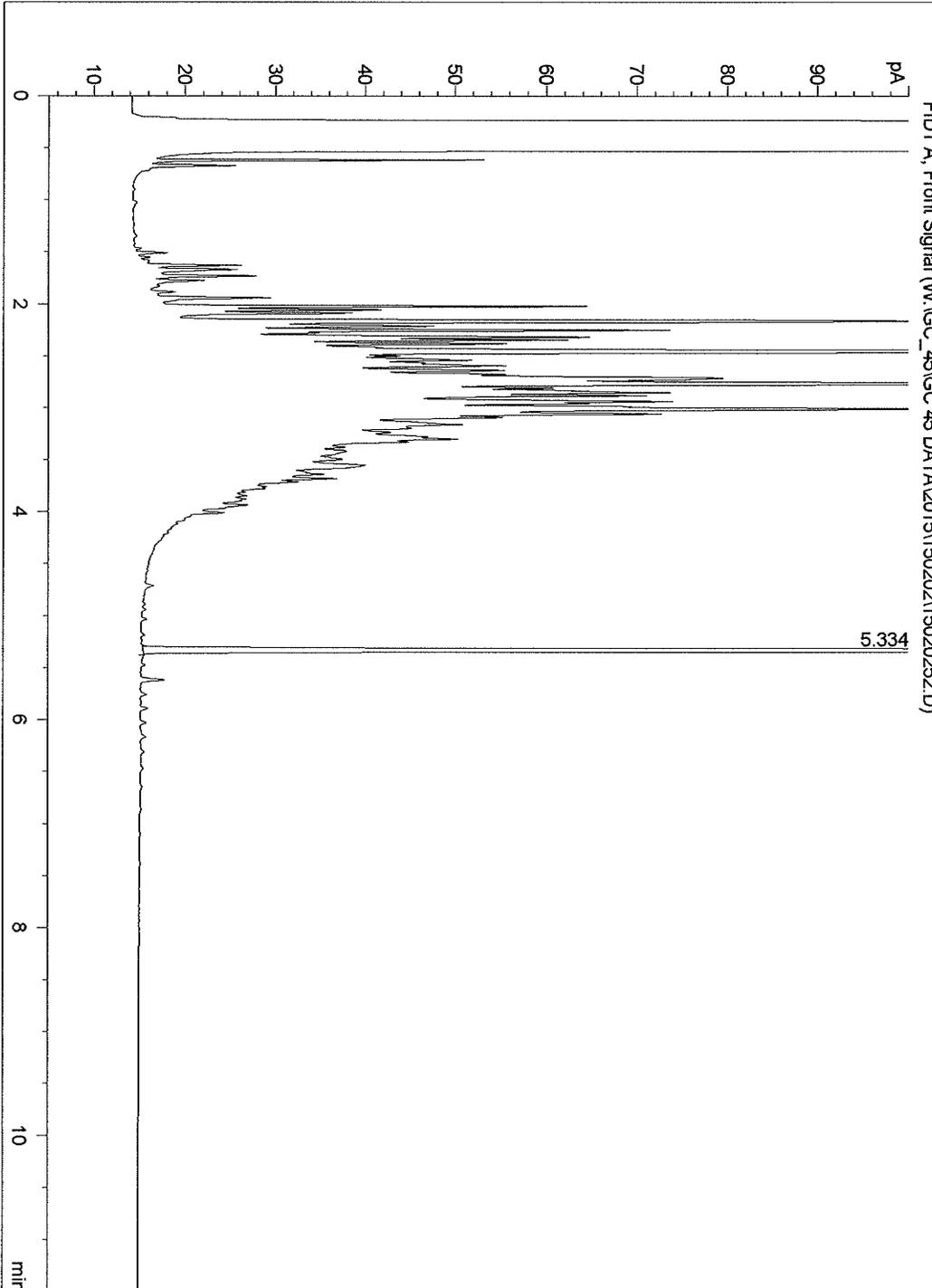
| Pk | Ret Time | Area | Height | Peak | Width | Response % |
|----|----------|--------|--------|------|-------|------------|
| 1 | 5.334 | 621.38 | 439 | BB | 0.022 | 100.000 |
| 2 | 0.000 | 0.00 | 0 | | 0.000 | 0.000 |

Total area = 621.38
0.00

=====
Area Percent Report
=====

Data File Name : W:\GC_45\GC 45 DATA\2015\150202\15020252.D
Page Number : 2
Operator : Vial Number : Vial 52
Instrument : GC 45 Injection Number : 1
Sample Name : 15-01-1810-2 Sequence Line : 52
Instrument Method: C:\CHEM32\1\METHODS\ ->
Acquired on : 03 Feb 15 03:03 am
Report Created on: 03 Feb 15 03:36 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies



=====
 Area Percent Report
 =====

Data File Name : W:\GC_45\GC 45 DATA\2015\150202\15020252.D
 Page Number : 1
 Operator : 682 Vial Number : Vial 52
 Instrument : GC 45 Injection Number : 1
 Sample Name : 15-01-1810-2 Sequence Line : 52
 Instrument Method: C:\CHEM32\1\METHODS\ ->
 Acquired on : 03 Feb 15 03:03 am
 Report Created on: 03 Feb 15 03:33 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Sig. 1 in W:\GC_45\GC 45 DATA\2015\150202\ ->

| Pk | Ret Time | Area | Height | Peak | Width | Response % |
|-----|----------|--------|--------|--------|-------|------------|
| --- | ----- | ----- | ----- | --- | ----- | ----- |
| 1 | 1.941 | 30.25 | | 15 VV | 0.029 | 0.626 |
| 2 | 2.022 | 61.89 | | 50 VV | 0.020 | 1.281 |
| 3 | 2.056 | 73.99 | | 27 VV | 0.035 | 1.531 |
| 4 | 2.164 | 169.21 | | 96 VV | 0.028 | 3.502 |
| 5 | 2.212 | 54.13 | | 33 VV | 0.024 | 1.120 |
| 6 | 2.252 | 103.98 | | 59 VV | 0.026 | 2.152 |
| 7 | 2.317 | 149.21 | | 50 VV | 0.041 | 3.088 |
| 8 | 2.381 | 67.99 | | 41 VV | 0.023 | 1.407 |
| 9 | 2.451 | 531.93 | | 328 VV | 0.024 | 11.008 |
| 10 | 2.539 | 80.44 | | 37 VV | 0.033 | 1.665 |
| 11 | 2.593 | 123.19 | | 41 VV | 0.041 | 2.549 |
| 12 | 2.638 | 84.03 | | 41 VV | 0.030 | 1.739 |
| 13 | 2.714 | 239.73 | | 65 VV | 0.053 | 4.961 |
| 14 | 2.766 | 280.11 | | 144 VV | 0.029 | 5.797 |
| 15 | 2.851 | 308.45 | | 59 VV | 0.065 | 6.383 |
| 16 | 2.939 | 173.99 | | 60 VV | 0.039 | 3.601 |
| 17 | 3.010 | 250.52 | | 102 VV | 0.035 | 5.184 |
| 18 | 3.058 | 201.49 | | 58 VV | 0.045 | 4.170 |
| 19 | 3.161 | 182.10 | | 36 VV | 0.065 | 3.768 |
| 20 | 3.238 | 54.84 | | 28 VV | 0.027 | 1.135 |
| 21 | 3.301 | 196.53 | | 36 VV | 0.068 | 4.067 |
| 22 | 3.379 | 43.41 | | 23 VV | 0.026 | 0.898 |
| 23 | 3.420 | 94.19 | | 23 VV | 0.053 | 1.949 |
| 24 | 3.497 | 70.89 | | 23 VV | 0.043 | 1.467 |
| 25 | 3.558 | 113.87 | | 25 VV | 0.067 | 2.357 |
| 26 | 3.641 | 62.78 | | 21 VV | 0.040 | 1.299 |
| 27 | 3.684 | 86.14 | | 22 VV | 0.050 | 1.783 |
| 28 | 3.764 | 65.73 | | 15 VV | 0.056 | 1.360 |
| 29 | 3.848 | 24.41 | | 12 VV | 0.027 | 0.505 |
| 30 | 3.882 | 37.99 | | 12 VV | 0.041 | 0.786 |
| 31 | 3.941 | 42.34 | | 12 VV | 0.046 | 0.876 |
| 32 | 4.013 | 117.83 | | 10 VV | 0.147 | 2.438 |
| 33 | 4.713 | 11.50 | | 2 VV | 0.079 | 0.238 |
| 34 | 4.892 | 2.55 | | 1 VV | 0.038 | 0.053 |
| 35 | 4.943 | 3.39 | | 1 VV | 0.047 | 0.070 |
| 36 | 5.036 | 5.04 | | 1 VV | 0.060 | 0.104 |
| 37 | 5.186 | 2.58 | | 1 VV | 0.040 | 0.053 |

=====
Area Percent Report
=====

Data File Name : W:\GC_45\GC 45 DATA\2015\150202\15020252.D
Page Number : 2
Operator : 682 Vial Number : Vial 52
Instrument : GC 45 Injection Number : 1
Sample Name : 15-01-1810-2 Sequence Line : 52
Instrument Method: C:\CHEM32\1\METHODS\ ->
Acquired on : 03 Feb 15 03:03 am
Report Created on: 03 Feb 15 03:33 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

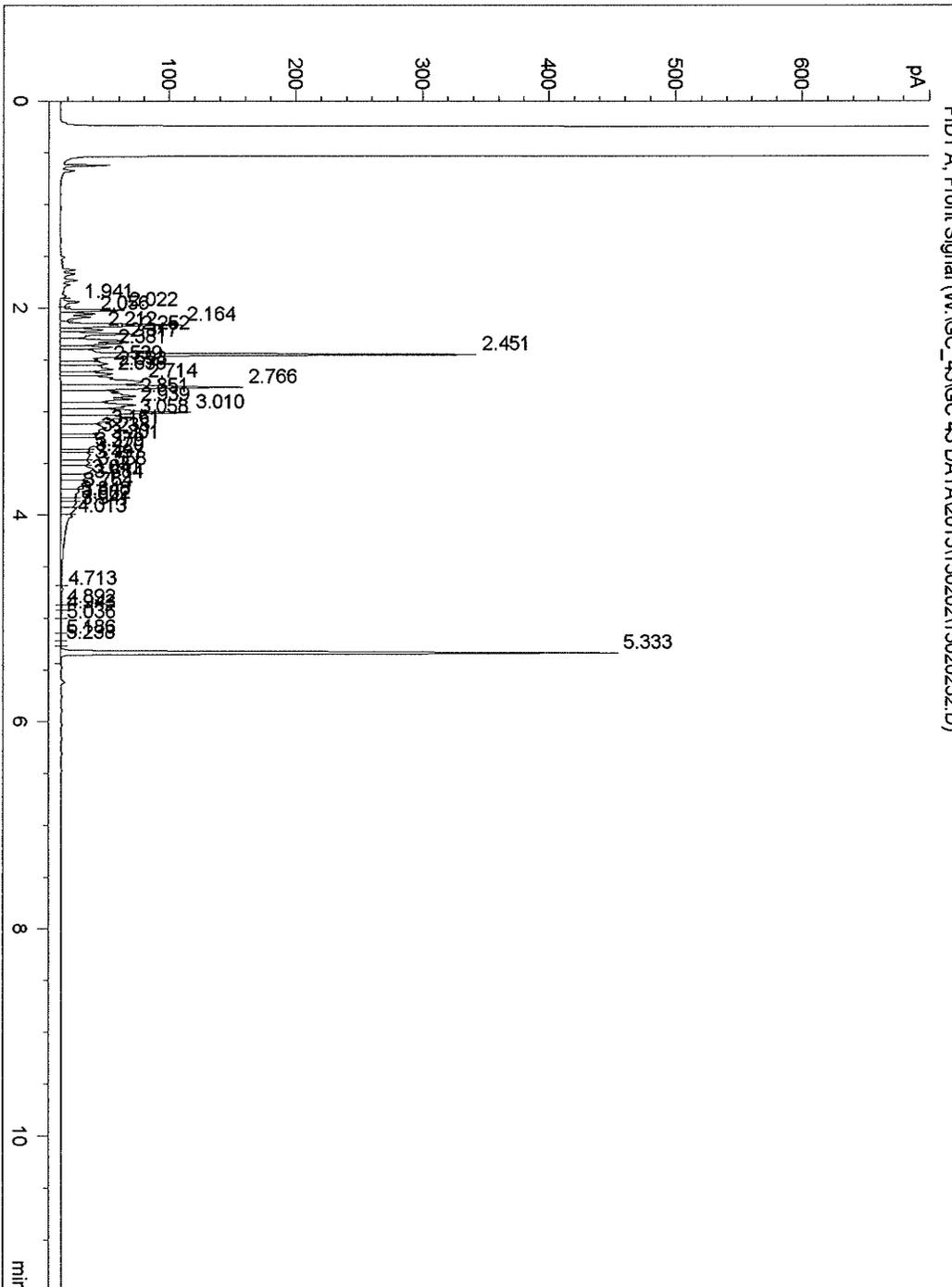
| Pk | Ret Time | Area | Height | Peak | Width | Response % |
|----|----------|--------|--------|--------|-------|------------|
| 38 | 5.238 | 1.68 | | 1 VV | 0.041 | 0.035 |
| 39 | 5.333 | 627.84 | | 440 VV | 0.024 | 12.993 |

Total area = 4832.19

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2015\150202\15020252.D
Page Number : 3
Operator : 682 Vial Number : Vial 52
Instrument : GC 45 Injection Number : 1
Sample Name : 15-01-1810-2 Sequence Line : 52
Instrument Method: C:\CHEM32\1\METHODS\ ->
Acquired on : 03 Feb 15 03:03 am
Report Created on: 03 Feb 15 03:33 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies



FID1 A, Front Signal (W:\GC_45\GC 45 DATA\2015\150202\15020252.D)

RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER: 15-01-1810
INSTRUMENT: GC 45
EXTRACTION: EPA 3510C
D/T EXTRACTED: 2015-02-02 00:00 ✓

ANALYZED BY: 960
D/T ANALYZED: 2015-02-03 03:24
REVIEWED BY:
D/T REVIEWED:

DATA FILE: W:\GC 45\GC 45 DATA\2015\150202\15020253.D\15020253 ✓

3 **CLIENT SAMPLE NUMBER: ES127**

LCS/MB BATCH: 150202B09 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 500.00 ml / ACTUAL: 500.00 ml ✓
MS/MSD BATCH: 150202S09 **FINAL VOLUME / WEIGHT:** DEFAULT: 5.00 ml / ACTUAL: 2.50 ml
UNITS: ug/L **ADJUSTMENT RATIO TO PF:** 0.50

COMMENT: Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

| <u>COMPOUND</u> | <u>INI. CONC</u> | <u>DF</u> | <u>CONC</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>QUAL</u> |
|-----------------|------------------|-----------|-------------|-----------|------------|------------|-------------|
| TPH as Diesel | 330000 | 1.00 ✓ | 1650 | 2.9 | 10 | 25 | b |

=====
Area Percent Report
=====

Data File Name : W:\GC_45\GC 45 DATA\2015\150202\15020253.D
Page Number : 1
Operator : Vial Number : Vial 53
Instrument : GC 45 Injection Number : 1
Sample Name : 15-01-1810-3 Sequence Line : 53
Instrument Method: C:\CHEM32\1\METHODS\ ->
Acquired on : 03 Feb 15 03:24 am
Report Created on: 03 Feb 15 03:36 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Sig. 1 in W:\GC_45\GC 45 DATA\2015\150202\ ->

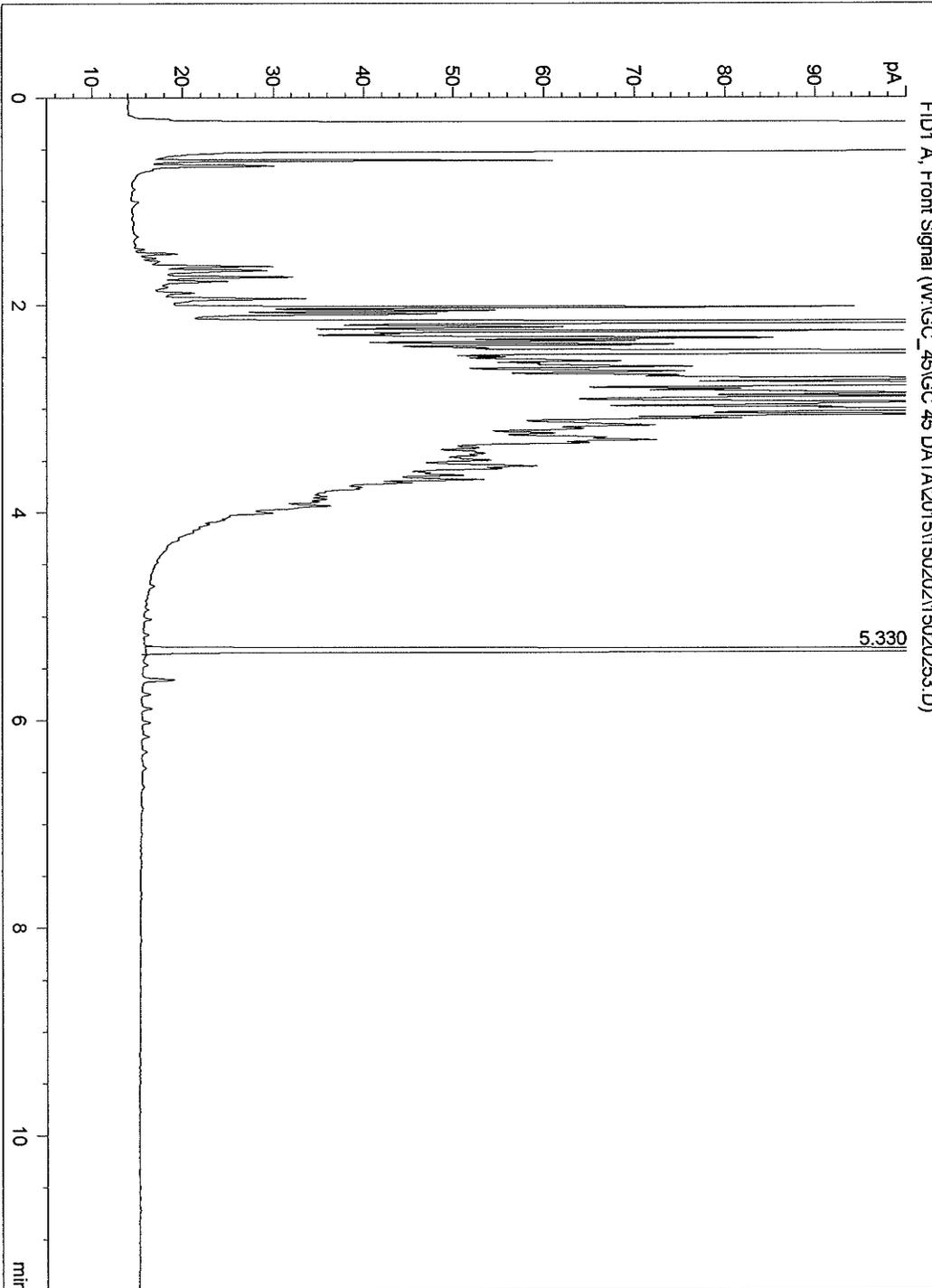
| Pk | Ret Time | Area | Height | Peak | Width | Response % |
|----|----------|--------|--------|------|-------|------------|
| 1 | 5.330 | 823.29 | 588 | BB | 0.021 | 100.000 |
| 2 | 0.000 | 0.00 | 0 | | 0.000 | 0.000 |

Total area = 823.29
0.00

=====
Area Percent Report
=====

Data File Name : W:\GC_45\GC 45 DATA\2015\150202\15020253.D
Page Number : 2
Operator : Vial Number : Vial 53
Instrument : GC 45 Injection Number : 1
Sample Name : 15-01-1810-3 Sequence Line : 53
Instrument Method: C:\CHEM32\1\METHODS\ ->
Acquired on : 03 Feb 15 03:24 am
Report Created on: 03 Feb 15 03:36 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies



=====
 Area Percent Report
 =====

Data File Name : W:\GC_45\GC 45 DATA\2015\150202\15020253.D
 Page Number : 1
 Operator : 682 Vial Number : Vial 53
 Instrument : GC 45 Injection Number : 1
 Sample Name : 15-01-1810-3 Sequence Line : 53
 Instrument Method: C:\CHEM32\1\METHODS\ ->
 Acquired on : 03 Feb 15 03:24 am
 Report Created on: 03 Feb 15 03:33 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Sig. 1 in W:\GC_45\GC 45 DATA\2015\150202\ ->

| Pk | Ret Time | Area | Height | Peak | Width | Response % |
|----|----------|--------|--------|------|-------|------------|
| 1 | 1.941 | 38.91 | 19 | VV | 0.029 | 0.545 |
| 2 | 2.021 | 139.02 | 80 | VV | 0.025 | 1.947 |
| 3 | 2.086 | 56.47 | 34 | VV | 0.024 | 0.791 |
| 4 | 2.165 | 242.37 | 135 | VV | 0.029 | 3.394 |
| 5 | 2.212 | 75.92 | 48 | VV | 0.024 | 1.063 |
| 6 | 2.251 | 147.24 | 86 | VV | 0.025 | 2.062 |
| 7 | 2.317 | 196.66 | 71 | VV | 0.038 | 2.754 |
| 8 | 2.381 | 97.69 | 60 | VV | 0.023 | 1.368 |
| 9 | 2.451 | 685.03 | 451 | VV | 0.024 | 9.592 |
| 10 | 2.539 | 162.78 | 54 | VV | 0.043 | 2.279 |
| 11 | 2.593 | 176.48 | 62 | VV | 0.038 | 2.471 |
| 12 | 2.638 | 127.59 | 61 | VV | 0.030 | 1.787 |
| 13 | 2.714 | 329.70 | 94 | VV | 0.050 | 4.617 |
| 14 | 2.766 | 386.94 | 193 | VV | 0.029 | 5.418 |
| 15 | 2.851 | 475.58 | 92 | VV | 0.064 | 6.660 |
| 16 | 2.939 | 264.81 | 93 | VV | 0.039 | 3.708 |
| 17 | 3.010 | 367.87 | 158 | VV | 0.032 | 5.151 |
| 18 | 3.057 | 314.31 | 90 | VV | 0.047 | 4.401 |
| 19 | 3.159 | 289.92 | 58 | VV | 0.065 | 4.060 |
| 20 | 3.237 | 90.15 | 47 | VV | 0.028 | 1.262 |
| 21 | 3.300 | 309.70 | 58 | VV | 0.068 | 4.337 |
| 22 | 3.377 | 76.47 | 38 | VV | 0.028 | 1.071 |
| 23 | 3.429 | 162.98 | 39 | VV | 0.053 | 2.282 |
| 24 | 3.496 | 114.48 | 39 | VV | 0.040 | 1.603 |
| 25 | 3.554 | 193.72 | 45 | VV | 0.057 | 2.713 |
| 26 | 3.641 | 105.07 | 36 | VV | 0.039 | 1.471 |
| 27 | 3.684 | 373.51 | 39 | VV | 0.118 | 5.230 |
| 28 | 3.939 | 74.80 | 22 | VV | 0.045 | 1.047 |
| 29 | 4.007 | 196.88 | 15 | VV | 0.157 | 2.757 |
| 30 | 4.711 | 15.09 | 2 | VV | 0.097 | 0.211 |
| 31 | 4.884 | 3.55 | 1 | VV | 0.047 | 0.050 |
| 32 | 4.938 | 4.38 | 1 | VV | 0.046 | 0.061 |
| 33 | 5.032 | 6.89 | 2 | VV | 0.058 | 0.097 |
| 34 | 5.181 | 4.24 | 1 | VV | 0.045 | 0.059 |
| 35 | 5.329 | 834.08 | 590 | VV | 0.023 | 11.680 |

Total area =

=====
Area Percent Report
=====

Data File Name : W:\GC_45\GC 45 DATA\2015\150202\15020253.D
Page Number : 2
Operator : 682 Vial Number : Vial 53
Instrument : GC 45 Injection Number : 1
Sample Name : 15-01-1810-3 Sequence Line : 53
Instrument Method: C:\CHEM32\1\METHODS\ ->
Acquired on : 03 Feb 15 03:24 am
Report Created on: 03 Feb 15 03:33 pm Analysis Method : 8015B.MTH

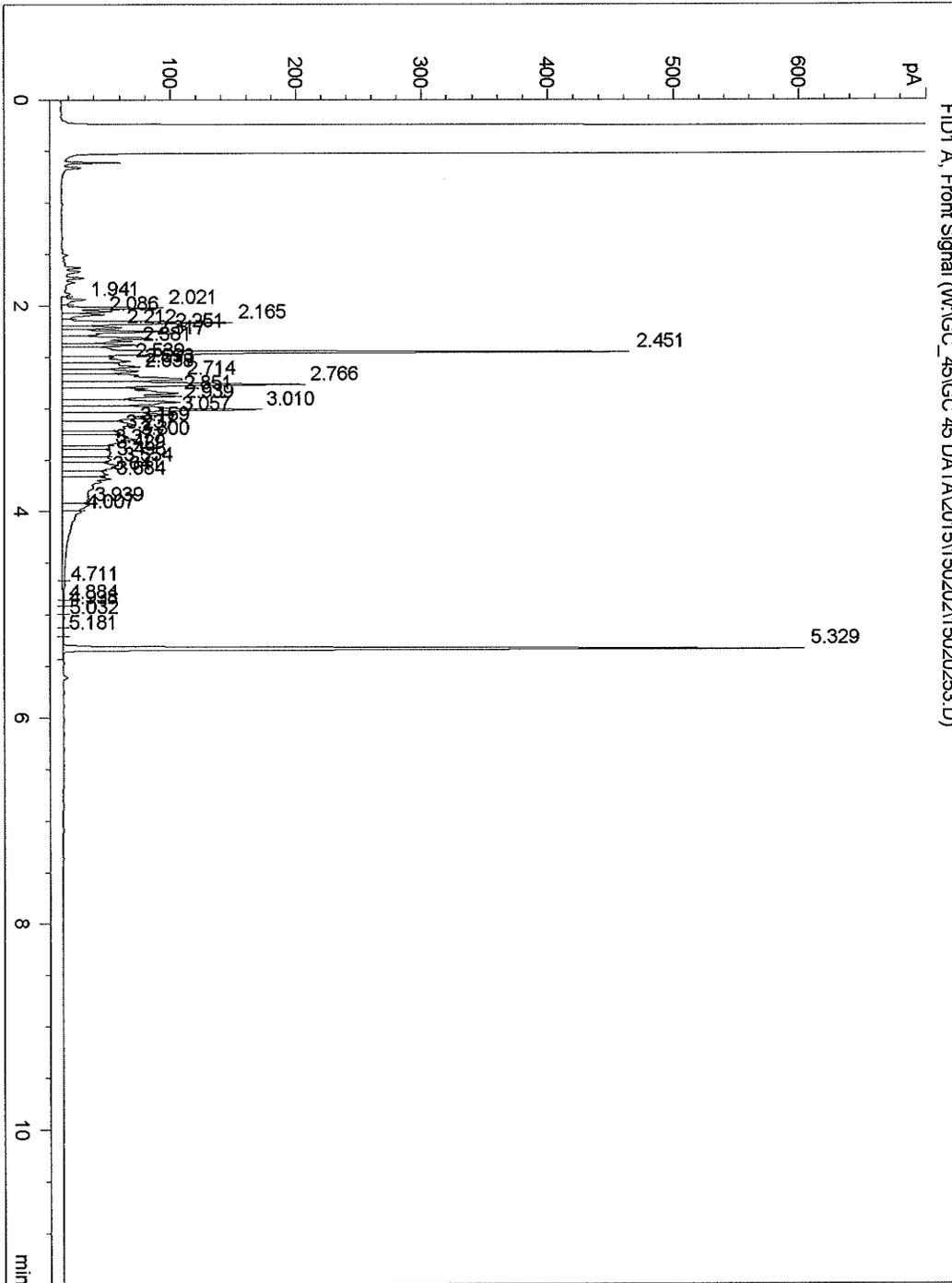
Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

7141.30

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2015\150202\15020253.D
Page Number : 3
Operator : 682 Vial Number : Vial 53
Instrument : GC 45 Injection Number : 1
Sample Name : 15-01-1810-3 Sequence Line : 53
Instrument Method: C:\CHEM32\1\METHODS\ ->
Acquired on : 03 Feb 15 03:24 am
Report Created on: 03 Feb 15 03:33 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies



RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER: 15-01-1810
INSTRUMENT: GC 45
EXTRACTION: EPA 3510C
D/T EXTRACTED: 2015-02-02 00:00 ✓

ANALYZED BY: 960
D/T ANALYZED: 2015-02-03 03:44
REVIEWED BY:
D/T REVIEWED:

DATA FILE: W:\GC 45\GC 45 DATA\2015\150202\15020254.D\15020254 ✓

4 **CLIENT SAMPLE NUMBER: ES123**

LCS/MB BATCH: 150202B09 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 500.00 ml / ACTUAL: 500.00 ml ✓
MS/MSD BATCH: 150202S09 **FINAL VOLUME / WEIGHT:** DEFAULT: 5.00 ml / ACTUAL: 2.50 ml
UNITS: ug/L **ADJUSTMENT RATIO TO PF:** 0.50

COMMENT: Results were evaluated to the MDL (DL), concentrations \geq to the MDL (DL) but $<$ RL (LOQ), if found, are qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

| <u>COMPOUND</u> | <u>INI. CONC</u> | <u>DF</u> | <u>CONC</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>QUAL</u> |
|-----------------|------------------|-----------|-------------|-----------|------------|------------|-------------|
| TPH as Diesel | 7750 | 1.00 ✓ | 38.7 | 2.9 | 10 | 25 | b |

=====
Area Percent Report
=====

Data File Name : W:\GC_45\GC 45 DATA\2015\150202\15020254.D
Page Number : 1
Operator : Vial Number : Vial 54
Instrument : GC 45 Injection Number : 1
Sample Name : 15-01-1810-4 Sequence Line : 54
Instrument Method: C:\CHEM32\1\METHODS\ ->
Acquired on : 03 Feb 15 03:44 am
Report Created on: 03 Feb 15 03:36 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Sig. 1 in W:\GC_45\GC 45 DATA\2015\150202\ ->

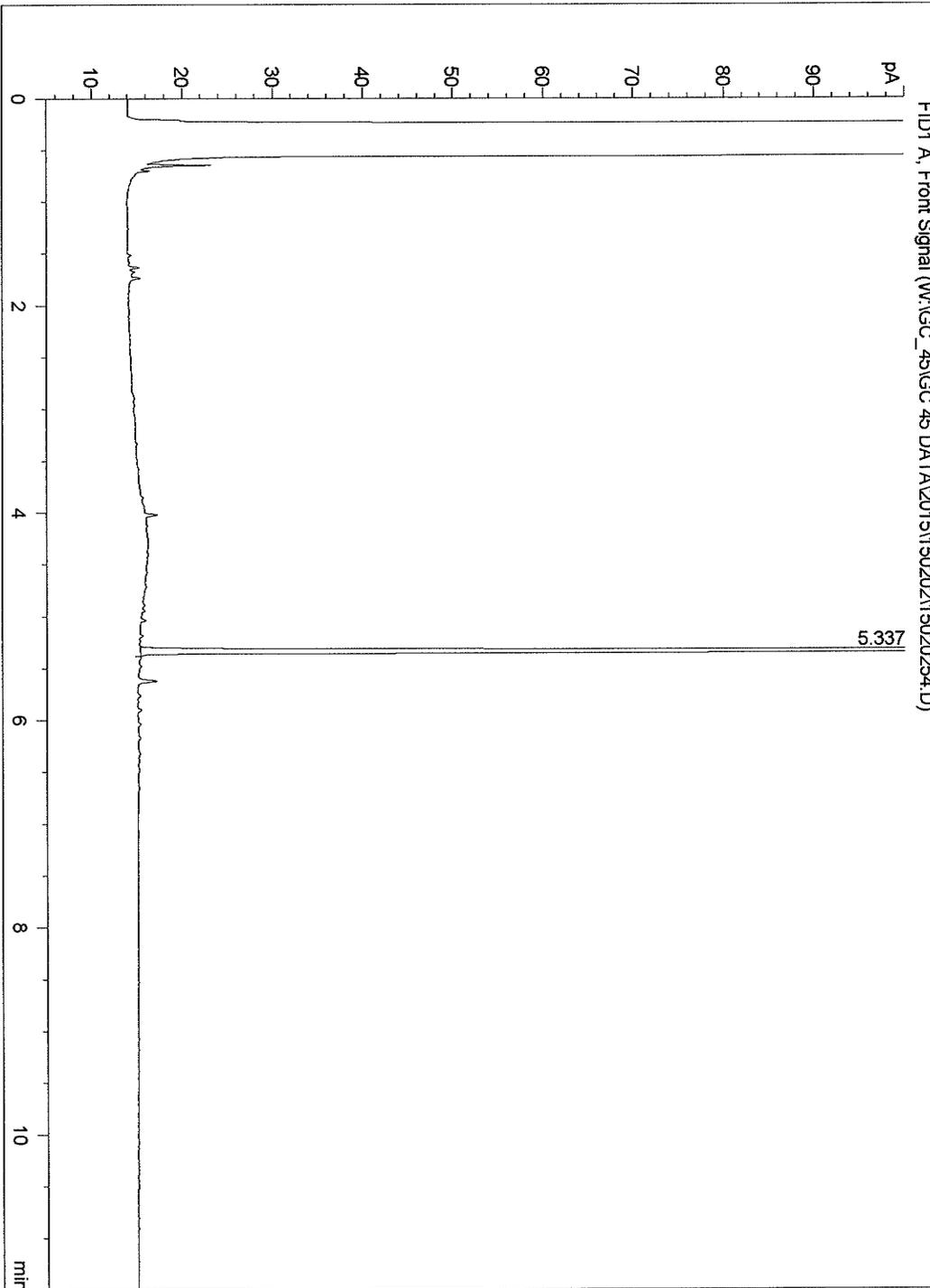
| Pk | Ret Time | Area | Height | Peak | Width | Response % |
|----|----------|--------|--------|------|-------|------------|
| 1 | 5.337 | 571.96 | 386 | BB | 0.023 | 100.000 |
| 2 | 0.000 | 0.00 | 0 | | 0.000 | 0.000 |

Total area = 571.96
0.00

=====
Area Percent Report
=====

Data File Name : W:\GC_45\GC 45 DATA\2015\150202\15020254.D
Page Number : 2
Operator : Vial Number : Vial 54
Instrument : GC 45 Injection Number : 1
Sample Name : 15-01-1810-4 Sequence Line : 54
Instrument Method: C:\CHEM32\1\METHODS\ ->
Acquired on : 03 Feb 15 03:44 am
Report Created on: 03 Feb 15 03:36 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies



=====
 Area Percent Report
 =====

Data File Name : W:\GC_45\GC 45 DATA\2015\150202\15020254.D
 Page Number : 1
 Operator : 682 Vial Number : Vial 54
 Instrument : GC 45 Injection Number : 1
 Sample Name : 15-01-1810-4 Sequence Line : 54
 Instrument Method: C:\CHEM32\1\METHODS\ ->
 Acquired on : 03 Feb 15 03:44 am
 Report Created on: 03 Feb 15 03:34 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Sig. 1 in W:\GC_45\GC 45 DATA\2015\150202\ ->

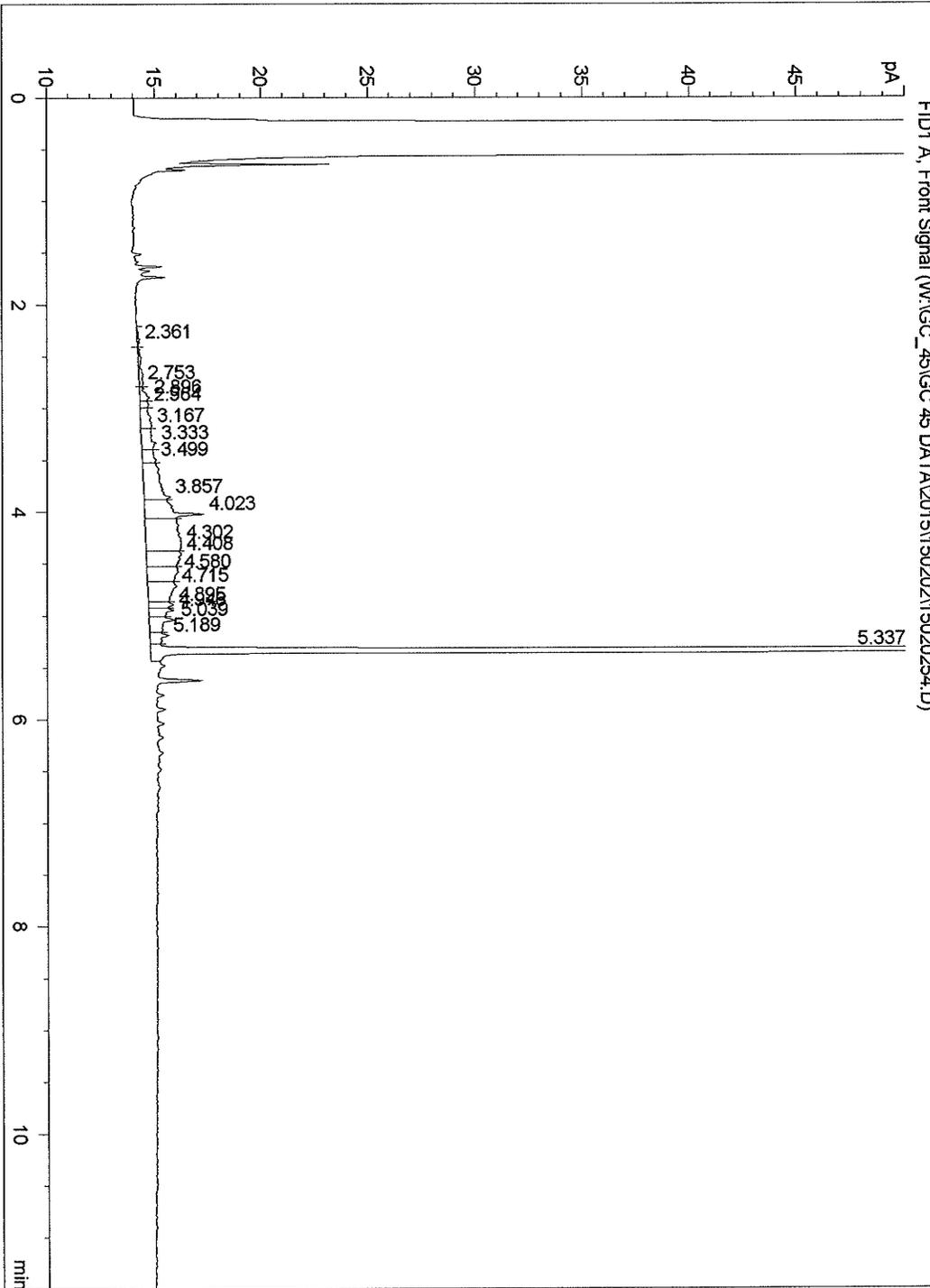
| Pk | Ret Time | Area | Height | Peak | Width | Response % |
|----|----------|--------|--------|--------|-------|------------|
| 1 | 2.361 | 0.67 | | 0 BV | 0.071 | 0.093 |
| 2 | 2.753 | 2.93 | | 0 VV | 0.184 | 0.407 |
| 3 | 2.896 | 2.31 | | 0 VV | 0.066 | 0.321 |
| 4 | 2.964 | 1.47 | | 0 VV | 0.053 | 0.204 |
| 5 | 3.167 | 5.22 | | 1 VV | 0.123 | 0.724 |
| 6 | 3.333 | 6.09 | | 1 VV | 0.110 | 0.845 |
| 7 | 3.499 | 4.24 | | 1 VV | 0.088 | 0.589 |
| 8 | 3.857 | 17.23 | | 1 VV | 0.166 | 2.394 |
| 9 | 4.023 | 15.81 | | 3 VV | 0.073 | 2.195 |
| 10 | 4.302 | 29.63 | | 2 VV | 0.215 | 4.115 |
| 11 | 4.408 | 13.29 | | 2 VV | 0.106 | 1.845 |
| 12 | 4.580 | 11.66 | | 1 VV | 0.097 | 1.619 |
| 13 | 4.715 | 13.16 | | 1 VB | 0.123 | 1.828 |
| 14 | 4.895 | 3.66 | | 1 BV | 0.043 | 0.508 |
| 15 | 4.946 | 4.49 | | 1 VV | 0.051 | 0.623 |
| 16 | 5.039 | 6.60 | | 1 VV | 0.068 | 0.916 |
| 17 | 5.189 | 3.94 | | 1 VV | 0.059 | 0.547 |
| 18 | 5.337 | 577.68 | | 386 VV | 0.024 | 80.226 |

Total area = 720.06

=====
Area Percent Report
=====

Data File Name : W:\GC_45\GC 45 DATA\2015\150202\15020254.D
Page Number : 2
Operator : 682 Vial Number : Vial 54
Instrument : GC 45 Injection Number : 1
Sample Name : 15-01-1810-4 Sequence Line : 54
Instrument Method: C:\CHEM32\1\METHODS\ ->
Acquired on : 03 Feb 15 03:44 am
Report Created on: 03 Feb 15 03:34 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies



RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER: 099-15-516
INSTRUMENT: GC 45
EXTRACTION: EPA 3510C ✓
D/T EXTRACTED: 2015-02-02 00:00 ✓

ANALYZED BY: 960
D/T ANALYZED: 2015-02-03 00:43 ✓
REVIEWED BY:
D/T REVIEWED:

DATA FILE: W:\GC 45\GC 45 DATA\2015\150202\15020245.D\15020245 ✓

MB **CLIENT SAMPLE NUMBER:** Method Blank

LCS/MB BATCH: 150202B09 **SAMPLE VOLUME / WEIGHT:** DEFAULT: 500.00 ml / ACTUAL: 500.00 ml
MS/MSD BATCH: **FINAL VOLUME / WEIGHT:** DEFAULT: 5.00 ml / ACTUAL: 2.50 ml
UNITS: ug/L **ADJUSTMENT RATIO TO PF:** 0.50

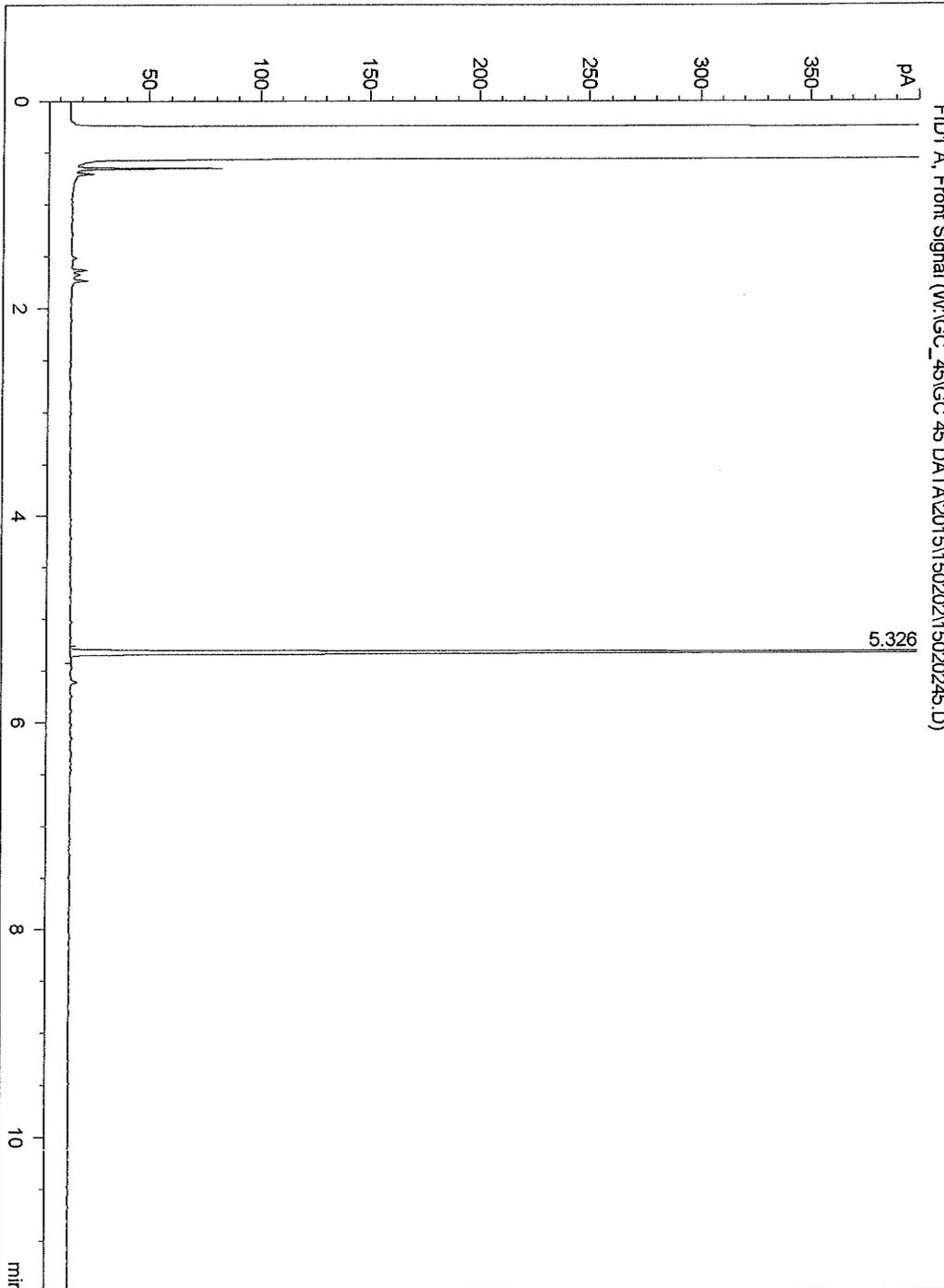
COMMENT:

| <u>COMPOUND</u> | <u>INI. CONC</u> | <u>DF</u> | <u>CONC</u> | <u>DL</u> | <u>LOD</u> | <u>LOQ</u> | <u>QUAL</u> |
|-----------------|------------------|-----------|-------------|-----------|------------|------------|-------------|
| TPH as Diesel | 0.000 | 1.00 ✓ | ND | 2.9 | 10 | 25 | |

=====
Area Percent Report
=====

Data File Name : W:\GC_45\GC 45 DATA\2015\150202\15020245.D
Page Number : 2
Operator : 682 Vial Number : Vial 45
Instrument : GC 45 Injection Number : 1
Sample Name : MB 15020209/10/11 Sequence Line : 45
Instrument Method: C:\CHEM32\1\METHODS\ ->
Acquired on : 03 Feb 15 00:43 am
Report Created on: 03 Feb 15 03:32 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies



**CONTINUING CALIBRATION VERIFICATION QUALITY CONTROL SHEET
FOR METHOD: EPA 8015B (M)**

CCV WORK ORDER: 099-15-515-363-6279

BATCH ID: 1502021018
INITIAL: 150202A090
CCV: GC 45
INSTRUMENT:

DATA FILE: W:\GC_45\GC 45 DATA\2015\150202\15020243.D\15020243 /

ANALYZED BY: 960

D/T ANALYZED: 2015-02-02 11:34
INITIAL: 2015-02-03 00:03
CCV:

REVIEWED BY:
D/T REVIEWED:

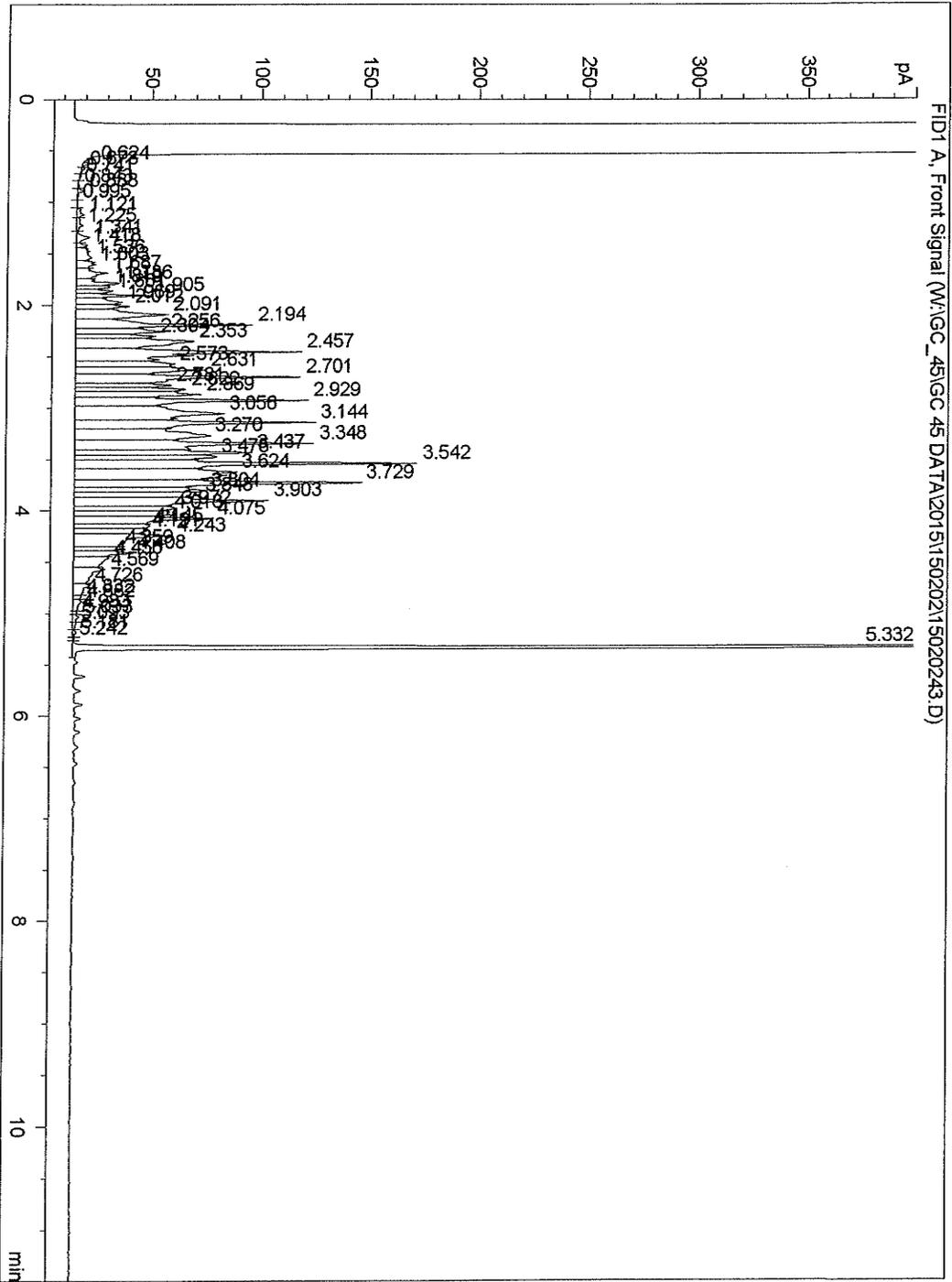
| <u>COMPOUND NAME</u> | <u>COMP TYPE</u> | <u>CALIB MODEL</u> | <u>MIN RF</u> | <u>AVG RF</u> | <u>CCV RF</u> | <u>AMOUNT</u> | <u>CCV CONC</u> | <u>CCV %D</u> | <u>CCV %D CL</u> | <u>STATUS</u> |
|----------------------|------------------|--------------------|---------------|---------------|---------------|---------------|-----------------|---------------|------------------|---------------|
| TPH as Diesel | C | Avg Resp | 0.00 | 0.019 | 0.019 | | | 1 | 0-15 | PASS / |

MIN RF: Method Specified Minimum Response Factor

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2015\150202\15020243.D
Page Number : 3
Operator : 682 Vial Number : Vial 43
Instrument : GC 45 Injection Number : 1
Sample Name : D400 C28 50 L012815D Sequence Line : 43
Instrument Method: C:\CHEM32\1\METHODS\ ->
Acquired on : 03 Feb 15 00:03 am
Report Created on: 03 Feb 15 03:32 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies



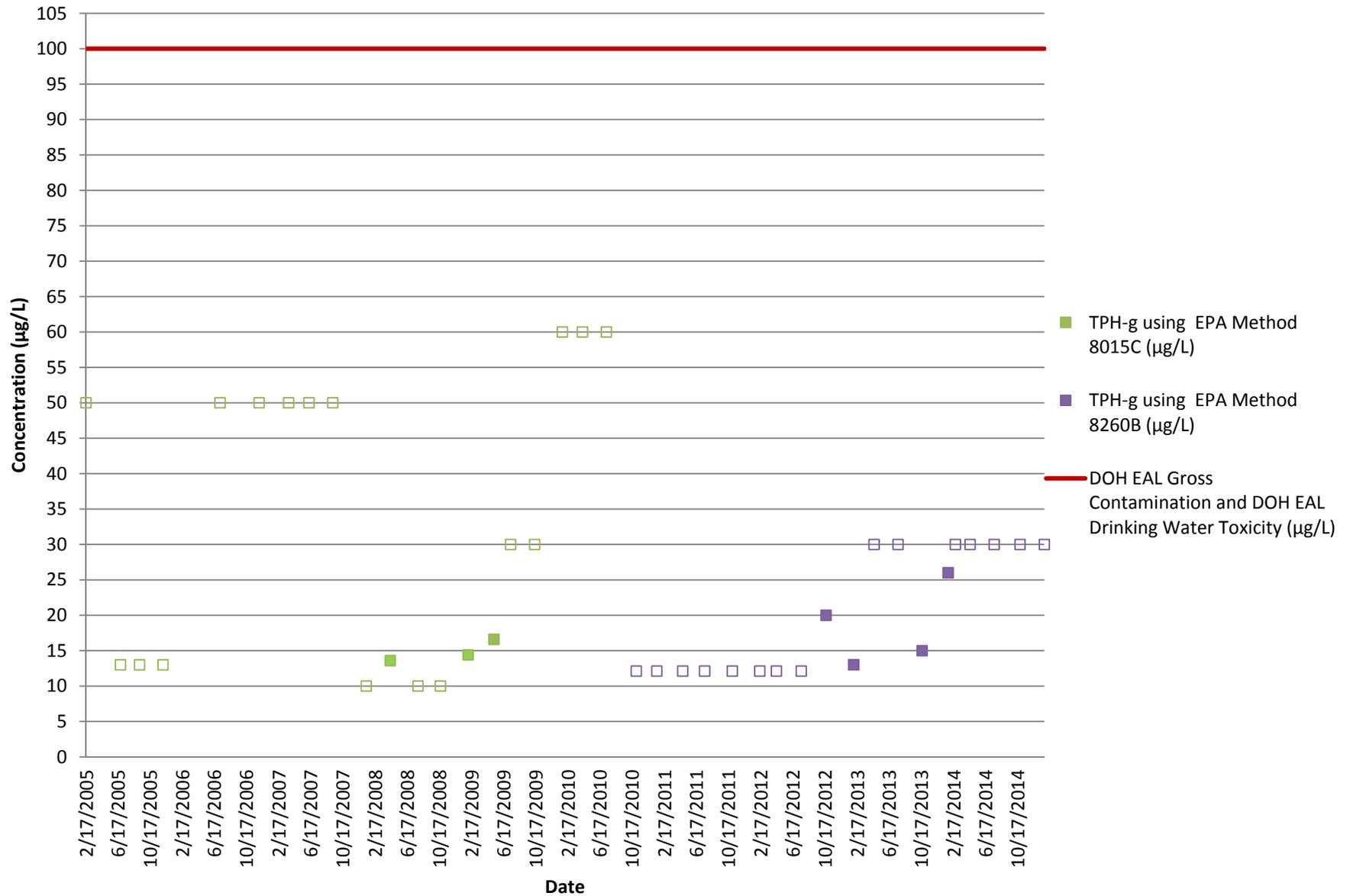
This Page Intentionally Left Blank.

APPENDIX D

Historical Groundwater Exceedance Trends

This Page Intentionally Left Blank

TPH-g Concentrations for RHMW01



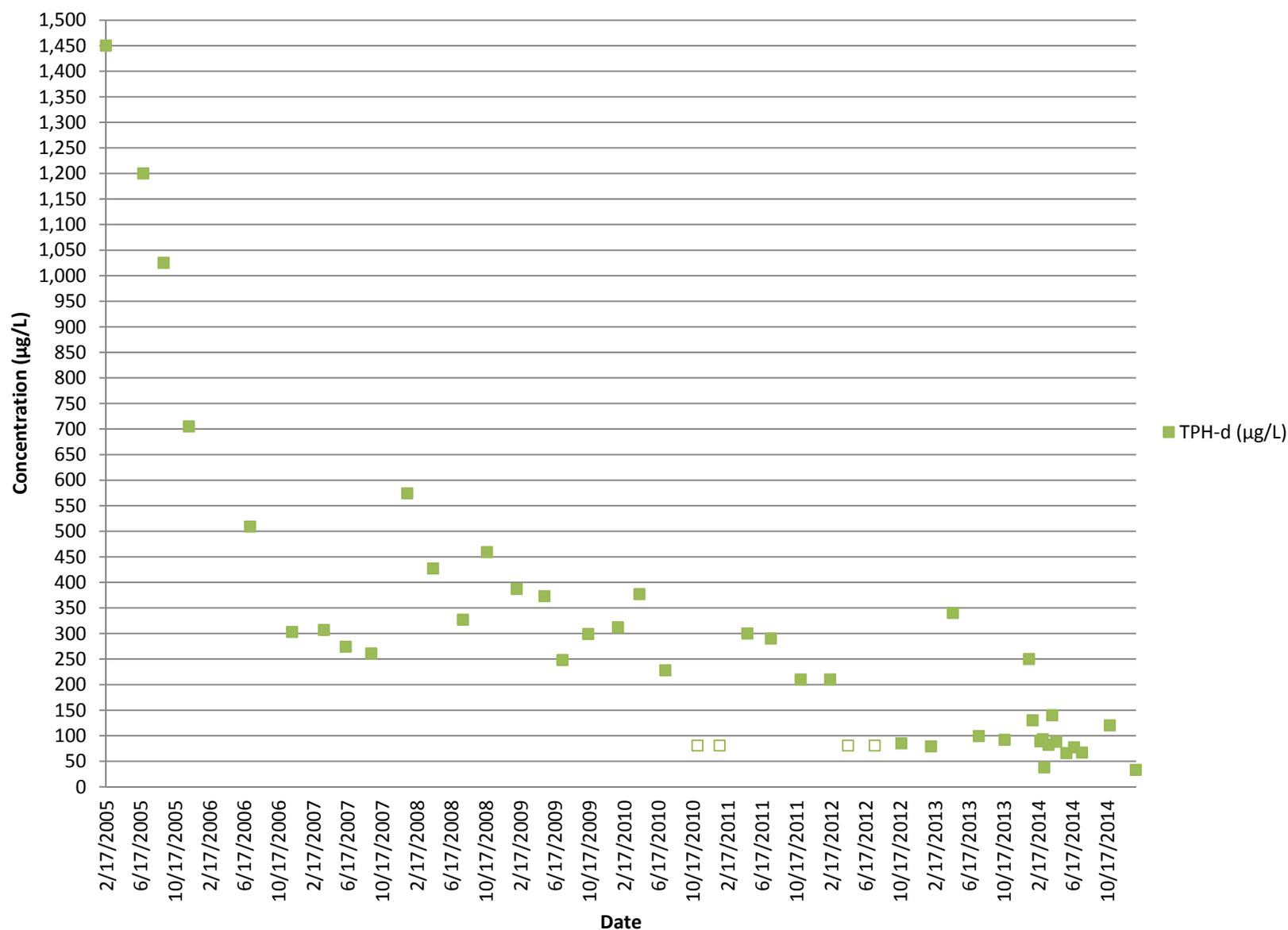
Data points for 2/17/2005 through 9/8/2005 and 12/6/2005 are the average of the primary and duplicate samples.

Possible laboratory contamination for 10/22/2012, 10/21/2013, and 1/28/2014 sampling events.

Unfilled boxes indicate non-detections. Method reporting limits (MRLs) are shown for February 2005, method detection limits (MDLs) are shown for June 2005 through October 2009, and limits of detection (LODs) are shown from January 2010 on.

This Page Intentionally Left Blank.

TPH-d Concentrations for RHMW01



Unfilled boxes indicate non-detections. LODs are shown.

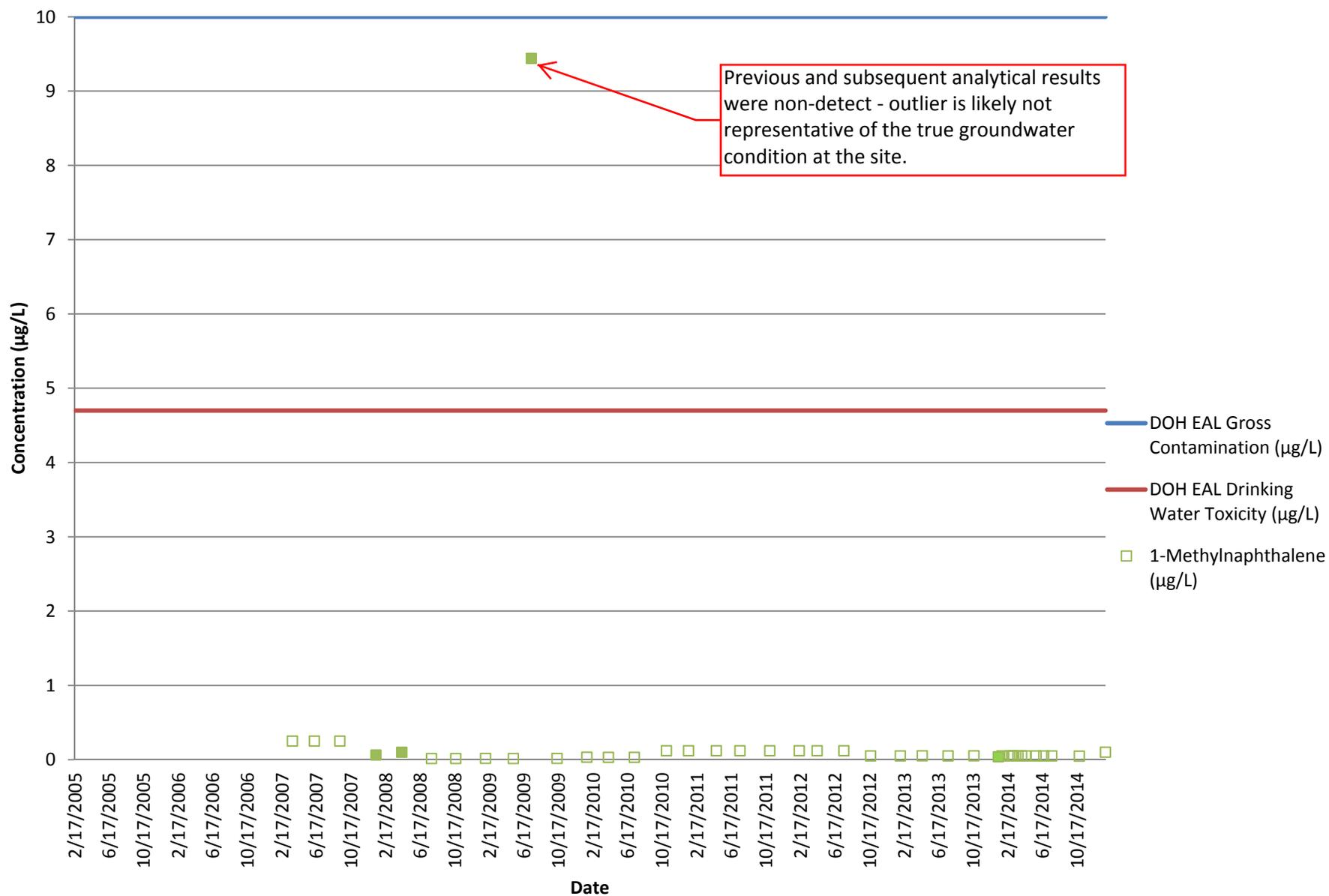
The Site-Specific Risk-Based Level (SSRBL) is 4,500 µg/L.

Numerous sample results had a chromatographic pattern that did not match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.

Data points for 2/17/2005 through 9/8/2005 and 12/6/2005 are the average of the primary and duplicate samples.

This Page Intentionally Left Blank.

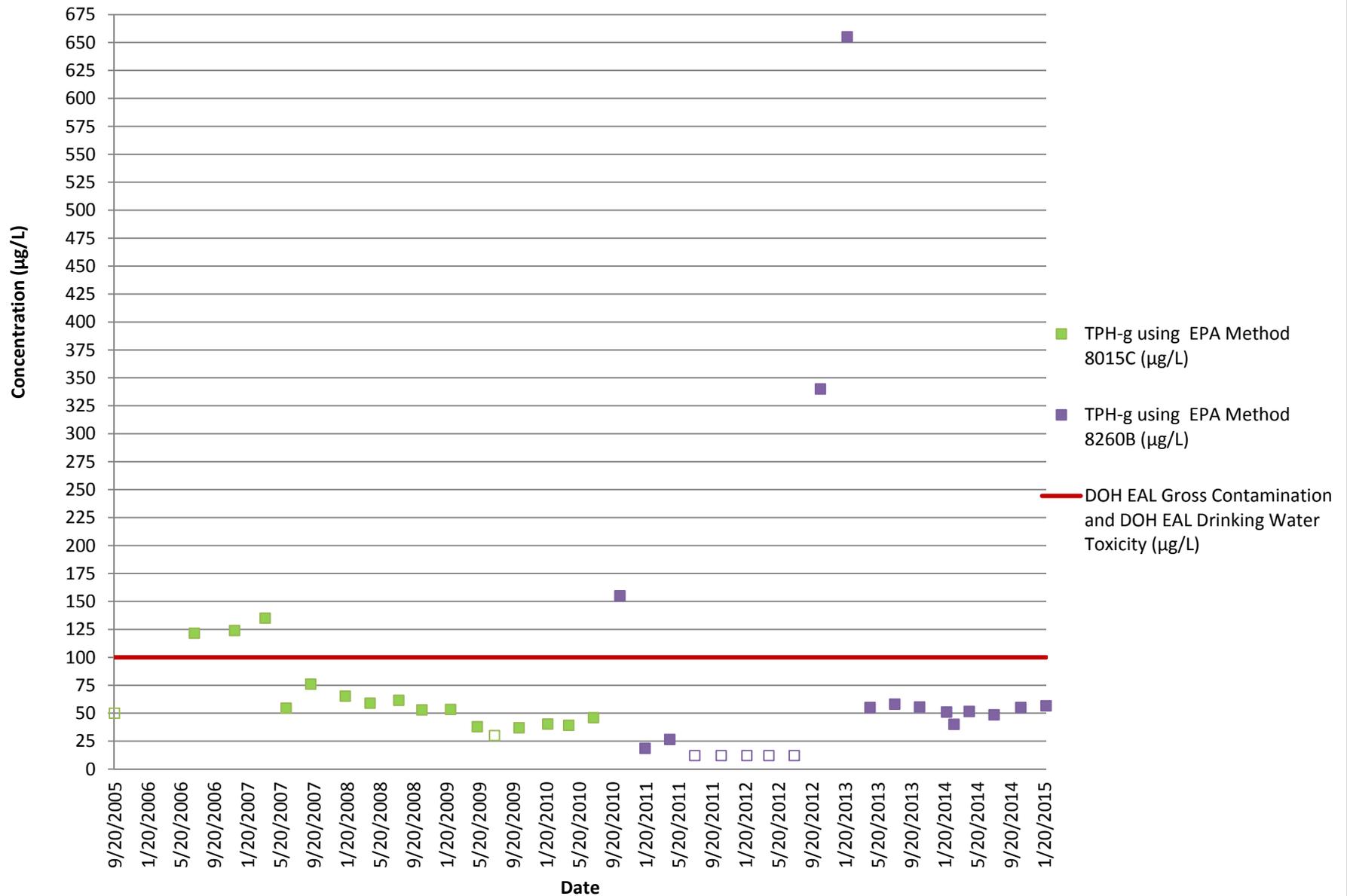
1-Methylnaphthalene Concentrations for RHMW01



Unfilled boxes indicate non-detections. MDLs are shown for June 2005 through October 2009, and LODs are shown from January 2010 on.

This Page Intentionally Left Blank.

TPH-g Concentrations for RHMW02



Data points for 9/20/2005 through 4/21/2014 are the average of the primary and duplicate samples.

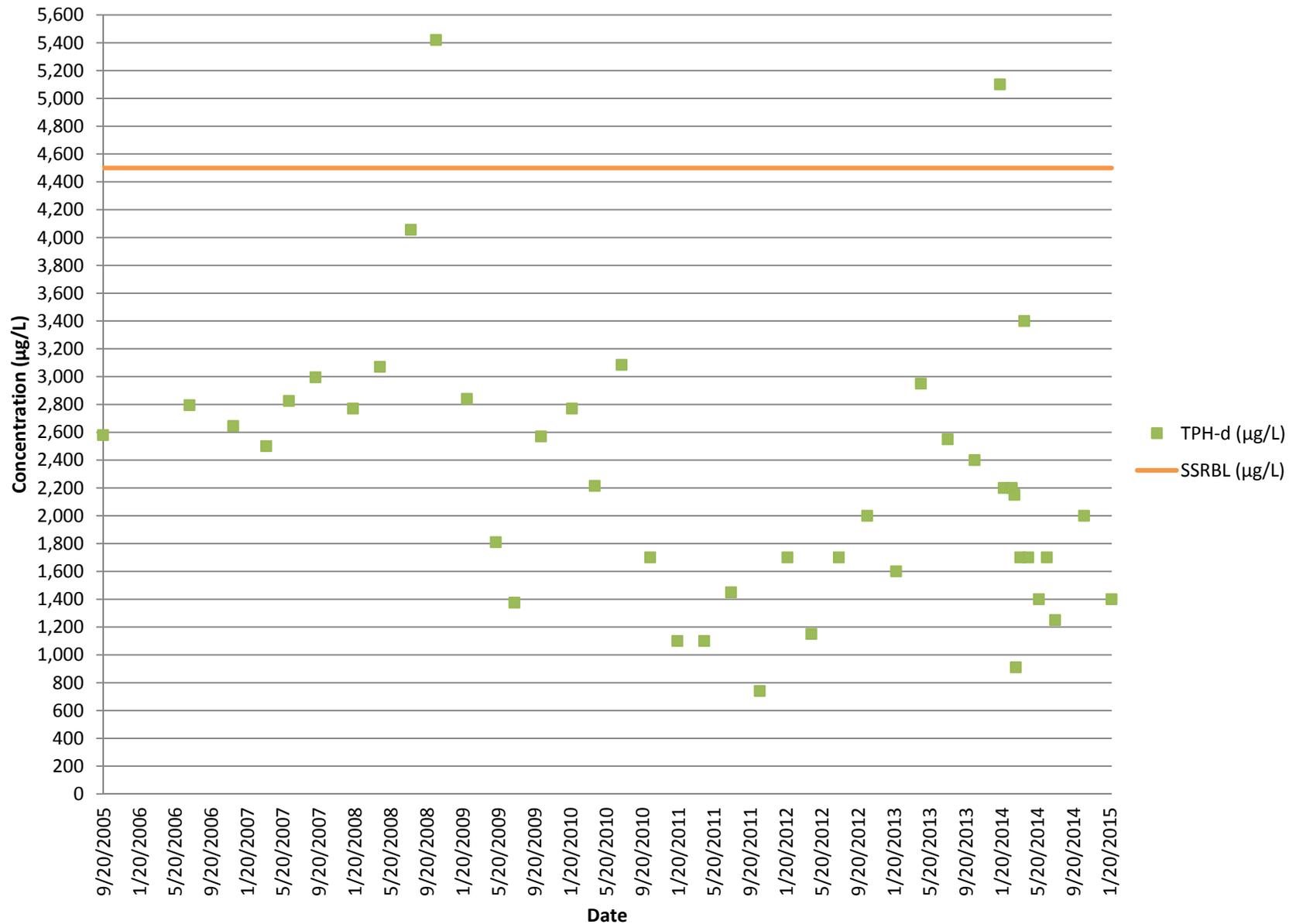
Possible laboratory contamination for 10/21/2013 and 1/28/2014 sampling events.

Unfilled boxes indicate non-detections. MDLs are shown for July 2009, and LODs are shown for September 2005 and from July 2011 on.

Primary sample results are shown for 1/26/2012 and 7/18/2012; all other concentrations are the average of the primary and duplicate sample results.

This Page Intentionally Left Blank.

TPH-d Concentrations for RHMW02

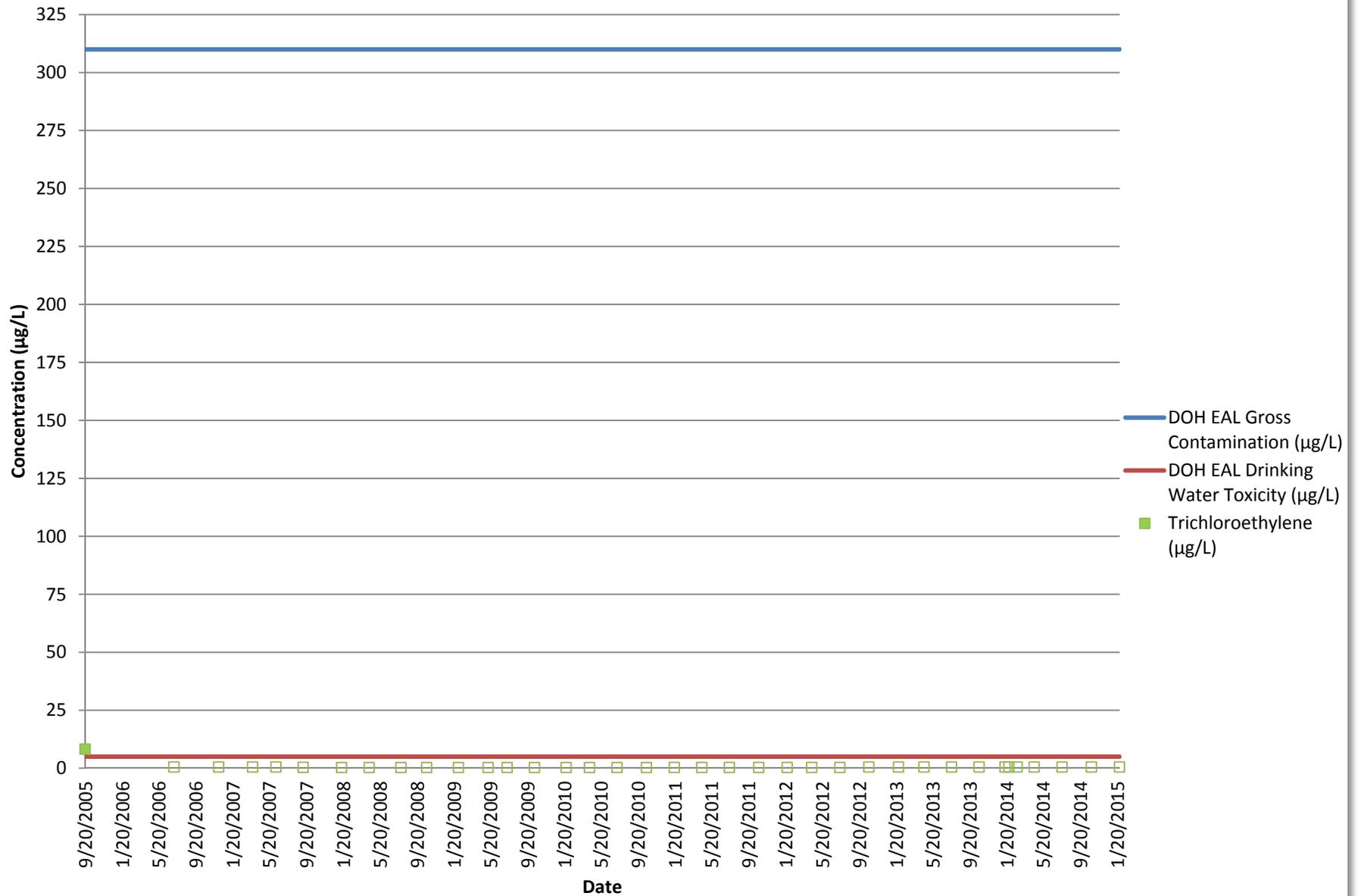


Data points for 9/20/2005 through 4/21/2014 are the average of the primary and duplicate samples.

Numerous sample results had a chromatographic pattern that did not match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.

This Page Intentionally Left Blank.

Trichloroethylene Concentrations for RHMW02

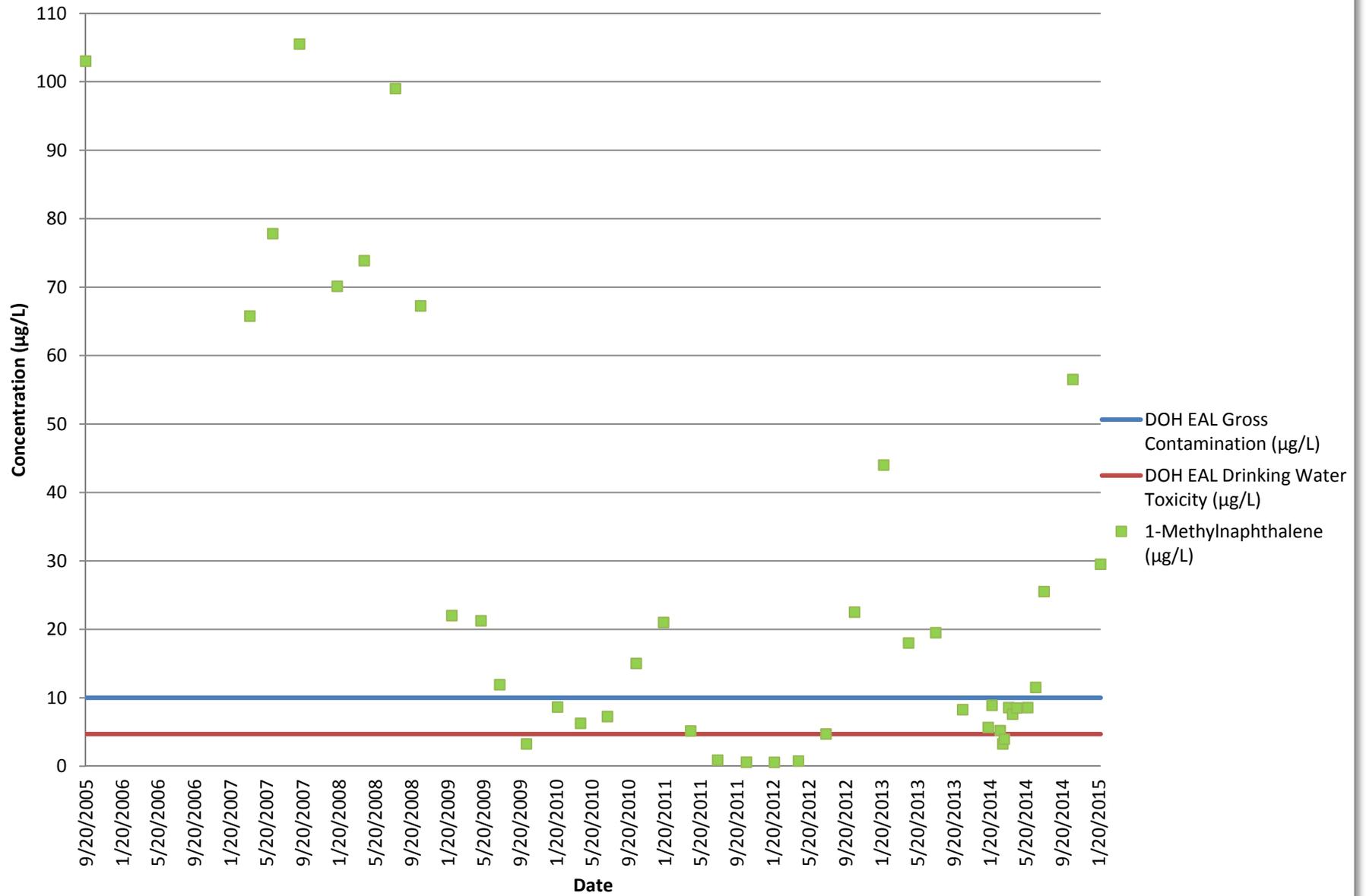


Data points for 9/20/2005 through 4/21/2014 are the average of the primary and duplicate samples.

Unfilled boxes indicate non-detections. MDLs are shown for July 2006 through October 2009, and LODs are shown from January 2010 on.

This Page Intentionally Left Blank.

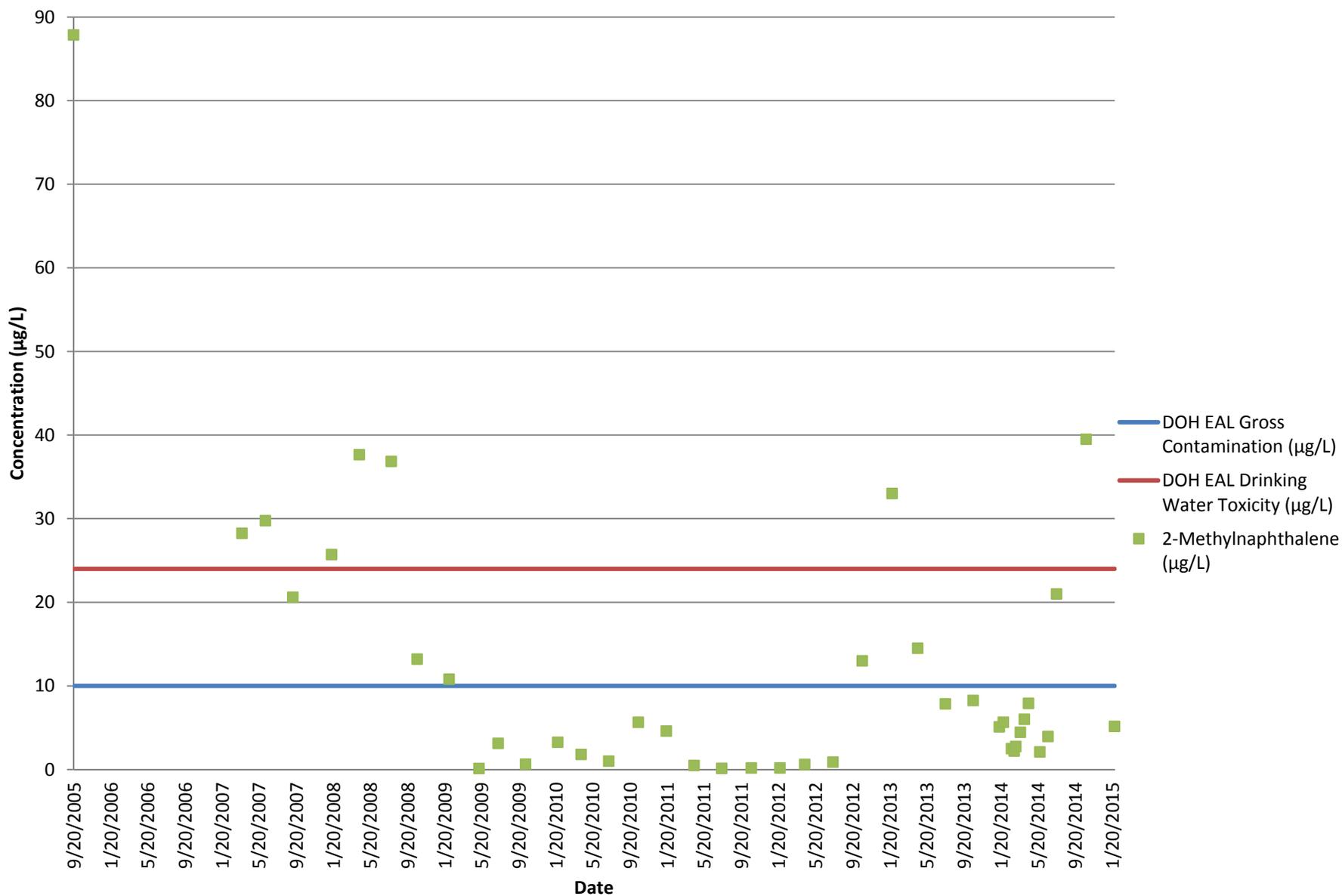
1-Methylnaphthalene Concentrations for RHMW02



Data points for 9/20/2005 and 3/27/2007 through 4/21/2014 are the average of the primary and duplicate samples.

This Page Intentionally Left Blank.

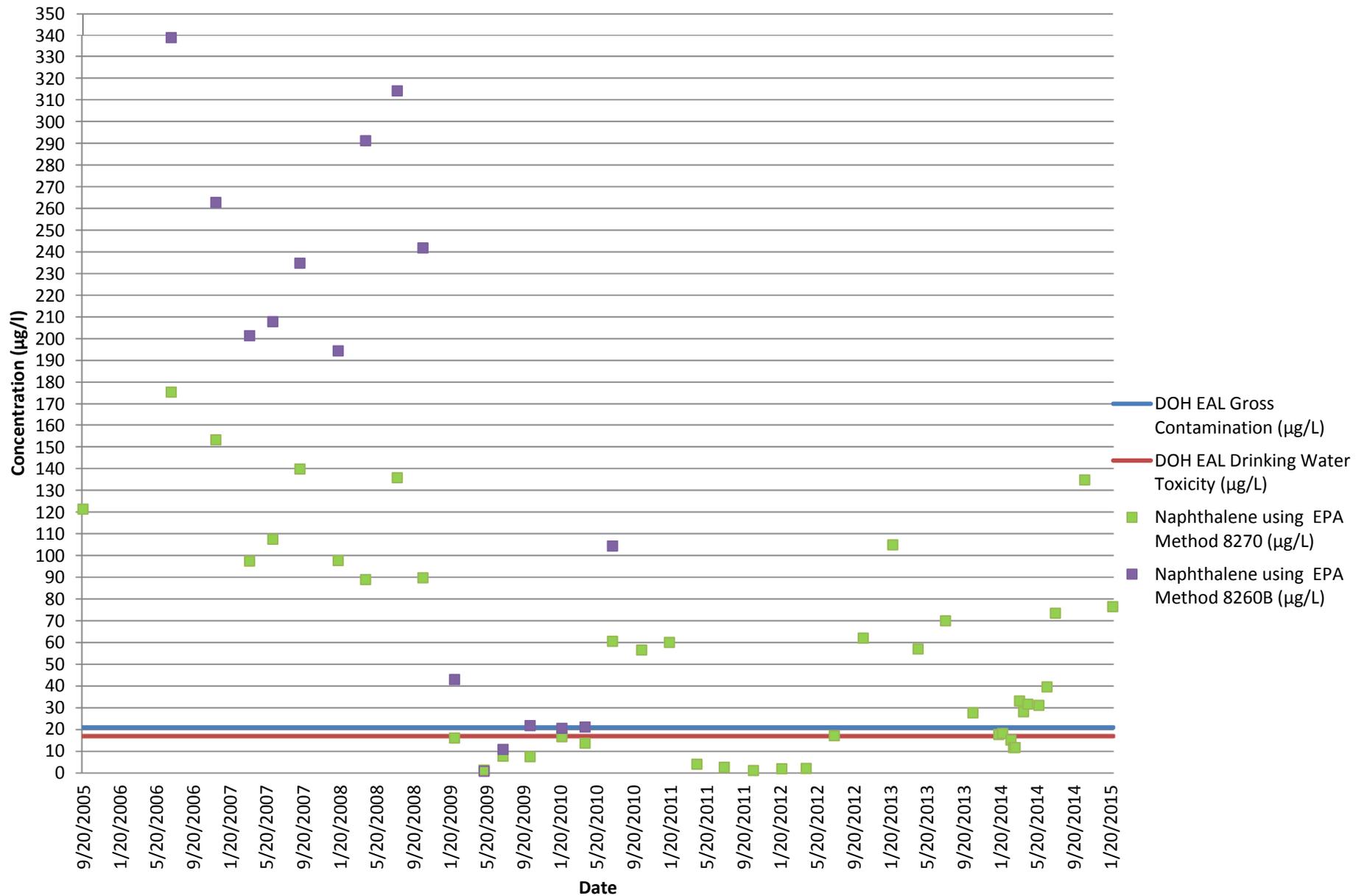
2-Methylnaphthalene Concentrations for RHMW02



Data points for 9/20/2005 and 3/27/2007 through 4/21/2014 are the average of the primary and duplicate samples.

This Page Intentionally Left Blank.

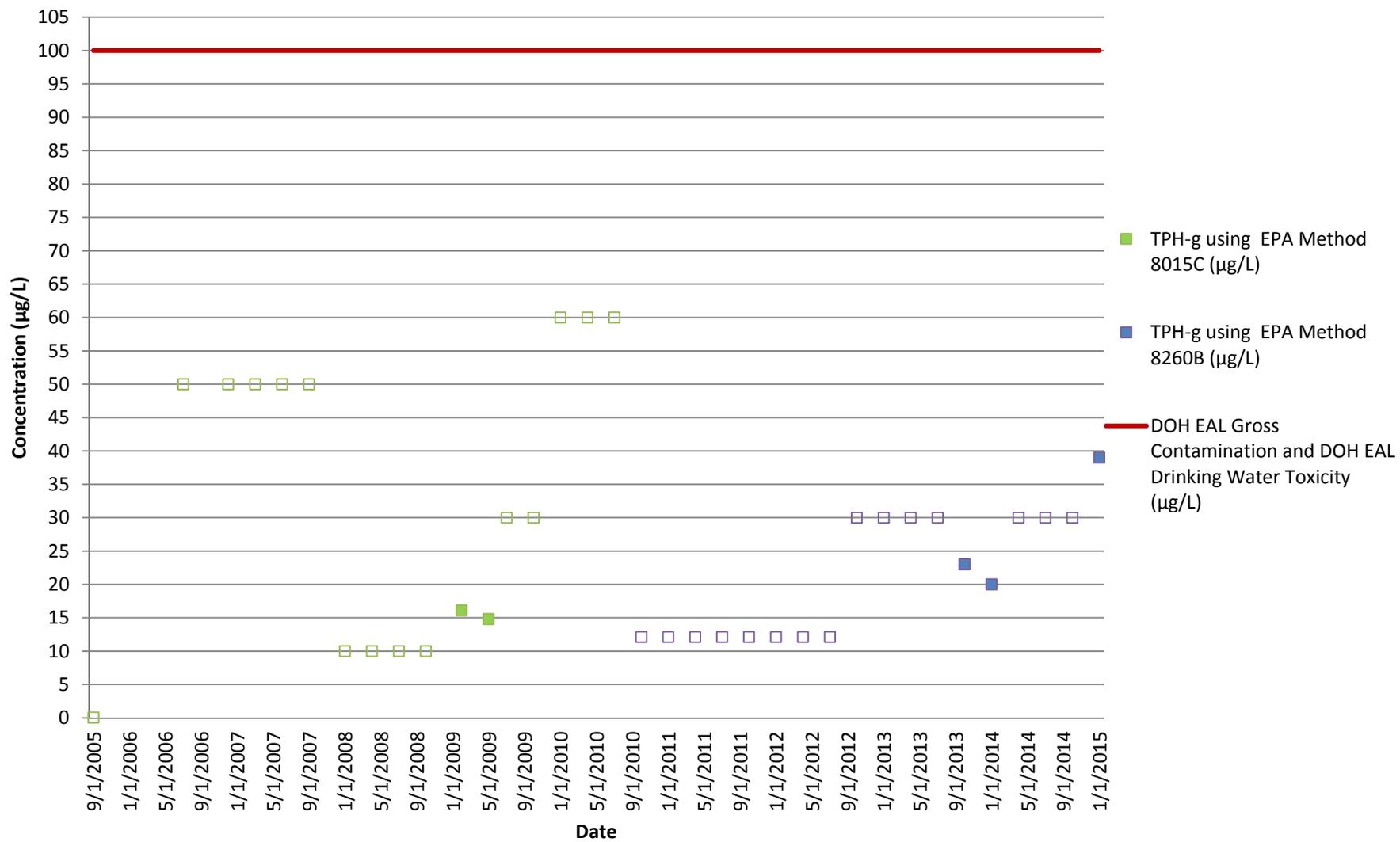
Naphthalene Concentrations for RHMW02



Possible laboratory contamination for 10/21/2013 and 1/28/2014 sampling events.
 Unfilled boxes indicate non-detections. LODs are shown.

This Page Intentionally Left Blank.

TPH-g Concentrations for RHMW03

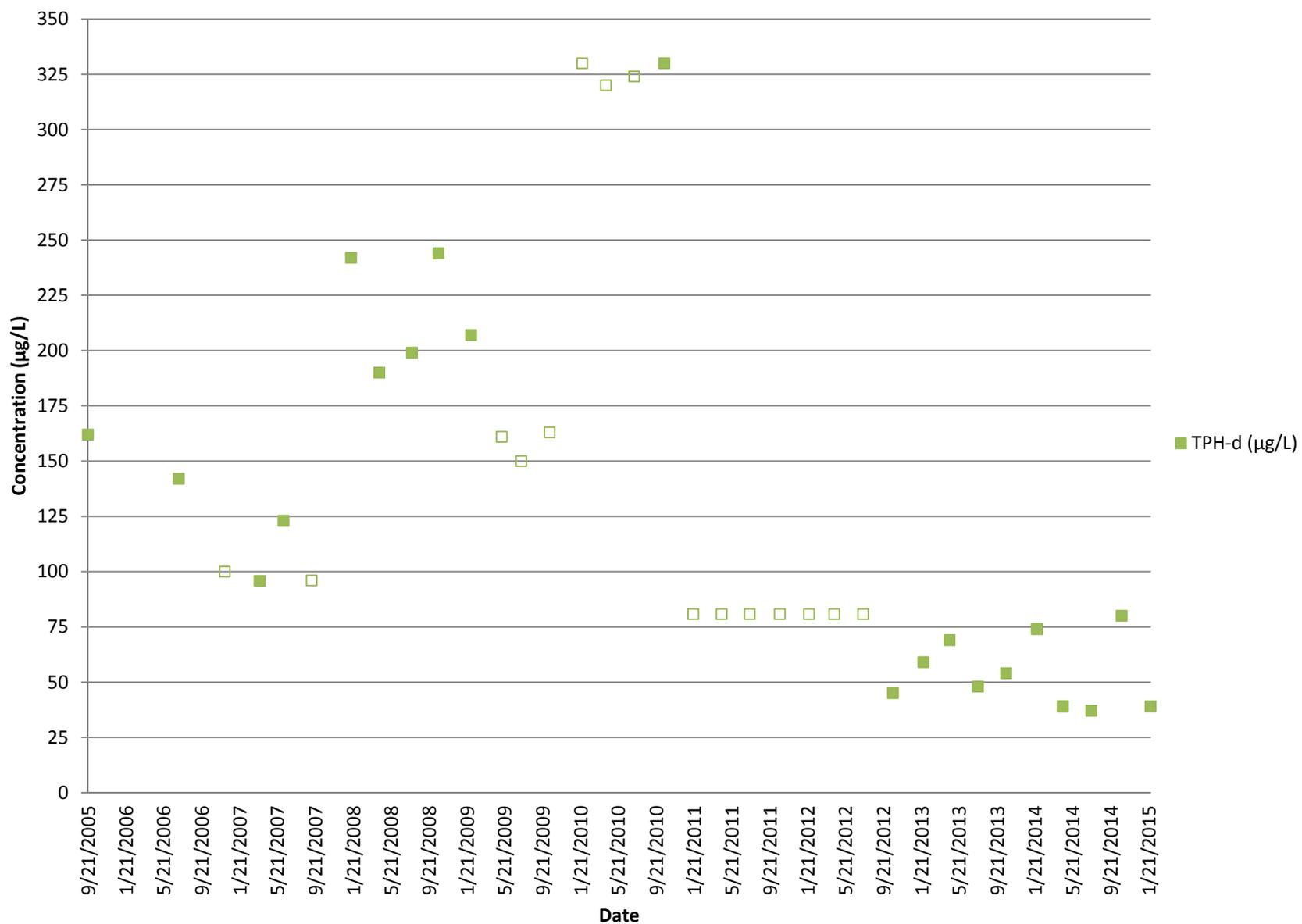


Possible laboratory contamination for 10/21/2013 and 1/28/2014 sampling events.

Unfilled boxes indicate non-detections. MDLs are shown for July 2006 through October 2009, and LODs are shown for September 2005 and from January 2010 on.

This Page Intentionally Left Blank.

TPH-d Concentrations for RHMW03



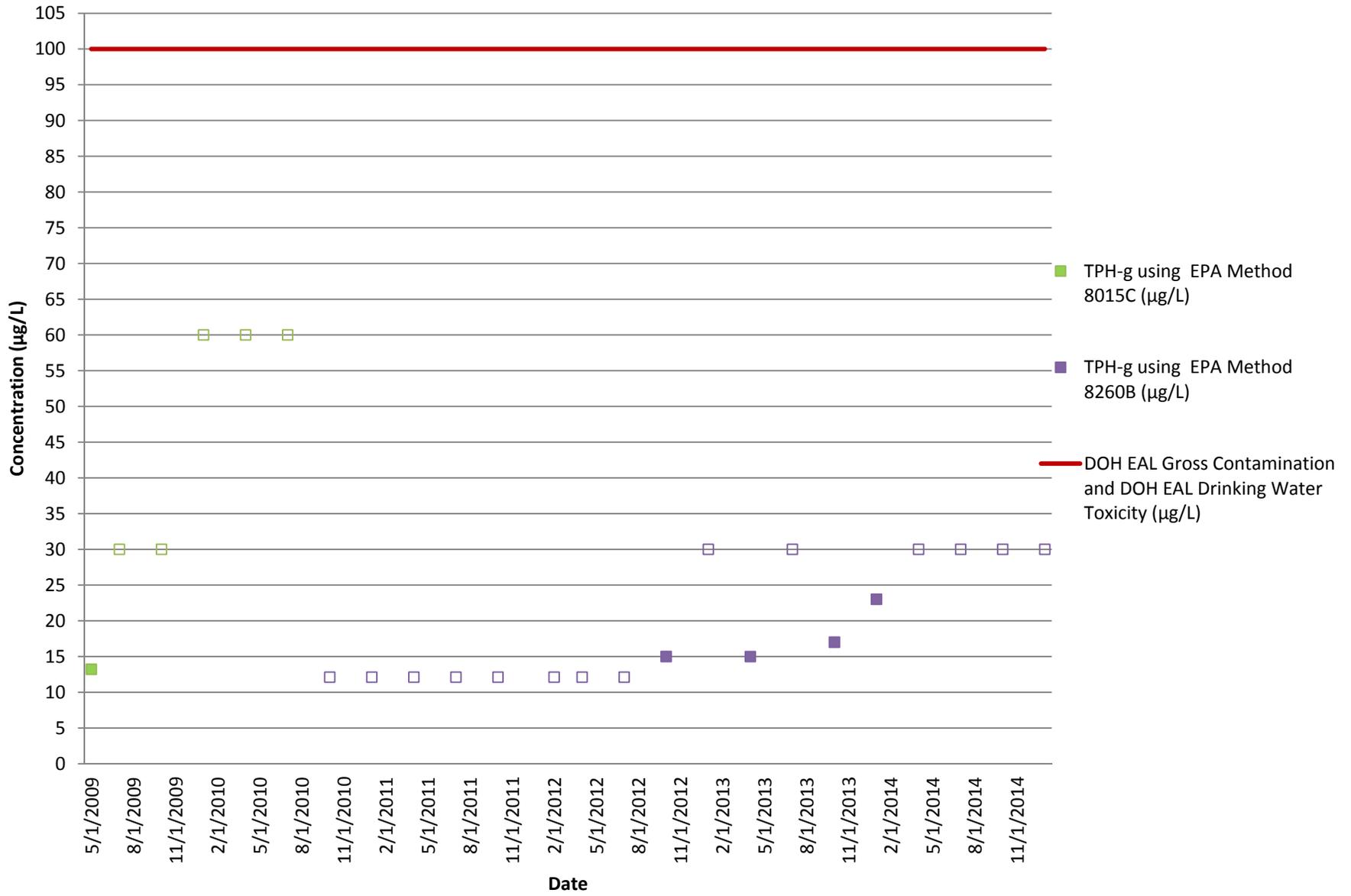
The Site-Specific Risk-Based Level (SSRBL) is 4,500 µg/L.

Unfilled boxes indicate non-detections. MDLs are shown for December 2006 through October 2009, and LODs are shown for September 2005 and from January 2010 on.

Numerous sample results had a chromatographic pattern that didn't match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.

This Page Intentionally Left Blank.

TPH-g Concentrations for RHMW05



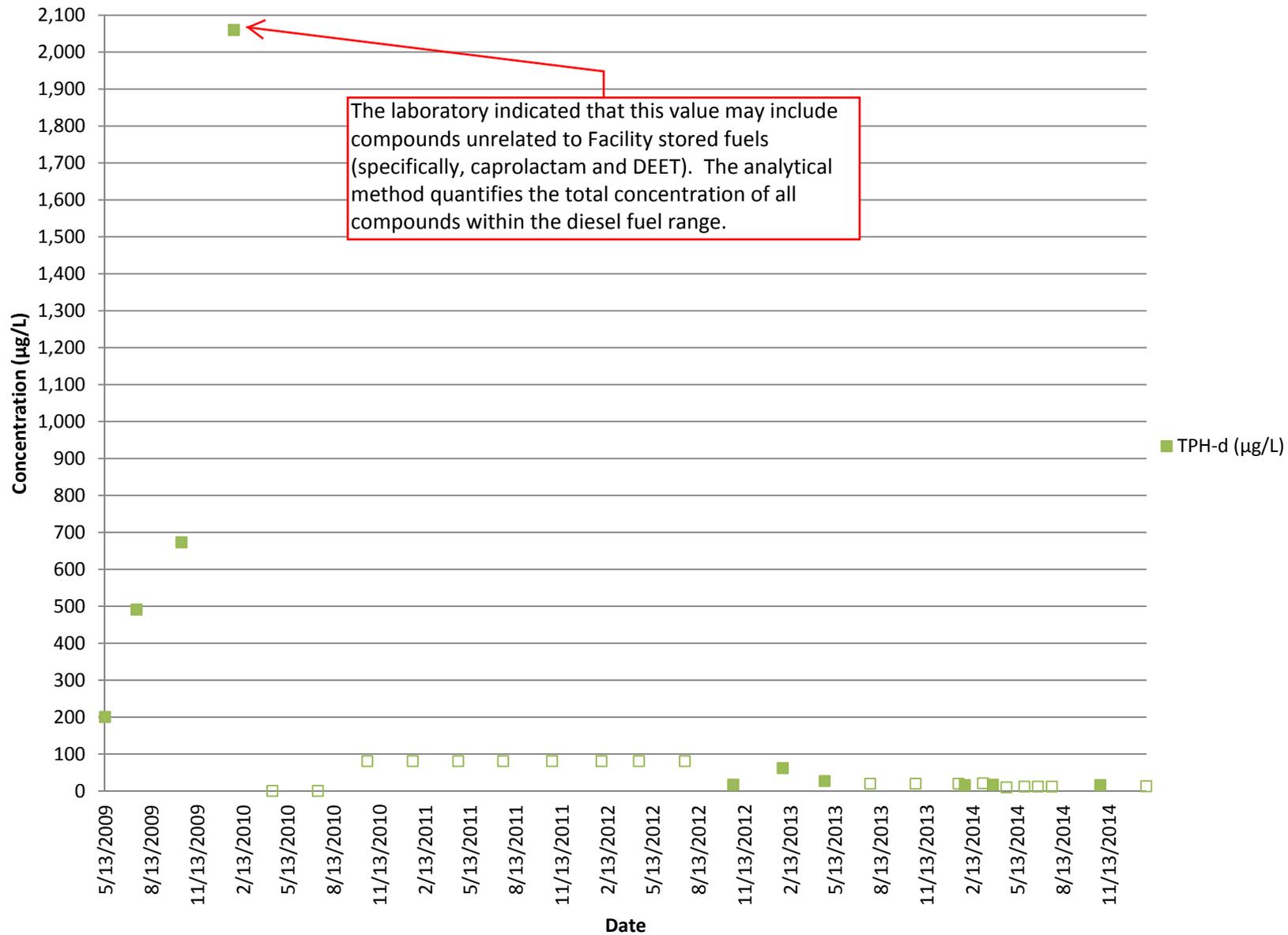
Data point for 7/17/2012 is the average of the primary and duplicate samples.

Unfilled boxes indicate non-detections. MDLs are shown for July and October 2009, and LODs are shown from January 2010 on.

Possible laboratory contamination for 10/22/2012, 10/22/2013, and 1/29/2014 sampling events.

This Page Intentionally Left Blank.

TPH-d Concentrations for RHMW05



The Site-Specific Risk-Based Level (SSRBL) is 4,500 µg/L.

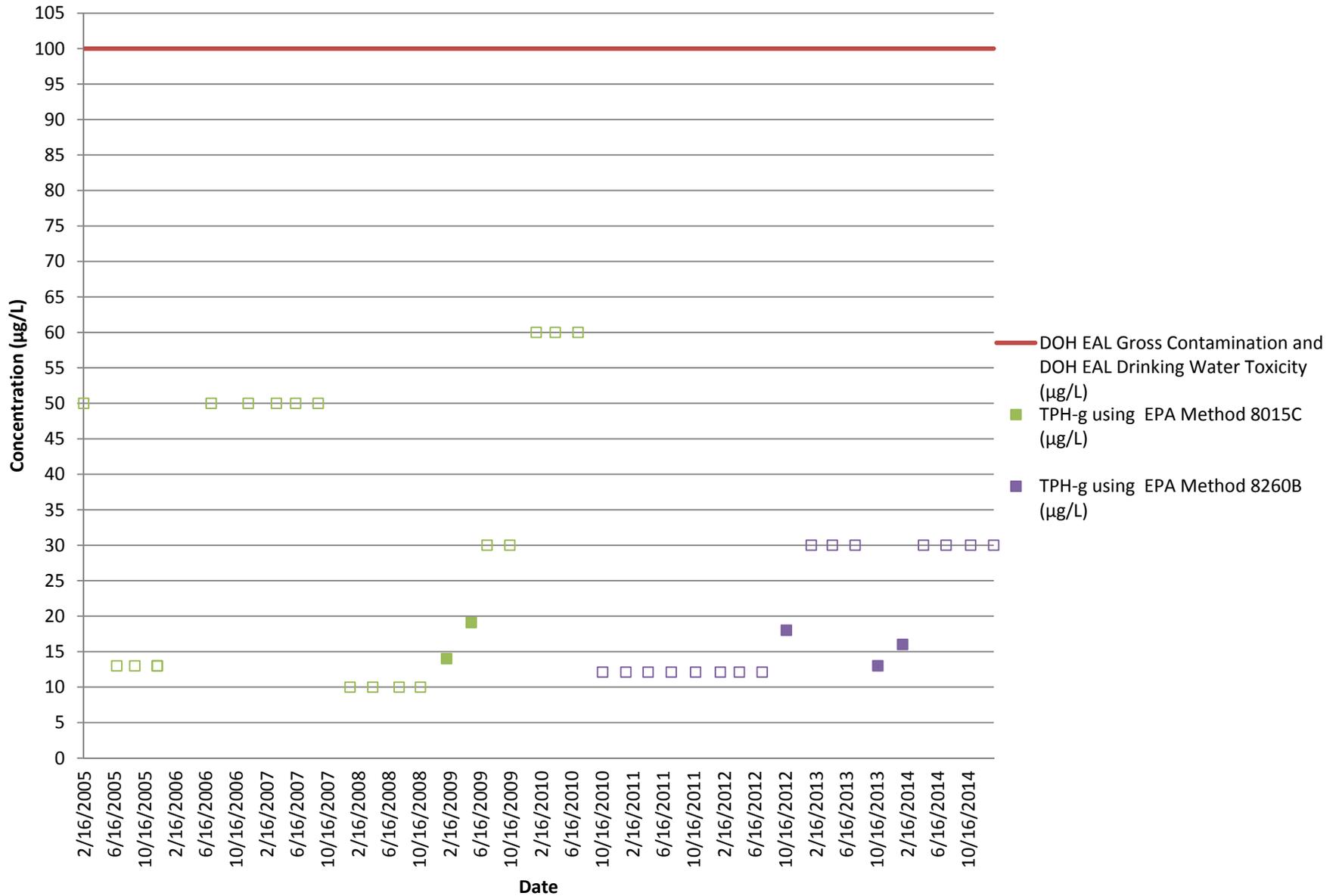
Data point for 7/17/2012 is the average of the primary and duplicate samples.

Unfilled boxes indicate non-detections. LODs are shown.

Numerous sample results had a chromatographic pattern that did not match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.

This Page Intentionally Left Blank.

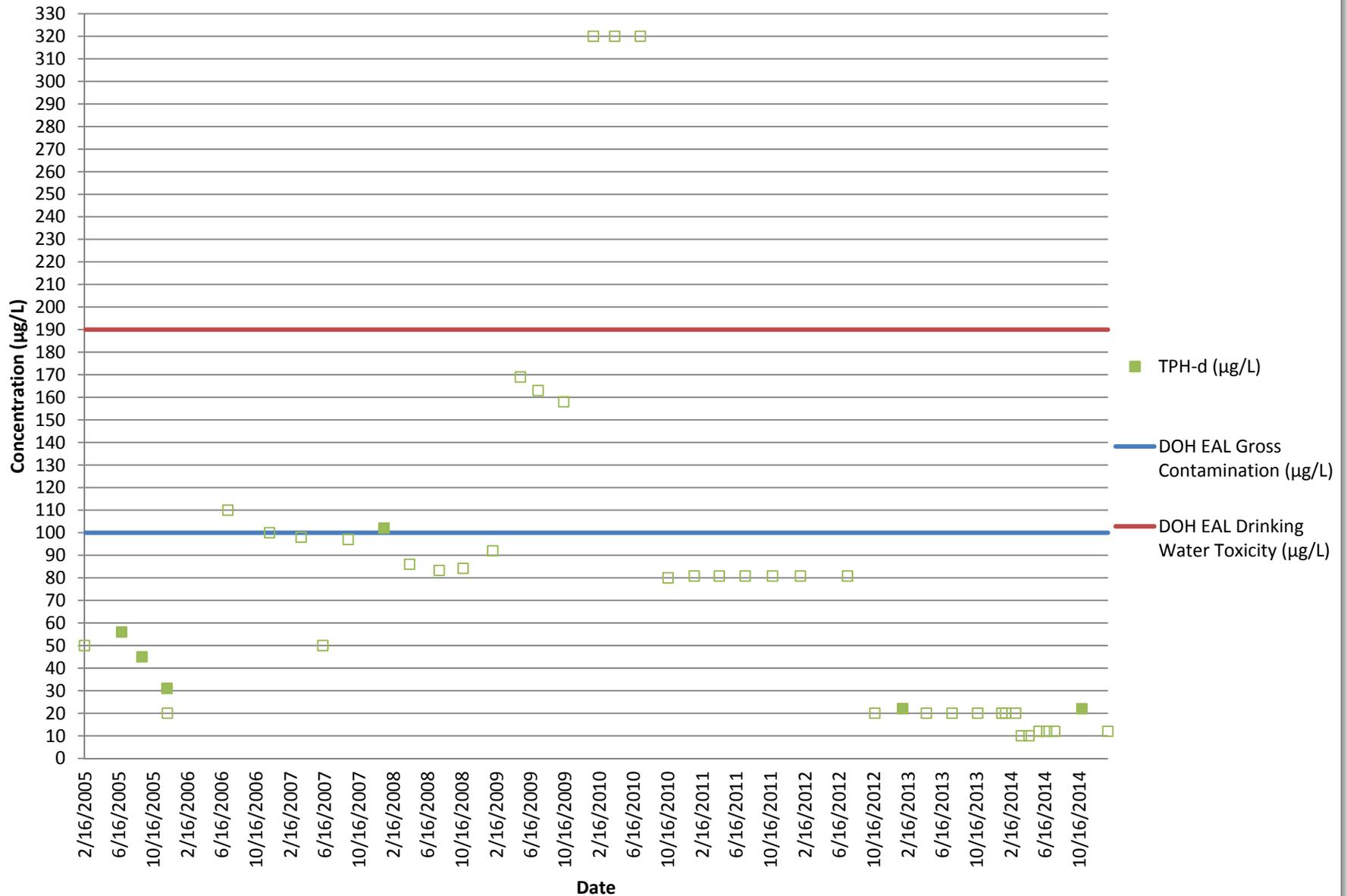
TPH-g Concentrations for RHMW2254-01



Unfilled boxes indicate non-detections. MRLs are shown for February 2005, MDLs are shown for June 2005 through October 2009, and LODs are shown from January 2010 on. Possible laboratory contamination for 10/22/2012, 10/22/2013, and 1/29/2014 sampling events.

This Page Intentionally Left Blank.

TPH-d Concentrations for RHMW2254-01



Unfilled boxes indicate non-detections. MRLs are shown for February 2005, MDLs are shown for December 2005 through October 2009, and LODs are shown from January 2010 on. Laboratory data rejected for 1/15/2008 sampling event.

Numerous sample results had a chromatographic pattern that did not match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.

This Page Intentionally Left Blank.

APPENDIX E

**Cumulative Groundwater Results
(included on attached CD)**

This Page Intentionally Left Blank.