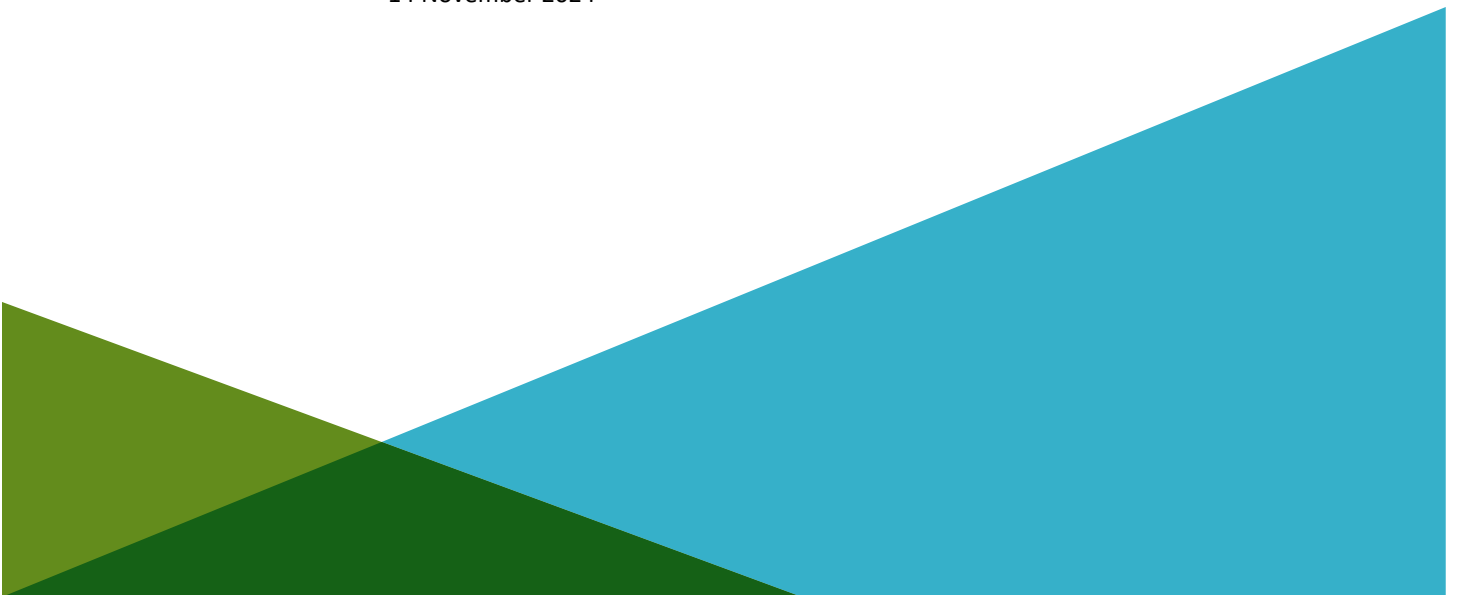


**REPORT ON
CUSTOM PLYWOOD GROUNDWATER MONITORING
ANACORTES, WASHINGTON**

by
Haley & Aldrich, Inc.
Seattle, Washington

for
Washington State Department of Ecology, Toxics Cleanup Program
Shoreline, Washington

File No. 0209325-000
14 November 2024





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14 November 2024
File No. 0209325-000

Washington State Department of Ecology, Toxics Cleanup Program
300 Desmond Drive SE
Lacey, Washington 98503

Attention: Josh Morman
Environmental Specialist, Toxics Cleanup Program

Subject: Custom Plywood Groundwater Monitoring
Anacortes, Washington

Dear Josh:

This groundwater monitoring report provides groundwater quality data on the upland portion of the Custom Plywood site (Site), located in the wetland mitigation area along Fidalgo Bay in Anacortes, Washington (Cleanup Site ID Number 4533, Facility Site ID Number 2685). The project location is shown on Figure 1. This report provides results of the two groundwater sampling events conducted in 2024.

Site Summary

The Site is one of several Anacortes-area, bay-wide priority sites for Fidalgo and Padilla Bays being addressed by the Toxics Cleanup Program (TCP) under the Puget Sound Initiative (PSI). The Site consists of a 6.6-acre upland portion of the larger former Custom Plywood Cleanup. This property is owned by GBH Investments and includes an additional 34 acres of intertidal and subtidal areas.

Custom Plywood operated as a lumber and planing mill between 1900 until burning down some time between 1925 and 1937. Through the mid-1900s, the property changed owners several times, was ultimately rebuilt and expanded until Custom Plywood was again operating some time prior to 1991. The facility was used for sawmill and plywood manufacturing until another fire occurred November 1992 and all operations ceased. Milling activities over the course of its industrial life produced wood-related wastes and chemical contaminants that impacted the soil, sediment, and groundwater.

PREVIOUS AND ONGOING CLEANUP ACTIONS

Past limited interim remedial actions were conducted under Washington Administrative Code 173-340-515 (Independent Remedial Actions) on the upland portion of the site beginning in 1998. To date, three interim remedial actions have been completed at the Site. Phase I of the remedial actions, completed in fall 2011, consisted of upland remediation. Phase II, completed in fall 2013, consisted of shoreline

restoration and cleanup in intertidal and limited subtidal areas. Phase III is currently underway and has consisted of subtidal sediment dredging and capping and eelgrass transplanting and monitoring.

As part of the Phase I Upland Cleanup Action, in September 2011 six groundwater monitoring wells were installed along the northern section of the buffer area and along the western Site boundary to further investigate historic exceedances in diesel- and lube oil-range total petroleum hydrocarbons (TPH), carcinogenic polycyclic aromatic hydrocarbons (cPAHs), and heavy metals associated with petroleum-related contaminants.

Groundwater Sampling and Results

FIELD INVESTIGATION ACTIVITIES AND OBSERVATIONS

The purpose of the groundwater monitoring is to generate data of sufficient quality characterize the nature and extent of potential environmental impacts on the Site as part of compliance monitoring recommendations and the Periodic Review process described in the Final Cleanup Action Plan for Phase I Upland Remediation (September 2011).

It has been several years since any groundwater monitoring has been performed. The objective of the compliance monitoring is to evaluate current shallow groundwater quality at the Site. Prior to mobilizing for the groundwater monitoring, a sampling and analysis plan (SAP) detailing the proposed field efforts, sampling and analysis procedures, and relevant laboratory information was prepared. The SAP include a site-specific Quality Assurance Project Plan (QAPP) and Health and Safety Plan (HASP) for the sampling effort (Haley & Aldrich, 2024).

Due to long period of inactivity at the Site, the monitoring wells were redeveloped over 12 hours prior to sampling by surging. Wells were redeveloped on 31 January and 1 February 2024. Details of the groundwater levels, development water characteristics, and development purge volumes can be found in Attachment A.

There were two sampling events involving the six compliance groundwater monitoring wells (MW-1 through MW-6), approximately six months apart. The two sampling events occurred 1 and 2 February and 12 August 2024. The locations of the wells are shown on Figure 2.

After depressurizing and accurately gauging the depth to groundwater, samples were collected from the wells using standard low-flow sampling techniques. Each well was purged until the field parameters of pH, temperature, and specific conductivity met the stability criteria (i.e., specific conductivity ± 10 percent, pH ± 0.1 pH units, and temperature ± 0.1 °C). Following stabilization, groundwater samples were collected for laboratory testing by directly filling laboratory-provided sample containers. The labeled sample containers were placed in coolers with ice. Samples were delivered under chain of custody protocol to OnSite Environmental, Inc. (OnSite) in Redmond, Washington, for laboratory analysis. During purging, visual and olfactory observations were recorded in addition to quantitative stabilization measurements. For each monitoring events, no visual or olfactory observations were noted that may signify impacted water quality.

GROUNDWATER SAMPLE CHEMICAL ANALYSIS AND RESULTS

Groundwater samples were collected and analyzed from six monitoring wells (MW-1 through MW-6) in each sampling event. The groundwater samples were submitted to OnSite and analyzed for TPH-D, TPH-O, cPAHs, and total and dissolved metals (arsenic, cadmium, chromium, lead, and mercury). Groundwater sample analytical results are summarized in Table 1 and the laboratory reports and data usability summary report (DUSR) are provided in Attachment B.

Cleanup levels are derived from the lowest concentration protective of human or ecological health from Model Toxics Control Act (MTCA) Method B, state surface water quality criteria (Chapter 173-201A WAC), Clean Water Act Section 304, or the National Toxics Rule (40 CFR 131). We compared results with Ecology's Groundwater Preliminary Cleanup Level (PCUL) that's protective of marine surface water, as updated in July 2024 (Table 1). Analytical results are summarized below and on Figure 3.

- TPH-D was detected in samples from MW-2 and MW-3 at concentrations up to 330 micrograms per liter ($\mu\text{g/L}$) between both sampling events, below the Groundwater PCUL of 500 $\mu\text{g/L}$. No other wells had detections of TPH-D at or above the laboratory reporting limit for either sampling event.
- TPH-O was detected in MW-1 and MW-3 in August, and MW-2 for both events at concentrations up to 350 $\mu\text{g/L}$, below the Groundwater PCUL of 500 $\mu\text{g/L}$. No other wells had detections of TPH-O at or above the laboratory reporting limit for either sampling event.
- TPH (the sum of TPH-D and TPH-O) was detected in MW-2 for both events at concentrations up to 680 $\mu\text{g/L}$, exceeding the Groundwater PCUL of 500 $\mu\text{g/L}$.
- cPAHs were not detected at any of the wells at or above laboratory reporting limits.
- Total and dissolved arsenic exceeded the Groundwater PCUL of 8 $\mu\text{g/L}$ in well MW-1 in both events, and in MW-2 and MW-5 in the February event. This is discussed in more detail in the section below.
- Total and dissolved mercury also exceeded the PCUL of 0.025 $\mu\text{g/L}$ in MW-1. This is discussed in more detail in the section below.
- MW-1 in the August sampling event had detections for all total and dissolved metals analyzed, and besides arsenic, were below the pertinent Groundwater PCULs for cadmium, chromium (tri-valent), lead, and mercury.
- Acenaphthene was detected in MW-2 and MW-5 and below the PCUL screening criteria (30 $\mu\text{g/L}$).
- A low-level (0.35 $\mu\text{g/L}$) of 1-methylnaphthalene was detected in MW-5, however there is no applicable PCUL screening criteria.

Conclusions and Recommendations

Dissolved arsenic concentrations are generally within the range of anticipated Puget Sound groundwater background levels except for samples collected from MW-1. A 2022 study by Ecology's Toxics Cleanup Program (TCP) indicates that natural background levels for arsenic ranged from <1 to 150 µg/L across the state. In Island County, an area located near the Site with comparable tidal influences and demands on well water, the calculated arsenic background threshold value is 13.3 µg/L (Ecology, 2022). Mercury also exceeded the PCUL in well MW-1 during August 2024 sampling event.

Diesel- and oil-range hydrocarbons were also detected at concentrations slightly above the PCUL in well MW-2. There was no evidence of petroleum hydrocarbons encountered in the samples (no sheen or odor) and concentrations of PAHs, which are often associated with petroleum products, were not detected at concentrations exceeding PCULs. Review of sample chromatograms and discussion with the laboratory indicates that the detected TPH-D and TPH-O concentrations do not appear to be associated with petroleum products and are likely due to the presence of natural organics. We recommend that groundwater samples analyzed for TPH-D extended in future sampling events should include testing with and without silica gel cleanup.

While it is unlikely that the Site poses a threat to human health and the environment, additional groundwater monitoring may be warranted to assess PCUL exceedances observed in several of the wells. We appreciate the opportunity to provide environmental consulting services on this project. Please do not hesitate to call if you have any questions or comments.

Sincerely yours,
HALEY & ALDRICH, INC.



Samantha Fisher
Senior Environmental Scientist



Andrew Kaparos, P.E.
Senior Associate Engineer



Mike Ehlebracht, L.H.G.
Principal Geochemist

Enclosures:

- Table 1 - Summary of Groundwater Quality Data
- Figure 1 - Vicinity Map
- Figure 2 - Groundwater Monitoring Well Locations
- Figure 3 - Chemical Analysis

Washington State Department of Ecology, Toxics Cleanup Program

14 November 2024

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Attachment A - Completed Field Forms

Attachment B - Laboratory Reports and Data Usability Summary Report

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References

1. San Juan, Charles. 2022. Natural Background Arsenic Concentrations in Washington State. Washington State Department of Ecology. January.
2. Hart Crowser, 2012. Final Construction Completion Report, Phase I, Upland Interim Remedial Action, Custom Plywood Site, Anacortes Washington. Prepared by Hart Crowser under Direction and Contract with the Washington State Department of Ecology under Agreement with GBH Investments, LLC. 22 October.
3. Haley & Aldrich, Inc., 2024. Sampling and Analysis Plan / Quality Assurance Project Plan, Custom Plywood Site, Anacortes, Washington. Prepared by Haley & Aldrich, Inc. for the Washington State Department of Ecology. January.

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TABLE

TABLE 1
SUMMARY OF GROUNDWATER QUALITY DATA
WASHINGTON STATE DEPARTMENT OF ECOLOGY
CUSTOM PLYWOOD
ANACORTES, WASHINGTON

Location Name	MW-1	MW-1	MW-2	MW-2	MW-3	MW-3	MW-4	MW-4	MW-5	MW-5	MW-6	MW-6
Sample Name	MW-1-20240201	MW-1-20240812	MW-2-20240202	MW-2-20240812	MW-3-20240201	MW-3-20240812	MW-4-20240202	MW-4-20240812	MW-5-20240202	MW-5-20240812	MW-6-20240202	MW-6-20240812
Sample Date	02/01/2024	08/12/2024	02/02/2024	08/12/2024	02/01/2024	08/12/2024	02/02/2024	08/12/2024	02/02/2024	08/12/2024	02/02/2024	08/12/2024
Lab Sample ID	2402-047-01	2408-148-05	2402-047-06	2408-148-03	2402-047-02	2408-148-02	2402-047-05	2408-148-01	2402-047-03	2408-148-04	2402-047-04	2408-148-06
Sample Depth (bgs)	Minimum	14.5 (ft)	14 (ft)	-	16 (ft)	14.5 (ft)	14 (ft)	13.5 (ft)	12 (ft)	13 (ft)	12 (ft)	13 (ft)
Total Petroleum Hydrocarbons (ug/L)												
Diesel Range Organics	50/500(a)	220 U	230 U	230	330	250	150 J-	200 U	220 U	210 U	200 U	220 U
PHC as Lube Oil	500	220 U	240	300	350	210 U	320 J-	200 U	220 U	210 U	200 U	220 U
Diesel Range + Oil Range Organics	500	220 U	240	530	680	250	470 J-	200 U	220 U	210 U	200 U	220 U
Inorganic Compounds (ug/L)												
Arsenic, Dissolved	8 (b)	20	27	7.8	3.6	4.4 U	1.4	4.4 U	2.4	7.7	0.67	4.4 U
Cadmium, Dissolved	7.9	4.4 U	0.31	4.4 U	0.22 U	4.4 U	0.22 U	4 U	0.2 U	4 U	0.2 U	4 U
Chromium, Dissolved	27 (c)	11 U	5.9	11 U	8.5	11 U	9.2	10 U	1 U	10 U	1 U	10 U
Lead, Dissolved	8.1	4.4 U	0.87	4.4 U	0.44 U	4.4 U	0.44 U	4 U	0.4 U	4 U	0.4 U	4 U
Mercury, Dissolved	0.025	0.5 U	0.16	0.5 U	0.13 U	0.5 U	0.13 U	0.5 U	0.13 U	0.5 U	0.13 U	0.5 U
Arsenic, Total	8 (b)	21	27	8.1	6.4	4.4 U	2.3	4.4 U	3.2	8.7	0.68	4.4 U
Cadmium, Total	7.9	4.4 U	0.6	4.4 U	0.22 U	4.4 U	0.22 U	4.4 U	0.22 U	4.4 U	0.22 U	4.4 U
Chromium, Total	27 (c)	11 U	12	11 U	12	11 U	9.2	11 U	1.1 U	11 U	1.1 U	11 U
Lead, Total	8.1	4.4 U	3.3	4.4 U	1.3	4.4 U	0.44 U	4.4 U	0.44 U	4.4 U	0.44 U	4.4 U
Mercury, Total	0.025	0.5 U	1	0.5 U	0.13 U	0.5 U	0.13 U	0.5 U	0.13 U	0.5 U	0.13 U	0.5 U
Semi-Volatile Organic Compounds (SIM) (ug/L)												
1-Methylnaphthalene	NA	0.095 U	0.1 U	0.098 U	0.095 U	0.095 U	0.098 U	0.095 U	0.1 U	0.098 U	0.35	0.095 U
2-Methylnaphthalene	NA	0.095 U	0.1 U	0.098 U	0.095 U	0.095 U	0.098 U	0.095 U	0.1 U	0.098 U	0.095 U	0.095 U
Acenaphthene	30	0.095 U	0.1 U	0.14	0.095 U	0.095 U	0.098 U	0.095 U	0.1 U	0.098 U	0.11	0.095 U
Acenaphthylene	NA	0.095 U	0.1 U	0.098 U	0.095 U	0.095 U	0.098 U	0.095 U	0.1 U	0.098 U	0.095 U	0.095 U
Anthracene	100	0.095 U	0.1 U	0.098 U	0.095 U	0.095 U	0.098 U	0.095 U	0.1 U	0.098 U	0.095 U	0.095 U
Benzo(a)anthracene	0.00016	0.0095 U	0.01 U	0.0098 U	0.0095 U	0.0095 U	0.0098 U	0.0095 U	0.01 U	0.0098 U	0.0095 U	0.0095 U
Benzo(a)pyrene	0.000016	0.0095 U	0.01 U	0.0098 U	0.0095 U	0.0095 U	0.0098 U	0.0095 U	0.01 U	0.0098 U	0.0095 U	0.0095 U
Benzo(b)fluoranthene	0.00016	0.0095 U	0.01 U	0.0098 U	0.0095 U	0.0095 U	0.0098 U	0.0095 U	0.01 U	0.0098 U	0.0095 U	0.0095 U
Benzo(g,h,i)perylene	NA	0.0095 U	0.01 U	0.0098 U	0.0095 U	0.0095 U	0.0098 U	0.0095 U	0.012 U	0.0098 U	0.0095 U	0.0095 U
Benzo(j,k)fluoranthene	0.0013	0.0095 U	0.01 U	0.0098 U	0.0095 U	0.0095 U	0.0098 U	0.0095 U	0.01 U	0.0098 U	0.0095 U	0.0095 U
Chrysene	0.016	0.0095 U	0.01 U	0.0098 U	0.0095 U	0.0095 U	0.0098 U	0.0095 U	0.01 U	0.0098 U	0.0095 U	0.0095 U
Dibenz(a,h)anthracene	0.000016	0.0095 U	0.01 U	0.0098 U	0.0095 U	0.0095 U	0.0098 U	0.0095 U	0.01 U	0.0098 U	0.0095 U	0.0095 U
Fluoranthene	6	0.095 U	0.1 U	0.098 U	0.095 U	0.095 U	0.098 U	0.095 U	0.1 U	0.098 U	0.095 U	0.095 U
Fluorene	10	0.095 U	0.1 U	0.098 U	0.095 U	0.095 U	0.098 U	0.095 U	0.1 U	0.098 U	0.095 U	0.095 U
Indeno(1,2,3-cd)pyrene	0.00016	0.0095 U	0.01 U	0.0098 U	0.0095 U	0.0095 U	0.0098 U	0.0095 U	0.014 U	0.0098 U	0.0095 U	0.0095 U
Naphthalene	1.4	0.095 U	0.1 U	0.098 U	0.095 U	0.095 U	0.098 U	0.095 U	0.1 U	0.098 U	0.095 U	0.095 U
Phenanthrene	NA	0.095 U	0.1 U	0.098 U	0.095 U	0.095 U	0.098 U	0.095 U	0.1 U	0.098 U	0.095 U	0.095 U
Pyrene	8	0.095 U	0.1 U	0.098 U	0.095 U	0.095 U	0.098 U	0.095 U	0.1 U	0.098 U	0.095 U	0.095 U
cPAHs-TEQ (d)	0.000016	0.007 U	0.008 U	0.007 U	0.007 U	0.007 U	0.007 U	0.007 U	0.008 U	0.007 U	0.007 U	0.007 U
Field Parameters												
Temperature (Deg C)	NA	9.1	16.1	10.6	14.6	11.3	12.8	9	13.4	10.1	13.2	8.6
Oxidation Reduction Potential (ORP), Field (mv)	NA	-91.6	-114.3	-200.4	-156	-259.4	-201.7	-66.8	-125	-150.7	-	-109.1
Turbidity, Field (NTU)	NA	34.8	38.52	30.57	24.6	40.39	9.87	3.95	2.94	6.31	2.25	2.92
pH, Field (pH units)	NA	7.33	7.57	6.9	6.6	6.91	6.99	7.01	7.22	7.3	-	7.85
Dissolved Oxygen, Field (mg/L)	NA	2	0.44	1.1	0.09	2.4	0	2.7	0.32	2.3	0.05	2.6
Conductivity, Field (uS/cm)	NA	16094	5363	7271	11652	20194	19318	5587	4781	3824	6125	845

ABBREVIATIONS AND NOTES:

--: Not analyzed, probe malfunction 8/12/2024

bgs: below ground surface

Deg C: Degrees Celsius

ft: feet

J-: value is estimated, biased low

mg/L: milligrams per liter

mv: millivolts

NA: Not Available

NTU: Nephelometric Turbidity Units

PCUL: Preliminary Cleanup Level

U: not detected above the indicated laboratory reporting limit

ug/L: micrograms per liter

uS/cm: microSiemen per centimeter

Groundwater PCUL Protect Surface Water Minimum, July 2024

(a) 50 ug/L if fresh, 500 ug/L if weathered

(b) PCUL below background levels. Puget Sound background level of 8 ug/L as determined by Ecology used.

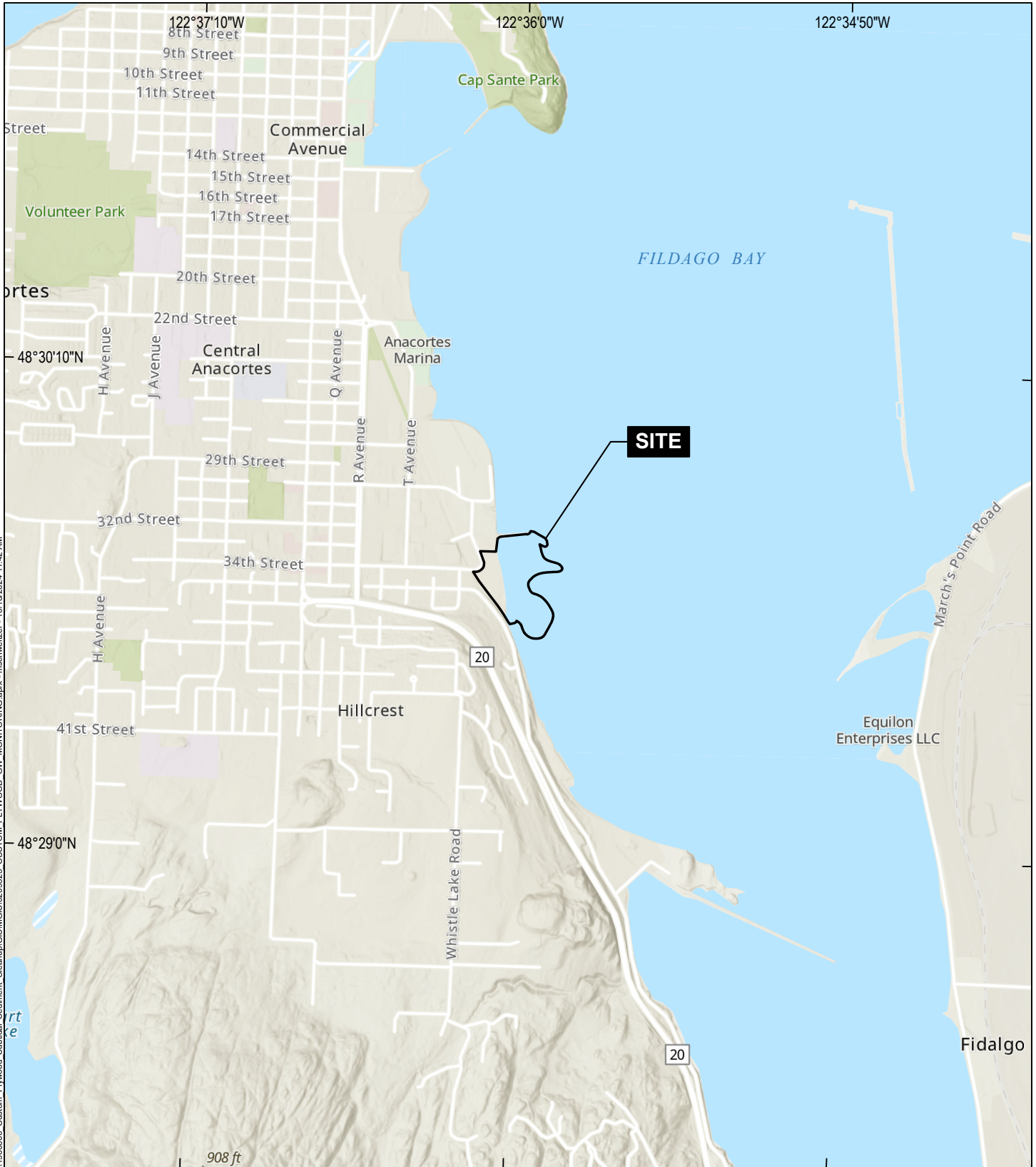
(c) PCUL of 27 ug/L for tri-valent chromium used; no suspected sources of hexavalent chromium at the site.

(d) Toxicity Equivalent using April 2015 (revised July 2021) Ecology's Implementation Memo No. 10: Evaluating the Human Health Toxicity of Carcinogenic PAHs (cPAHs) Using Toxicity Equivalency Factors (TEFs), Publication No. 15-09-049. For non-detects (<RL), a value of one-half the RL has been used for TEQ calculation. cPAH calculated for samples with individual PAHs identified as a non-detect will be qualified with U.

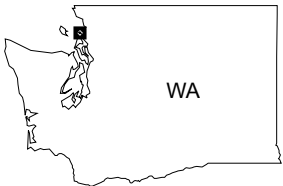
Bold denotes a detected concentration.

Blue shading denotes a detected analyte concentration exceeding a Groundwater PCUL Protect Surface Water Minimum specific to the Lower Duwamish Waterway.

FIGURES



GIS: \\haleyaldrich.com\share\sea_projects\notebooks\1960000_Custom_Plywood_Subtidal_Sediment_Cleanup\GIS\GIS\020925_CUSTOM_PLYWOOD_GW_MONITORING.aprx - mschweitzer - 10/18/2024 11:42 AM



**HALEY
ALDRICH**

CUSTOM PLYWOOD
ANACORTES, WASHINGTON

VICINITY MAP

MAP SOURCE: ESRI
SITE COORDINATES: 48°29'40"N, 122°36'01"W




APPROXIMATE SCALE: 1 IN = 2000 FT
OCTOBER 2024

FIGURE 1

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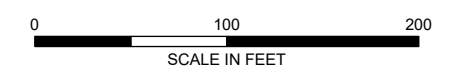


LEGEND

-  GROUNDWATER MONITORING WELL
-  FENCE LINE
-  SITE BOUNDARY

NOTES

1. ALL LOCATIONS AND DIMENSIONS ARE APPROXIMATE.
2. AERIAL IMAGERY SOURCE: NEARMAP, 14 MAY 2024



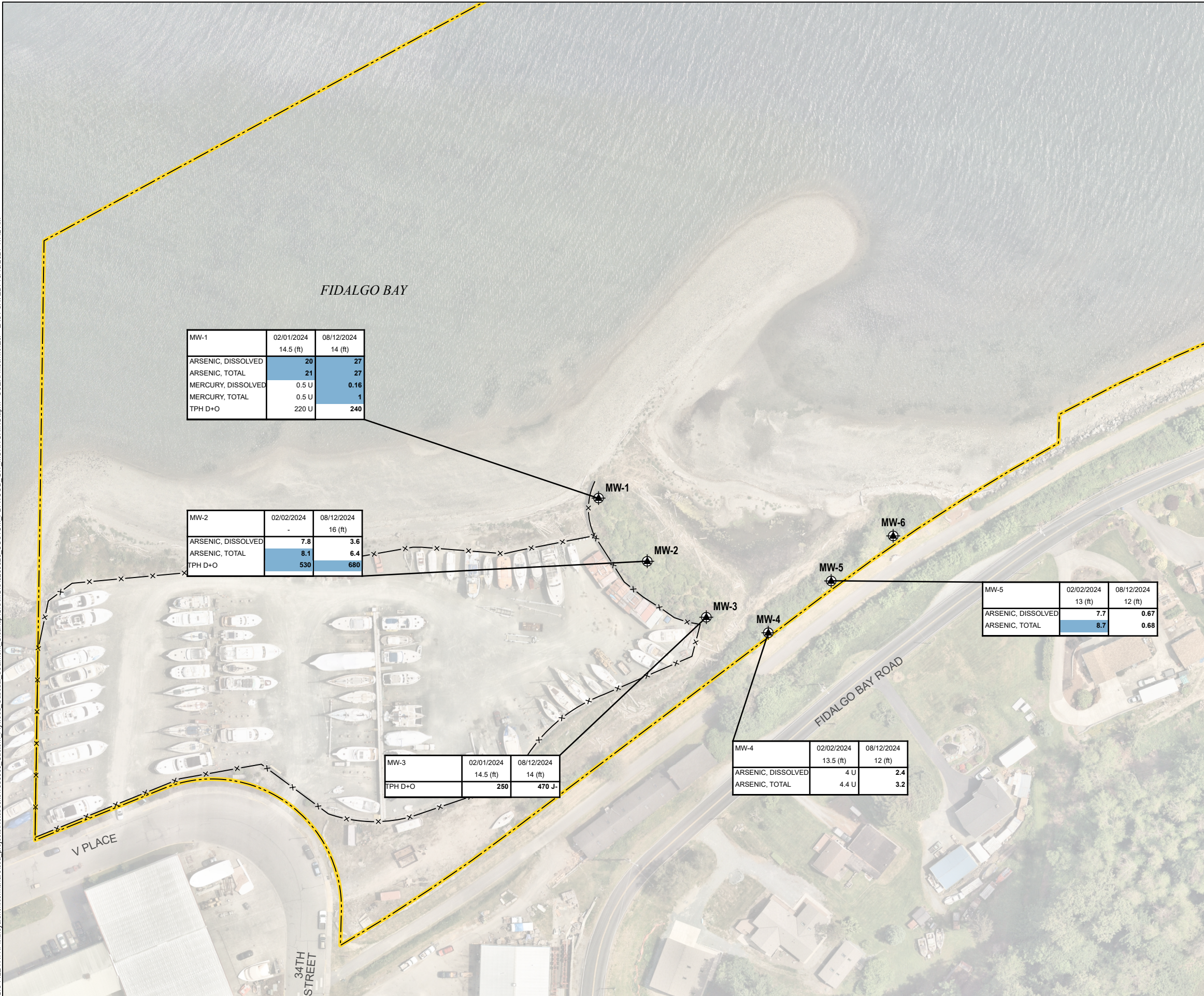
HALEY ALDRICH CUSTOM PLYWOOD ANACORTES, WASHINGTON

GROUNDWATER MONITORING WELL LOCATIONS

OCTOBER 2024

FIGURE 2

GIS FILE PATH: \\haleyaldrich.com\share\sea_projects\Natebooks\1960000_Custom_Plywood_Sublida_Sediment_Cleanup\GIS\MGIS\0209325_CUSTOM_PLYWOOD_GW_MONITORING.aprx - USER: mschwelzer - LAST SAVED: 10/18/2024 11:42 AM



MW-1	02/01/2024 14.5 (ft)	08/12/2024 14 (ft)
ARSENIC, DISSOLVED	20	27
ARSENIC, TOTAL	21	27
MERCURY, DISSOLVED	0.5 U	0.16
MERCURY, TOTAL	0.5 U	1
TPH D+O	220 U	240




MW-2	02/02/2024	08/12/2024 16 (ft)
ARSENIC, DISSOLVED	7.8	3.6
ARSENIC, TOTAL	8.1	6.4
TPH D+O	530	680

MW-3	02/01/2024 14.5 (ft)	08/12/2024 14 (ft)
TPH D+O	250	470 J-

MW-4	02/02/2024 13.5 (ft)	08/12/2024 12 (ft)
ARSENIC, DISSOLVED	4 U	2.4
ARSENIC, TOTAL	4.4 U	3.2

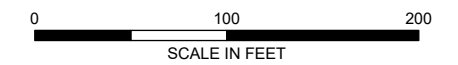
MW-5	02/02/2024 13 (ft)	08/12/2024 12 (ft)
ARSENIC, DISSOLVED	7.7	0.67
ARSENIC, TOTAL	8.7	0.68

LEGEND

-  GROUNDWATER MONITORING WELL
-  FENCE LINE
-  SITE BOUNDARY

NOTES

1. ALL LOCATIONS AND DIMENSIONS ARE APPROXIMATE.
2. CONCENTRATIONS ARE IN MICROGRAMS PER LITER (ug/L)
3. DEFINITIONS:
 FT = FEET
 J- = RESULT IS AN ESTIMATE
 J- = RESULT IS AN ESTIMATE, BIASED LOW
 TPH D+O = SUM OF DIESEL- AND OIL-RANGE TOTAL PETROLEUM HYDROCARBONS
 U = NOT DETECTED ABOVE THE INDICATED LABORATORY REPORTING LIMIT
4. VALUES IN **BOLD** DENOTES A DETECTED CONCENTRATION.
5. VALUES WITH **BLUE SHADING** DENOTES A DETECTED ANALYTE CONCENTRATION EXCEEDING A GROUNDWATER PCUL PROTECT SURFACE WATER MINIMUM SPECIFIC TO THE LOWER DUWAMISH WATERWAY, ECOLOGY (JULY 2024).
6. AERIAL IMAGERY SOURCE: NEARMAP, 14 MAY 2024



HALEY ALDRICH CUSTOM PLYWOOD ANACORTES, WASHINGTON

CHEMICAL ANALYSIS

OCTOBER 2024

FIGURE 3

APPENDIX A
Completed Field Forms

APPENDIX B
Laboratory Reports and Data Usability Summary Report



14648 NE 95th Street, Redmond, WA 98052 • (425) 883-3881

March 5, 2024

Andrew Kaparos
Hart Crowser, Inc.
A Division of Haley & Aldrich, Inc.
3131 Elliott Avenue, Suite 600
Seattle, WA 98121

Re: Analytical Data for Project 0209325-000
Laboratory Reference No. 2402-047

Dear Andrew:

Enclosed are the analytical results and associated quality control data for samples submitted on February 7, 2024.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal flourish extending to the right.

David Baumeister
Project Manager

Enclosures



OnSite Environmental, Inc. 14648 NE 95th Street, Redmond, WA 98052 (425) 883-3881

This report pertains to the samples analyzed in accordance with the chain of custody, and is intended only for the use of the individual or company to whom it is addressed.

Date of Report: March 5, 2024
Samples Submitted: February 7, 2024
Laboratory Reference: 2402-047
Project: 0209325-000

Case Narrative

Samples were collected on February 1 and 2, 2024 and received by the laboratory on February 7, 2024. They were maintained at the laboratory at a temperature of 2°C to 6°C.

Please note that any and all soil sample results are reported on a dry-weight basis, unless otherwise noted below. However the soil results for the QA/QC samples are reported on a wet-weight basis.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

NWTPH-Dx Analysis

The surrogate percent recovery in sample MW-3 was below the control limit of 50% due to matrix effects. The sample was re-extracted with the same result. The percent recovery for the silica gel cleanup for this sample was within control limits due to the natural tendency for extracts to concentrate slightly during the cleanup process.

Dissolved Metals by EPA 200.8/7470A Analysis

The dissolved field filtered samples MW-1, MW-3 and MW-2 were received containing solid material. The samples were digested according to the OnSite Environmental standard operating procedure.

Please note that any other QA/QC issues associated with these extractions and analyses will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.



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**DIESEL AND HEAVY OIL RANGE ORGANICS
 NWTPH-Dx**

Matrix: Water
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-1					
Laboratory ID:	02-047-01					
Diesel Range Organics	ND	0.22	NWTPH-Dx	2-6-24	2-6-24	
Lube Oil Range Organics	ND	0.22	NWTPH-Dx	2-6-24	2-6-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	69	50-150				

Client ID:	MW-1					
Laboratory ID:	02-047-01					
Diesel Range Organics	ND	0.22	NWTPH-Dx	2-6-24	2-6-24	X2
Lube Oil Range Organics	ND	0.22	NWTPH-Dx	2-6-24	2-6-24	X2
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	70	50-150				

Client ID:	MW-3					
Laboratory ID:	02-047-02					
Diesel Range Organics	0.25	0.21	NWTPH-Dx	2-6-24	2-7-24	
Lube Oil Range Organics	ND	0.21	NWTPH-Dx	2-6-24	2-7-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	50	50-150				Q

Client ID:	MW-3					
Laboratory ID:	02-047-02					
Diesel Range Organics	ND	0.21	NWTPH-Dx	2-6-24	2-7-24	X2
Lube Oil Range Organics	ND	0.21	NWTPH-Dx	2-6-24	2-7-24	X2
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	52	50-150				

Client ID:	MW-5					
Laboratory ID:	02-047-03					
Diesel Range Organics	ND	0.21	NWTPH-Dx	2-5-24	2-5-24	
Lube Oil Range Organics	ND	0.21	NWTPH-Dx	2-5-24	2-5-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	70	50-150				

Client ID:	MW-5					
Laboratory ID:	02-047-03					
Diesel Range Organics	ND	0.21	NWTPH-Dx	2-5-24	2-5-24	X2
Lube Oil Range Organics	ND	0.21	NWTPH-Dx	2-5-24	2-5-24	X2
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	78	50-150				



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**DIESEL AND HEAVY OIL RANGE ORGANICS
 NWTPH-Dx**

Matrix: Water
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-6					
Laboratory ID:	02-047-04					
Diesel Range Organics	ND	0.20	NWTPH-Dx	2-5-24	2-5-24	
Lube Oil Range Organics	ND	0.20	NWTPH-Dx	2-5-24	2-5-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	68	50-150				

Client ID:	MW-6					
Laboratory ID:	02-047-04					
Diesel Range Organics	ND	0.20	NWTPH-Dx	2-5-24	2-5-24	X2
Lube Oil Range Organics	ND	0.20	NWTPH-Dx	2-5-24	2-5-24	X2
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	67	50-150				

Client ID:	MW-4					
Laboratory ID:	02-047-05					
Diesel Range Organics	ND	0.20	NWTPH-Dx	2-5-24	2-5-24	
Lube Oil Range Organics	ND	0.20	NWTPH-Dx	2-5-24	2-5-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	91	50-150				

Client ID:	MW-4					
Laboratory ID:	02-047-05					
Diesel Range Organics	ND	0.20	NWTPH-Dx	2-5-24	2-5-24	X2
Lube Oil Range Organics	ND	0.20	NWTPH-Dx	2-5-24	2-5-24	X2
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	95	50-150				

Client ID:	MW-2					
Laboratory ID:	02-047-06					
Diesel Range Organics	0.23	0.23	NWTPH-Dx	2-6-24	2-6-24	
Lube Oil Range Organics	0.30	0.23	NWTPH-Dx	2-6-24	2-6-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	52	50-150				

Client ID:	MW-2					
Laboratory ID:	02-047-06					
Diesel Range Organics	ND	0.23	NWTPH-Dx	2-6-24	2-6-24	X2
Lube Oil Range Organics	ND	0.23	NWTPH-Dx	2-6-24	2-6-24	X2
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	55	50-150				



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**DIESEL AND HEAVY OIL RANGE ORGANICS
 NWTPH-Dx
 QUALITY CONTROL**

Matrix: Water
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0205W1					
Diesel Range Organics	ND	0.16	NWTPH-Dx	2-5-24	2-5-24	
Lube Oil Range Organics	ND	0.16	NWTPH-Dx	2-5-24	2-5-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	94	50-150				
Laboratory ID:	MB0205W1					
Diesel Range Organics	ND	0.16	NWTPH-Dx	2-5-24	2-5-24	X2
Lube Oil Range Organics	ND	0.16	NWTPH-Dx	2-5-24	2-5-24	X2
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	117	50-150				
Laboratory ID:	MB0206W1					
Diesel Range Organics	ND	0.16	NWTPH-Dx	2-6-24	2-6-24	
Lube Oil Range Organics	ND	0.16	NWTPH-Dx	2-6-24	2-6-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	88	50-150				
Laboratory ID:	MB0206W1					
Diesel Range Organics	ND	0.16	NWTPH-Dx	2-6-24	2-6-24	X2
Lube Oil Range Organics	ND	0.16	NWTPH-Dx	2-6-24	2-6-24	X2
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	99	50-150				



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**DIESEL AND HEAVY OIL RANGE ORGANICS
 NWTPH-Dx
 QUALITY CONTROL**

Matrix: Water
 Units: mg/L (ppm)

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
DUPLICATE								
Laboratory ID:	SB0205W1							
	ORIG	DUP						
Diesel Fuel #2	0.398	0.387	NA	NA	NA	NA	3	40
Surrogate:								
<i>o</i> -Terphenyl				95	91	50-150		
Laboratory ID:	SB0205W1							
	ORIG	DUP						
Diesel Fuel #2	0.401	0.373	NA	NA	NA	NA	7	40
Surrogate:								
<i>o</i> -Terphenyl				99	92	50-150		
Laboratory ID:	SB0206W1							
	ORIG	DUP						
Diesel Fuel #2	0.405	0.373	NA	NA	NA	NA	8	40
Surrogate:								
<i>o</i> -Terphenyl				101	100	50-150		
Laboratory ID:	SB0206W1							
	ORIG	DUP						
Diesel Fuel #2	0.474	0.366	NA	NA	NA	NA	26	40
Surrogate:								
<i>o</i> -Terphenyl				116	100	50-150		



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**DIESEL AND HEAVY OIL RANGE ORGANICS
 NWTPH-Dx
 QUALITY CONTROL**

Matrix: Water
 Units: mg/L (ppm)

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
SPIKE BLANK								
Laboratory ID:	SB0205W1							
Diesel Fuel #2	0.398	0.500	NA	80	53-126	NA	NA	
Surrogate: o-Terphenyl				95	50-150			
Laboratory ID:	SB0205W1 SG							
Diesel Fuel #2	0.401	0.500	NA	80	53-126	NA	NA	X2
Surrogate: o-Terphenyl				99	50-150			
Laboratory ID:	SB0206W1							
Diesel Fuel #2	0.405	0.500	NA	81	53-126	NA	NA	
Surrogate: o-Terphenyl				101	50-150			
Laboratory ID:	SB0206W1 SG							
Diesel Fuel #2	0.474	0.500	NA	95	53-126	NA	NA	X2
Surrogate: o-Terphenyl				116	50-150			



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PAHs EPA 8270E/SIM

Matrix: Water
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-1					
Laboratory ID:	02-047-01					
Naphthalene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
2-Methylnaphthalene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
1-Methylnaphthalene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
Acenaphthylene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
Acenaphthene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
Fluorene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
Phenanthrene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
Anthracene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
Fluoranthene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
Pyrene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[a]anthracene	ND	0.0095	EPA 8270E/SIM	2-5-24	2-6-24	
Chrysene	ND	0.0095	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[b]fluoranthene	ND	0.0095	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo(j,k)fluoranthene	ND	0.0095	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[a]pyrene	ND	0.0095	EPA 8270E/SIM	2-5-24	2-6-24	
Indeno(1,2,3-c,d)pyrene	ND	0.0095	EPA 8270E/SIM	2-5-24	2-6-24	
Dibenz[a,h]anthracene	ND	0.0095	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[g,h,i]perylene	ND	0.0095	EPA 8270E/SIM	2-5-24	2-6-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>55</i>	<i>26-106</i>				
<i>Pyrene-d10</i>	<i>72</i>	<i>45-104</i>				
<i>Terphenyl-d14</i>	<i>93</i>	<i>43-114</i>				



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PAHs EPA 8270E/SIM

Matrix: Water
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-3					
Laboratory ID:	02-047-02					
Naphthalene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
2-Methylnaphthalene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
1-Methylnaphthalene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
Acenaphthylene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
Acenaphthene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
Fluorene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
Phenanthrene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
Anthracene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
Fluoranthene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
Pyrene	ND	0.095	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[a]anthracene	ND	0.0095	EPA 8270E/SIM	2-5-24	2-6-24	
Chrysene	ND	0.0095	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[b]fluoranthene	ND	0.0095	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo(j,k)fluoranthene	ND	0.0095	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[a]pyrene	ND	0.0095	EPA 8270E/SIM	2-5-24	2-6-24	
Indeno(1,2,3-c,d)pyrene	ND	0.0095	EPA 8270E/SIM	2-5-24	2-6-24	
Dibenz[a,h]anthracene	ND	0.0095	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[g,h,i]perylene	ND	0.0095	EPA 8270E/SIM	2-5-24	2-6-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>49</i>	<i>26-106</i>				
<i>Pyrene-d10</i>	<i>52</i>	<i>45-104</i>				
<i>Terphenyl-d14</i>	<i>65</i>	<i>43-114</i>				



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PAHs EPA 8270E/SIM

Matrix: Water
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-5					
Laboratory ID:	02-047-03					
Naphthalene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
2-Methylnaphthalene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
1-Methylnaphthalene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
Acenaphthylene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
Acenaphthene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
Fluorene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
Phenanthrene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
Anthracene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
Fluoranthene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
Pyrene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[a]anthracene	ND	0.0098	EPA 8270E/SIM	2-5-24	2-6-24	
Chrysene	ND	0.0098	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[b]fluoranthene	ND	0.0098	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo(j,k)fluoranthene	ND	0.0098	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[a]pyrene	ND	0.0098	EPA 8270E/SIM	2-5-24	2-6-24	
Indeno(1,2,3-c,d)pyrene	ND	0.0098	EPA 8270E/SIM	2-5-24	2-6-24	
Dibenz[a,h]anthracene	ND	0.0098	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[g,h,i]perylene	ND	0.0098	EPA 8270E/SIM	2-5-24	2-6-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>44</i>	<i>26-106</i>				
<i>Pyrene-d10</i>	<i>61</i>	<i>45-104</i>				
<i>Terphenyl-d14</i>	<i>70</i>	<i>43-114</i>				



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PAHs EPA 8270E/SIM

Matrix: Water
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-6					
Laboratory ID:	02-047-04					
Naphthalene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
2-Methylnaphthalene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
1-Methylnaphthalene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
Acenaphthylene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
Acenaphthene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
Fluorene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
Phenanthrene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
Anthracene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
Fluoranthene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
Pyrene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
Benzo[a]anthracene	ND	0.0095	EPA 8270E/SIM	2-7-24	2-7-24	
Chrysene	ND	0.0095	EPA 8270E/SIM	2-7-24	2-7-24	
Benzo[b]fluoranthene	ND	0.0095	EPA 8270E/SIM	2-7-24	2-7-24	
Benzo(j,k)fluoranthene	ND	0.0095	EPA 8270E/SIM	2-7-24	2-7-24	
Benzo[a]pyrene	ND	0.0095	EPA 8270E/SIM	2-7-24	2-7-24	
Indeno(1,2,3-c,d)pyrene	ND	0.0095	EPA 8270E/SIM	2-7-24	2-7-24	
Dibenz[a,h]anthracene	ND	0.0095	EPA 8270E/SIM	2-7-24	2-7-24	
Benzo[g,h,i]perylene	ND	0.0095	EPA 8270E/SIM	2-7-24	2-7-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>56</i>	<i>26-106</i>				
<i>Pyrene-d10</i>	<i>60</i>	<i>45-104</i>				
<i>Terphenyl-d14</i>	<i>63</i>	<i>43-114</i>				



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PAHs EPA 8270E/SIM

Matrix: Water
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-4					
Laboratory ID:	02-047-05					
Naphthalene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
2-Methylnaphthalene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
1-Methylnaphthalene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
Acenaphthylene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
Acenaphthene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
Fluorene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
Phenanthrene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
Anthracene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
Fluoranthene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
Pyrene	ND	0.095	EPA 8270E/SIM	2-7-24	2-7-24	
Benzo[a]anthracene	ND	0.0095	EPA 8270E/SIM	2-7-24	2-7-24	
Chrysene	ND	0.0095	EPA 8270E/SIM	2-7-24	2-7-24	
Benzo[b]fluoranthene	ND	0.0095	EPA 8270E/SIM	2-7-24	2-7-24	
Benzo(j,k)fluoranthene	ND	0.0095	EPA 8270E/SIM	2-7-24	2-7-24	
Benzo[a]pyrene	ND	0.0095	EPA 8270E/SIM	2-7-24	2-7-24	
Indeno(1,2,3-c,d)pyrene	ND	0.0095	EPA 8270E/SIM	2-7-24	2-7-24	
Dibenz[a,h]anthracene	ND	0.0095	EPA 8270E/SIM	2-7-24	2-7-24	
Benzo[g,h,i]perylene	ND	0.0095	EPA 8270E/SIM	2-7-24	2-7-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>67</i>	<i>26-106</i>				
<i>Pyrene-d10</i>	<i>75</i>	<i>45-104</i>				
<i>Terphenyl-d14</i>	<i>83</i>	<i>43-114</i>				



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PAHs EPA 8270E/SIM

Matrix: Water
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-2					
Laboratory ID:	02-047-06					
Naphthalene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
2-Methylnaphthalene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
1-Methylnaphthalene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
Acenaphthylene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
Acenaphthene	0.14	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
Fluorene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
Phenanthrene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
Anthracene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
Fluoranthene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
Pyrene	ND	0.098	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[a]anthracene	ND	0.0098	EPA 8270E/SIM	2-5-24	2-6-24	
Chrysene	ND	0.0098	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[b]fluoranthene	ND	0.0098	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo(j,k)fluoranthene	ND	0.0098	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[a]pyrene	ND	0.0098	EPA 8270E/SIM	2-5-24	2-6-24	
Indeno(1,2,3-c,d)pyrene	ND	0.0098	EPA 8270E/SIM	2-5-24	2-6-24	
Dibenz[a,h]anthracene	ND	0.0098	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[g,h,i]perylene	ND	0.0098	EPA 8270E/SIM	2-5-24	2-6-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>52</i>	<i>26-106</i>				
<i>Pyrene-d10</i>	<i>71</i>	<i>45-104</i>				
<i>Terphenyl-d14</i>	<i>92</i>	<i>43-114</i>				



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**PAHs EPA 8270E/SIM
 QUALITY CONTROL**

Matrix: Water
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0205W1					
Naphthalene	ND	0.10	EPA 8270E/SIM	2-5-24	2-6-24	
2-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	2-5-24	2-6-24	
1-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	2-5-24	2-6-24	
Acenaphthylene	ND	0.10	EPA 8270E/SIM	2-5-24	2-6-24	
Acenaphthene	ND	0.10	EPA 8270E/SIM	2-5-24	2-6-24	
Fluorene	ND	0.10	EPA 8270E/SIM	2-5-24	2-6-24	
Phenanthrene	ND	0.10	EPA 8270E/SIM	2-5-24	2-6-24	
Anthracene	ND	0.10	EPA 8270E/SIM	2-5-24	2-6-24	
Fluoranthene	ND	0.10	EPA 8270E/SIM	2-5-24	2-6-24	
Pyrene	ND	0.10	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[a]anthracene	ND	0.010	EPA 8270E/SIM	2-5-24	2-6-24	
Chrysene	ND	0.010	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[b]fluoranthene	ND	0.010	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[a]pyrene	ND	0.010	EPA 8270E/SIM	2-5-24	2-6-24	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270E/SIM	2-5-24	2-6-24	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270E/SIM	2-5-24	2-6-24	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270E/SIM	2-5-24	2-6-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>63</i>	<i>26-106</i>				
<i>Pyrene-d10</i>	<i>75</i>	<i>45-104</i>				
<i>Terphenyl-d14</i>	<i>74</i>	<i>43-114</i>				



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**PAHs EPA 8270E/SIM
 QUALITY CONTROL**

Matrix: Water
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0207W1					
Naphthalene	ND	0.10	EPA 8270E/SIM	2-7-24	2-7-24	
2-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	2-7-24	2-7-24	
1-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	2-7-24	2-7-24	
Acenaphthylene	ND	0.10	EPA 8270E/SIM	2-7-24	2-7-24	
Acenaphthene	ND	0.10	EPA 8270E/SIM	2-7-24	2-7-24	
Fluorene	ND	0.10	EPA 8270E/SIM	2-7-24	2-7-24	
Phenanthrene	ND	0.10	EPA 8270E/SIM	2-7-24	2-7-24	
Anthracene	ND	0.10	EPA 8270E/SIM	2-7-24	2-7-24	
Fluoranthene	ND	0.10	EPA 8270E/SIM	2-7-24	2-7-24	
Pyrene	ND	0.10	EPA 8270E/SIM	2-7-24	2-7-24	
Benzo[a]anthracene	ND	0.010	EPA 8270E/SIM	2-7-24	2-7-24	
Chrysene	ND	0.010	EPA 8270E/SIM	2-7-24	2-7-24	
Benzo[b]fluoranthene	ND	0.010	EPA 8270E/SIM	2-7-24	2-7-24	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270E/SIM	2-7-24	2-7-24	
Benzo[a]pyrene	ND	0.010	EPA 8270E/SIM	2-7-24	2-7-24	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270E/SIM	2-7-24	2-7-24	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270E/SIM	2-7-24	2-7-24	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270E/SIM	2-7-24	2-7-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>51</i>	<i>26-106</i>				
<i>Pyrene-d10</i>	<i>66</i>	<i>45-104</i>				
<i>Terphenyl-d14</i>	<i>65</i>	<i>43-114</i>				



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**PAHs EPA 8270E/SIM
 QUALITY CONTROL**

Matrix: Water
 Units: ug/L

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD	RPD	Flags
					Recovery	Limits	Limit			
SPIKE BLANKS										
Laboratory ID:	SB0205W1									
	SB	SBD	SB	SBD	SB	SBD				
Naphthalene	0.268	0.304	0.500	0.500	54	61	35 - 84	13	34	
Acenaphthylene	0.297	0.352	0.500	0.500	59	70	44 - 97	17	29	
Acenaphthene	0.274	0.327	0.500	0.500	55	65	40 - 93	18	29	
Fluorene	0.279	0.333	0.500	0.500	56	67	46 - 97	18	24	
Phenanthrene	0.295	0.342	0.500	0.500	59	68	49 - 102	15	21	
Anthracene	0.296	0.355	0.500	0.500	59	71	50 - 99	18	21	
Fluoranthene	0.312	0.345	0.500	0.500	62	69	53 - 107	10	21	
Pyrene	0.317	0.376	0.500	0.500	63	75	52 - 111	17	23	
Benzo[a]anthracene	0.320	0.387	0.500	0.500	64	77	51 - 119	19	20	
Chrysene	0.311	0.358	0.500	0.500	62	72	52 - 113	14	21	
Benzo[b]fluoranthene	0.387	0.459	0.500	0.500	77	92	50 - 116	17	24	
Benzo(j,k)fluoranthene	0.292	0.345	0.500	0.500	58	69	54 - 113	17	22	
Benzo[a]pyrene	0.339	0.399	0.500	0.500	68	80	52 - 110	16	21	
Indeno(1,2,3-c,d)pyrene	0.373	0.423	0.500	0.500	75	85	55 - 114	13	21	
Dibenz[a,h]anthracene	0.340	0.405	0.500	0.500	68	81	55 - 111	17	19	
Benzo[g,h,i]perylene	0.334	0.395	0.500	0.500	67	79	52 - 111	17	20	
<i>Surrogate:</i>										
2-Fluorobiphenyl					58	68	26-106			
Pyrene-d10					70	79	45-104			
Terphenyl-d14					69	77	43-114			



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**PAHs EPA 8270E/SIM
 QUALITY CONTROL**

Matrix: Water
 Units: ug/L

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD	RPD	Flags
					Recovery	Limits	Limit			
SPIKE BLANKS										
Laboratory ID:	SB0207W1									
	SB	SBD	SB	SBD	SB	SBD				
Naphthalene	0.343	0.358	0.500	0.500	69	72	35 - 84	4	34	
Acenaphthylene	0.414	0.375	0.500	0.500	83	75	44 - 97	10	29	
Acenaphthene	0.368	0.339	0.500	0.500	74	68	40 - 93	8	29	
Fluorene	0.393	0.365	0.500	0.500	79	73	46 - 97	7	24	
Phenanthrene	0.385	0.374	0.500	0.500	77	75	49 - 102	3	21	
Anthracene	0.414	0.391	0.500	0.500	83	78	50 - 99	6	21	
Fluoranthene	0.464	0.437	0.500	0.500	93	87	53 - 107	6	21	
Pyrene	0.442	0.436	0.500	0.500	88	87	52 - 111	1	23	
Benzo[a]anthracene	0.480	0.438	0.500	0.500	96	88	51 - 119	9	20	
Chrysene	0.479	0.436	0.500	0.500	96	87	52 - 113	9	21	
Benzo[b]fluoranthene	0.527	0.445	0.500	0.500	105	89	50 - 116	17	24	
Benzo(j,k)fluoranthene	0.484	0.460	0.500	0.500	97	92	54 - 113	5	22	
Benzo[a]pyrene	0.513	0.453	0.500	0.500	103	91	52 - 110	12	21	
Indeno(1,2,3-c,d)pyrene	0.540	0.464	0.500	0.500	108	93	55 - 114	15	21	
Dibenz[a,h]anthracene	0.514	0.460	0.500	0.500	103	92	55 - 111	11	19	
Benzo[g,h,i]perylene	0.510	0.456	0.500	0.500	102	91	52 - 111	11	20	
<i>Surrogate:</i>										
2-Fluorobiphenyl					83	73	26-106			
Pyrene-d10					90	85	45-104			
Terphenyl-d14					90	85	43-114			



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TOTAL METALS
EPA 200.8/7470A

Matrix: Water
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-1					
Laboratory ID:	02-047-01					
Arsenic	21	4.4	EPA 200.8	2-26-24	3-5-24	
Cadmium	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Chromium	ND	11	EPA 200.8	2-26-24	3-5-24	
Lead	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Mercury	ND	0.50	EPA 7470A	2-15-24	2-15-24	

Client ID:	MW-3					
Laboratory ID:	02-047-02					
Arsenic	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Cadmium	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Chromium	ND	11	EPA 200.8	2-26-24	3-5-24	
Lead	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Mercury	ND	0.50	EPA 7470A	2-15-24	2-15-24	

Client ID:	MW-5					
Laboratory ID:	02-047-03					
Arsenic	8.7	4.4	EPA 200.8	2-26-24	3-5-24	
Cadmium	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Chromium	ND	11	EPA 200.8	2-26-24	3-5-24	
Lead	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Mercury	ND	0.50	EPA 7470A	2-15-24	2-15-24	

Client ID:	MW-6					
Laboratory ID:	02-047-04					
Arsenic	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Cadmium	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Chromium	ND	11	EPA 200.8	2-26-24	3-5-24	
Lead	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Mercury	ND	0.50	EPA 7470A	2-15-24	2-15-24	



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TOTAL METALS
EPA 200.8/7470A

Matrix: Water
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-4					
Laboratory ID:	02-047-05					
Arsenic	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Cadmium	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Chromium	ND	11	EPA 200.8	2-26-24	3-5-24	
Lead	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Mercury	ND	0.50	EPA 7470A	2-15-24	2-15-24	

Client ID:	MW-2					
Laboratory ID:	02-047-06					
Arsenic	8.1	4.4	EPA 200.8	2-26-24	3-5-24	
Cadmium	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Chromium	ND	11	EPA 200.8	2-26-24	3-5-24	
Lead	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Mercury	ND	0.50	EPA 7470A	2-15-24	2-15-24	



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**TOTAL METALS
 EPA 200.8/7470A
 QUALITY CONTROL**

Matrix: Water
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0226WM1					
Arsenic	ND	3.3	EPA 200.8	2-26-24	2-26-24	
Cadmium	ND	4.4	EPA 200.8	2-26-24	2-26-24	
Chromium	ND	11	EPA 200.8	2-26-24	2-26-24	
Lead	ND	1.1	EPA 200.8	2-26-24	2-26-24	

Laboratory ID:	MB0215W1					
Mercury	ND	0.50	EPA 7470A	2-15-24	2-15-24	

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
DUPLICATE								
Laboratory ID:	01-067-08							
	ORIG	DUP						
Arsenic	ND	ND	NA	NA	NA	NA	NA	20
Cadmium	ND	ND	NA	NA	NA	NA	NA	20
Chromium	ND	ND	NA	NA	NA	NA	NA	20
Lead	ND	ND	NA	NA	NA	NA	NA	20

Laboratory ID:	02-186-01							
Mercury	ND	ND	NA	NA	NA	NA	NA	20



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**TOTAL METALS
 EPA 200.8/7470A
 QUALITY CONTROL**

Matrix: Water
 Units: ug/L (ppb)

Analyte	Result		Spike Level		Source	Percent	Recovery	RPD		Flags
					Result	Recovery	Limits	RPD	Limit	
MATRIX SPIKES										
Laboratory ID:	01-067-08									
	MS	MSD	MS	MSD		MS	MSD			
Arsenic	113	112	111	111	ND	102	101	75-125	1	20
Cadmium	111	109	111	111	ND	100	98	75-125	2	20
Chromium	112	111	111	111	ND	101	100	75-125	1	20
Lead	109	109	111	111	ND	99	98	75-125	0	20
SPIKE BLANK										
Laboratory ID:	02-186-01									
Mercury	11.9	12.1	12.5	12.5	ND	95	97	75-125	1	20
Laboratory ID:	SB0226WM1									
Arsenic	112		111		N/A	101		85-115		
Cadmium	111		111		N/A	100		85-115		
Chromium	110		111		N/A	99		85-115		
Lead	112		111		N/A	101		85-115		
Laboratory ID:	SB0221W1									
Mercury	11.9		12.5		N/A	95		80-120		



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DISSOLVED METALS
EPA 200.8/7470A

Matrix: Water
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-1					
Laboratory ID:	02-047-01					
Arsenic	20	4.4	EPA 200.8	2-26-24	3-5-24	
Cadmium	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Chromium	ND	11	EPA 200.8	2-26-24	3-5-24	
Lead	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Mercury	ND	0.50	EPA 7470A		2-15-24	

Client ID:	MW-3					
Laboratory ID:	02-047-02					
Arsenic	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Cadmium	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Chromium	ND	11	EPA 200.8	2-26-24	3-5-24	
Lead	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Mercury	ND	0.50	EPA 7470A		2-15-24	

Client ID:	MW-5					
Laboratory ID:	02-047-03					
Arsenic	7.7	4.0	EPA 200.8		3-5-24	
Cadmium	ND	4.0	EPA 200.8		3-5-24	
Chromium	ND	10	EPA 200.8		3-5-24	
Lead	ND	4.0	EPA 200.8		3-5-24	
Mercury	ND	0.50	EPA 7470A		2-15-24	

Client ID:	MW-6					
Laboratory ID:	02-047-04					
Arsenic	ND	4.0	EPA 200.8		3-5-24	
Cadmium	ND	4.0	EPA 200.8		3-5-24	
Chromium	ND	10	EPA 200.8		3-5-24	
Lead	ND	4.0	EPA 200.8		3-5-24	
Mercury	ND	0.50	EPA 7470A		2-15-24	



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DISSOLVED METALS
EPA 200.8/7470A

Matrix: Water
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-4					
Laboratory ID:	02-047-05					
Arsenic	ND	4.0	EPA 200.8		3-5-24	
Cadmium	ND	4.0	EPA 200.8		3-5-24	
Chromium	ND	10	EPA 200.8		3-5-24	
Lead	ND	4.0	EPA 200.8		3-5-24	
Mercury	ND	0.50	EPA 7470A		2-15-24	

Client ID:	MW-2					
Laboratory ID:	02-047-06					
Arsenic	7.8	4.4	EPA 200.8	2-26-24	3-5-24	
Cadmium	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Chromium	ND	11	EPA 200.8	2-26-24	3-5-24	
Lead	ND	4.4	EPA 200.8	2-26-24	3-5-24	
Mercury	ND	0.50	EPA 7470A		2-15-24	



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**DISSOLVED METALS
 EPA 200.8/7470A
 QUALITY CONTROL**

Matrix: Water
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0226WM1					
Arsenic	ND	3.3	EPA 200.8	2-26-24	2-26-24	
Cadmium	ND	4.4	EPA 200.8	2-26-24	2-26-24	
Chromium	ND	11	EPA 200.8	2-26-24	2-26-24	
Lead	ND	1.1	EPA 200.8	2-26-24	2-26-24	

Laboratory ID:	MB0215D1					
Mercury	ND	0.50	EPA 7470A	2-15-24	2-15-24	

Laboratory ID:	MB0305D1					
Arsenic	ND	3.0	EPA 200.8		3-5-24	
Cadmium	ND	4.0	EPA 200.8		3-5-24	
Chromium	ND	10	EPA 200.8		3-5-24	
Lead	ND	1.0	EPA 200.8		3-5-24	

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
DUPLICATE								
Laboratory ID:	01-067-08							
	ORIG	DUP						
Arsenic	ND	ND	NA	NA	NA	NA	NA	20
Cadmium	ND	ND	NA	NA	NA	NA	NA	20
Chromium	ND	ND	NA	NA	NA	NA	NA	20
Lead	ND	ND	NA	NA	NA	NA	NA	20

Laboratory ID:	02-047-01							
Mercury	ND	ND	NA	NA	NA	NA	NA	20

Laboratory ID:	02-045-02							
	ORIG	DUP						
Arsenic	6.78	6.36	NA	NA	NA	NA	6	20
Cadmium	ND	ND	NA	NA	NA	NA	NA	20
Chromium	ND	ND	NA	NA	NA	NA	NA	20
Lead	ND	ND	NA	NA	NA	NA	NA	20



Date of Report: March 5, 2024
 Samples Submitted: February 7, 2024
 Laboratory Reference: 2402-047
 Project: 0209325-000

**DISSOLVED METALS
 EPA 200.8/7470A
 QUALITY CONTROL**

Matrix: Water
 Units: ug/L (ppb)

Analyte	Result		Spike Level		Source	Percent	Recovery	RPD		Flags
					Result	Recovery	Limits	RPD	Limit	
MATRIX SPIKES										
Laboratory ID:		01-067-08								
	MS	MSD	MS	MSD		MS	MSD			
Arsenic	113	112	111	111	ND	102	101	75-125	1	20
Cadmium	111	109	111	111	ND	100	98	75-125	2	20
Chromium	112	111	111	111	ND	101	100	75-125	1	20
Lead	109	109	111	111	ND	99	98	75-125	0	20
Laboratory ID:		02-047-01								
Mercury	12.1	12.1	12.5	12.5	ND	97	97	75-125	0	20
Laboratory ID:		02-045-02								
	MS	MSD	MS	MSD		MS	MSD			
Arsenic	89.6	90.8	80.0	80.0	6.78	104	105	75-125	1	20
Cadmium	76.2	77.2	80.0	80.0	ND	95	97	75-125	1	20
Chromium	80.4	79.8	80.0	80.0	ND	101	100	75-125	1	20
Lead	63.2	64.4	80.0	80.0	ND	79	81	75-125	2	20
SPIKE BLANK										
Laboratory ID:		SB0226WM1								
Arsenic	112		111		N/A	101		85-115		
Cadmium	111		111		N/A	100		85-115		
Chromium	110		111		N/A	99		85-115		
Lead	112		111		N/A	101		85-115		
Laboratory ID:		SB0215D1								
Mercury	12.0		12.5		N/A	96		80-120		
Laboratory ID:		SB0305D1								
Arsenic	79.8		80.0		N/A	100		85-115		
Cadmium	81.2		80.0		N/A	102		85-115		
Chromium	81.2		80.0		N/A	102		85-115		
Lead	80.8		80.0		N/A	101		85-115		





Data Qualifiers and Abbreviations

- A - Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
 - B - The analyte indicated was also found in the blank sample.
 - C - The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
 - E - The value reported exceeds the quantitation range and is an estimate.
 - F - Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
 - H - The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
 - I - Compound recovery is outside of the control limits.
 - J - The value reported was below the practical quantitation limit. The value is an estimate.
 - K - Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
 - L - The RPD is outside of the control limits.
 - M - Hydrocarbons in the gasoline range are impacting the diesel range result.
 - M1 - Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
 - N - Hydrocarbons in the lube oil range are impacting the diesel range result.
 - N1 - Hydrocarbons in diesel range are impacting lube oil range results.
 - O - Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
 - P - The RPD of the detected concentrations between the two columns is greater than 40.
 - Q - Surrogate recovery is outside of the control limits.
 - S - Surrogate recovery data is not available due to the necessary dilution of the sample.
 - T - The sample chromatogram is not similar to a typical _____.
 - U - The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
 - U1 - The practical quantitation limit is elevated due to interferences present in the sample.
 - V - Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
 - W - Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
 - X - Sample extract treated with a mercury cleanup procedure.
 - X1 - Sample extract treated with a sulfuric acid/silica gel cleanup procedure.
 - X2 - Sample extract treated with a silica gel cleanup procedure.
 - Y - The calibration verification for this analyte exceeded the 20% drift specified in methods 8260 & 8270, and therefore the reported result should be considered an estimate. The overall performance of the calibration verification standard met the acceptance criteria of the method.
 - Y1 - Negative effects of the matrix from this sample on the instrument caused values for this analyte in the bracketing continuing calibration verification standard (CCVs) to be outside of 20% acceptance criteria. Because of this, quantitation limits and sample concentrations should be considered estimates.
 - Z -
- ND - Not Detected at PQL
 PQL - Practical Quantitation Limit
 RPD - Relative Percent Difference





OnSite Environmental Inc.
 Analytical Laboratory Testing Services
 14648 NE 95th Street • Redmond, WA 98052
 Phone: (425) 883-3881 • www.onsite-env.com

Chain of Custody

Turnaround Request
 (in working days)
 (Check One)

- Same Day 1 Day
- 2 Days 3 Days
- Standard (7 Days)
- _____ (other)

Laboratory Number: **02-047**

Company: **HALEY 3 ANDRICH**
 Project Number: **0209325-600**
 Project Name: **CUSTOM PLUMBING**
 Project Manager: **A-KAPAROS**
 Sampled by: **C. DOVITHITI, S. FISHER**

Lab ID	Sample Identification	Date Sampled	Time Sampled	Matrix
1	MW-1	2/1/24	15:40	GW
2	MW-3	2/1/24	16:10	GW
3	MW-5	2/2/24	10:20	GW
4	MW-6	2/2/24	10:45	GW
5	MW-4	2/2/24	12:15	GW
6	MW-2	2/2/24	13:00	GW

Number of Containers	
NWTPH-HCID	
NWTPH-Gx/BTEX (8021 <input type="checkbox"/> 8260 <input type="checkbox"/>	
NWTPH-Gx	
NWTPH-Dx (SG Clean-up) <input checked="" type="checkbox"/>	WITH + WITHOUT
Volatiles 8260	
Halogenated Volatiles 8260	
EDB EPA 8011 (Waters Only)	
Semivolatiles 8270/SIM (with low-level PAHs)	
PAHs 8270/SIM (low-level)	C PAHs
PCBs 8082	
Organochlorine Pesticides 8081	
Organophosphorus Pesticides 8270/SIM	
Chlorinated Acid Herbicides 8151	
Total RCRA Metals	
Total MTCA Metals	
TCLP Metals	
HEM (oil and grease) 1664	
DISSOLVED METALS *	
% Moisture	

Signature	Company	Date	Time	Comments/Special Instructions
	HALEY 3 ANDRICH	2/2/24	15:42	DISCOVERED METALS FIELD FILTERED Please send results to skypavos@halcyonenvironmental.com AND update me @ halcyon@halcyonenv.com
	CD	2/2/24	15:42	

Data Package: Standard Level III Level IV
 Chromatograms with final report Electronic Data Deliverables (EDDs)



14648 NE 95th Street, Redmond, WA 98052 • (425) 883-3881

August 23, 2024

Andrew Kaparos
Hart Crowser, Inc.
A Division of Haley & Aldrich, Inc.
3131 Elliott Avenue, Suite 600
Seattle, WA 98121

Re: Analytical Data for Project 0209325-000
Laboratory Reference No. 2408-148

Dear Andrew:

Enclosed are the analytical results and associated quality control data for samples submitted on August 13, 2024.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal flourish extending to the right.

David Baumeister
Project Manager

Enclosures



OnSite Environmental, Inc. 14648 NE 95th Street, Redmond, WA 98052 (425) 883-3881

This report pertains to the samples analyzed in accordance with the chain of custody, and is intended only for the use of the individual or company to whom it is addressed.

Date of Report: August 23, 2024
Samples Submitted: August 13, 2024
Laboratory Reference: 2408-148
Project: 0209325-000

Case Narrative

Samples were collected on August 12, 2024 and received by the laboratory on August 13, 2024. They were maintained at the laboratory at a temperature of 2°C to 6°C.

Please note that any and all soil sample results are reported on a dry-weight basis, unless otherwise noted below. However the soil results for the QA/QC samples are reported on a wet-weight basis.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

NWTPH Dx Analysis

The surrogate percent recovery in sample MW-3 was below the control limit of 50% due to matrix effects. The sample was re-extracted with the same result.

PAHs EPA 8270E/SIM Analysis

Sample MW-1 had one surrogate recovery outside of control limits. This is within allowance of our standard operating procedure as long as the recovery is above 10%.

Dissolved Metals EPA 200.8/7470A Analysis

The dissolved field filter samples MW-3, MW-2, and MW-1 were received containing solid material. The samples were digested according to OnSite Environmental standard operating procedure.

Please note that any other QA/QC issues associated with these extractions and analyses will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.



Date of Report: August 23, 2024
 Samples Submitted: August 13, 2024
 Laboratory Reference: 2408-148
 Project: 0209325-000

**DIESEL AND HEAVY OIL RANGE ORGANICS
 NWTPH-Dx**

Matrix: Water
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-4					
Laboratory ID:	08-148-01					
Diesel Range Organics	ND	0.22	NWTPH-Dx	8-16-24	8-16-24	
Lube Oil Range Organics	ND	0.22	NWTPH-Dx	8-16-24	8-16-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	112	50-150				
Client ID:	MW-3					
Laboratory ID:	08-148-02					
Diesel Range Organics	0.15	0.14	NWTPH-Dx	8-16-24	8-16-24	
Lube Oil Range Organics	0.32	0.22	NWTPH-Dx	8-16-24	8-16-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	33	50-150				Q
Client ID:	MW-2					
Laboratory ID:	08-148-03					
Diesel Range Organics	0.33	0.21	NWTPH-Dx	8-16-24	8-17-24	
Lube Oil Range Organics	0.35	0.21	NWTPH-Dx	8-16-24	8-17-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	85	50-150				



Date of Report: August 23, 2024
 Samples Submitted: August 13, 2024
 Laboratory Reference: 2408-148
 Project: 0209325-000

**DIESEL AND HEAVY OIL RANGE ORGANICS
 NWTPH-Dx**

Matrix: Water
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-5					
Laboratory ID:	08-148-04					
Diesel Range Organics	ND	0.20	NWTPH-Dx	8-16-24	8-16-24	
Lube Oil Range Organics	ND	0.20	NWTPH-Dx	8-16-24	8-16-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	104	50-150				

Client ID:	MW-1					
Laboratory ID:	08-148-05					
Diesel Range Organics	ND	0.23	NWTPH-Dx	8-16-24	8-16-24	
Lube Oil Range Organics	0.24	0.23	NWTPH-Dx	8-16-24	8-16-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	84	50-150				

Client ID:	MW-6					
Laboratory ID:	08-148-06					
Diesel Range Organics	ND	0.22	NWTPH-Dx	8-16-24	8-16-24	
Lube Oil Range Organics	ND	0.22	NWTPH-Dx	8-16-24	8-16-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	117	50-150				



Date of Report: August 23, 2024
 Samples Submitted: August 13, 2024
 Laboratory Reference: 2408-148
 Project: 0209325-000

**DIESEL AND HEAVY OIL RANGE ORGANICS
 NWTPH-Dx
 QUALITY CONTROL**

Matrix: Water
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0816W1					
Diesel Range Organics	ND	0.10	NWTPH-Dx	8-16-24	8-16-24	
Lube Oil Range Organics	ND	0.16	NWTPH-Dx	8-16-24	8-16-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	94	50-150				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
DUPLICATE								
Laboratory ID:	SB0816W1							
	ORIG	DUP						
Diesel Fuel #2	0.552	0.533	NA	NA	NA	NA	4	40
<i>Surrogate:</i>								
<i>o-Terphenyl</i>				125	128	50-150		

SPIKE BLANK								
Laboratory ID:	SB0816W1							
Diesel Fuel #2	0.552		0.500	NA	110	50-129	NA	NA
<i>Surrogate:</i>								
<i>o-Terphenyl</i>					125	50-150		



Date of Report: August 23, 2024
 Samples Submitted: August 13, 2024
 Laboratory Reference: 2408-148
 Project: 0209325-000

PAHs EPA 8270E/SIM

Matrix: Water
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-4					
Laboratory ID:	08-148-01					
Naphthalene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
2-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
1-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
Acenaphthylene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
Acenaphthene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
Fluorene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
Phenanthrene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
Anthracene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
Fluoranthene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
Pyrene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[a]anthracene	ND	0.010	EPA 8270E/SIM	8-19-24	8-20-24	
Chrysene	ND	0.010	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[b]fluoranthene	ND	0.010	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[a]pyrene	ND	0.010	EPA 8270E/SIM	8-19-24	8-20-24	
Indeno(1,2,3-c,d)pyrene	ND	0.014	EPA 8270E/SIM	8-19-24	8-20-24	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[g,h,i]perylene	ND	0.012	EPA 8270E/SIM	8-19-24	8-20-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>51</i>	<i>27-106</i>				
<i>Pyrene-d10</i>	<i>76</i>	<i>37-125</i>				
<i>Terphenyl-d14</i>	<i>75</i>	<i>37-110</i>				



Date of Report: August 23, 2024
 Samples Submitted: August 13, 2024
 Laboratory Reference: 2408-148
 Project: 0209325-000

PAHs EPA 8270E/SIM

Matrix: Water
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-3					
Laboratory ID:	08-148-02					
Naphthalene	ND	0.098	EPA 8270E/SIM	8-19-24	8-20-24	
2-Methylnaphthalene	ND	0.098	EPA 8270E/SIM	8-19-24	8-20-24	
1-Methylnaphthalene	ND	0.098	EPA 8270E/SIM	8-19-24	8-20-24	
Acenaphthylene	ND	0.098	EPA 8270E/SIM	8-19-24	8-20-24	
Acenaphthene	ND	0.098	EPA 8270E/SIM	8-19-24	8-20-24	
Fluorene	ND	0.098	EPA 8270E/SIM	8-19-24	8-20-24	
Phenanthrene	ND	0.098	EPA 8270E/SIM	8-19-24	8-20-24	
Anthracene	ND	0.098	EPA 8270E/SIM	8-19-24	8-20-24	
Fluoranthene	ND	0.098	EPA 8270E/SIM	8-19-24	8-20-24	
Pyrene	ND	0.098	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[a]anthracene	ND	0.0098	EPA 8270E/SIM	8-19-24	8-20-24	
Chrysene	ND	0.0098	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[b]fluoranthene	ND	0.0098	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo(j,k)fluoranthene	ND	0.0098	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[a]pyrene	ND	0.0098	EPA 8270E/SIM	8-19-24	8-20-24	
Indeno(1,2,3-c,d)pyrene	ND	0.0098	EPA 8270E/SIM	8-19-24	8-20-24	
Dibenz[a,h]anthracene	ND	0.0098	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[g,h,i]perylene	ND	0.0098	EPA 8270E/SIM	8-19-24	8-20-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorobiphenyl	59	27-106				
Pyrene-d10	64	37-125				
Terphenyl-d14	77	37-110				



Date of Report: August 23, 2024
 Samples Submitted: August 13, 2024
 Laboratory Reference: 2408-148
 Project: 0209325-000

PAHs EPA 8270E/SIM

Matrix: Water
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-2					
Laboratory ID:	08-148-03					
Naphthalene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
2-Methylnaphthalene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
1-Methylnaphthalene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Acenaphthylene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Acenaphthene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Fluorene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Phenanthrene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Anthracene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Fluoranthene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Pyrene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[a]anthracene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Chrysene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[b]fluoranthene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo(j,k)fluoranthene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[a]pyrene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Indeno(1,2,3-c,d)pyrene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Dibenz[a,h]anthracene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[g,h,i]perylene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>27</i>	<i>27-106</i>				
<i>Pyrene-d10</i>	<i>54</i>	<i>37-125</i>				
<i>Terphenyl-d14</i>	<i>65</i>	<i>37-110</i>				



Date of Report: August 23, 2024
 Samples Submitted: August 13, 2024
 Laboratory Reference: 2408-148
 Project: 0209325-000

PAHs EPA 8270E/SIM

Matrix: Water
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-5					
Laboratory ID:	08-148-04					
Naphthalene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
2-Methylnaphthalene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
1-Methylnaphthalene	0.35	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Acenaphthylene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Acenaphthene	0.11	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Fluorene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Phenanthrene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Anthracene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Fluoranthene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Pyrene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[a]anthracene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Chrysene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[b]fluoranthene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo(j,k)fluoranthene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[a]pyrene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Indeno(1,2,3-c,d)pyrene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Dibenz[a,h]anthracene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[g,h,i]perylene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorobiphenyl	53	27-106				
Pyrene-d10	69	37-125				
Terphenyl-d14	77	37-110				



Date of Report: August 23, 2024
 Samples Submitted: August 13, 2024
 Laboratory Reference: 2408-148
 Project: 0209325-000

PAHs EPA 8270E/SIM

Matrix: Water
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-1					
Laboratory ID:	08-148-05					
Naphthalene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
2-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
1-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
Acenaphthylene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
Acenaphthene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
Fluorene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
Phenanthrene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
Anthracene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
Fluoranthene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
Pyrene	ND	0.10	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[a]anthracene	ND	0.010	EPA 8270E/SIM	8-19-24	8-20-24	
Chrysene	ND	0.010	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[b]fluoranthene	ND	0.010	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[a]pyrene	ND	0.010	EPA 8270E/SIM	8-19-24	8-20-24	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270E/SIM	8-19-24	8-20-24	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270E/SIM	8-19-24	8-20-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorobiphenyl	68	27-106				
Pyrene-d10	74	37-125				
Terphenyl-d14	155	37-110				Q



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PAHs EPA 8270E/SIM

Matrix: Water
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-6					
Laboratory ID:	08-148-06					
Naphthalene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
2-Methylnaphthalene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
1-Methylnaphthalene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Acenaphthylene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Acenaphthene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Fluorene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Phenanthrene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Anthracene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Fluoranthene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Pyrene	ND	0.095	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[a]anthracene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Chrysene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[b]fluoranthene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo(j,k)fluoranthene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[a]pyrene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Indeno(1,2,3-c,d)pyrene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Dibenz[a,h]anthracene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
Benzo[g,h,i]perylene	ND	0.0095	EPA 8270E/SIM	8-19-24	8-20-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorobiphenyl	57	27-106				
Pyrene-d10	82	37-125				
Terphenyl-d14	78	37-110				



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**PAHs EPA 8270E/SIM
 QUALITY CONTROL**

Matrix: Water
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0819W1					
Naphthalene	ND	0.10	EPA 8270E/SIM	8-19-24	8-19-24	
2-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	8-19-24	8-19-24	
1-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	8-19-24	8-19-24	
Acenaphthylene	ND	0.10	EPA 8270E/SIM	8-19-24	8-19-24	
Acenaphthene	ND	0.10	EPA 8270E/SIM	8-19-24	8-19-24	
Fluorene	ND	0.10	EPA 8270E/SIM	8-19-24	8-19-24	
Phenanthrene	ND	0.10	EPA 8270E/SIM	8-19-24	8-19-24	
Anthracene	ND	0.10	EPA 8270E/SIM	8-19-24	8-19-24	
Fluoranthene	ND	0.10	EPA 8270E/SIM	8-19-24	8-19-24	
Pyrene	ND	0.10	EPA 8270E/SIM	8-19-24	8-19-24	
Benzo[a]anthracene	ND	0.010	EPA 8270E/SIM	8-19-24	8-19-24	
Chrysene	ND	0.010	EPA 8270E/SIM	8-19-24	8-19-24	
Benzo[b]fluoranthene	ND	0.010	EPA 8270E/SIM	8-19-24	8-19-24	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270E/SIM	8-19-24	8-19-24	
Benzo[a]pyrene	ND	0.010	EPA 8270E/SIM	8-19-24	8-19-24	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270E/SIM	8-19-24	8-19-24	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270E/SIM	8-19-24	8-19-24	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270E/SIM	8-19-24	8-19-24	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>55</i>	<i>27-106</i>				
<i>Pyrene-d10</i>	<i>67</i>	<i>37-125</i>				
<i>Terphenyl-d14</i>	<i>67</i>	<i>37-110</i>				



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**PAHs EPA 8270E/SIM
 QUALITY CONTROL**

Matrix: Water
 Units: ug/L

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD		Flags
					Recovery	Limits	RPD	Limit		
SPIKE BLANKS										
Laboratory ID:	SB0819W1									
	SB	SBD	SB	SBD	SB	SBD				
Naphthalene	0.281	0.292	0.500	0.500	56	58	36-95	4	34	
Acenaphthylene	0.330	0.365	0.500	0.500	66	73	51-103	10	26	
Acenaphthene	0.312	0.341	0.500	0.500	62	68	47-97	9	25	
Fluorene	0.340	0.374	0.500	0.500	68	75	51-105	10	23	
Phenanthrene	0.422	0.458	0.500	0.500	84	92	52-110	8	24	
Anthracene	0.360	0.388	0.500	0.500	72	78	55-104	7	24	
Fluoranthene	0.427	0.403	0.500	0.500	85	81	59-111	6	24	
Pyrene	0.404	0.433	0.500	0.500	81	87	59-110	7	22	
Benzo[a]anthracene	0.442	0.443	0.500	0.500	88	89	55-116	0	22	
Chrysene	0.388	0.405	0.500	0.500	78	81	59-111	4	23	
Benzo[b]fluoranthene	0.445	0.442	0.500	0.500	89	88	62-115	1	27	
Benzo(j,k)fluoranthene	0.397	0.409	0.500	0.500	79	82	59-117	3	23	
Benzo[a]pyrene	0.434	0.444	0.500	0.500	87	89	64-109	2	24	
Indeno(1,2,3-c,d)pyrene	0.483	0.476	0.500	0.500	97	95	58-114	1	22	
Dibenz[a,h]anthracene	0.503	0.509	0.500	0.500	101	102	63-114	1	24	
Benzo[g,h,i]perylene	0.409	0.421	0.500	0.500	82	84	61-110	3	24	
<i>Surrogate:</i>										
2-Fluorobiphenyl					59	63	27-106			
Pyrene-d10					83	86	37-125			
Terphenyl-d14					82	84	37-110			



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TOTAL METALS
EPA 200.8/7470A

Matrix: Water
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-4					
Laboratory ID:	08-148-01					
Arsenic	3.2	0.56	EPA 200.8	8-19-24	8-20-24	
Cadmium	ND	0.22	EPA 200.8	8-19-24	8-20-24	
Chromium	ND	1.1	EPA 200.8	8-19-24	8-20-24	
Lead	ND	0.44	EPA 200.8	8-19-24	8-20-24	
Mercury	ND	0.13	EPA 7470A	8-22-24	8-22-24	

Client ID:	MW-3					
Laboratory ID:	08-148-02					
Arsenic	2.3	0.56	EPA 200.8	8-19-24	8-20-24	
Cadmium	ND	0.22	EPA 200.8	8-19-24	8-20-24	
Chromium	9.2	1.1	EPA 200.8	8-19-24	8-20-24	
Lead	ND	0.44	EPA 200.8	8-19-24	8-20-24	
Mercury	ND	0.13	EPA 7470A	8-22-24	8-22-24	

Client ID:	MW-2					
Laboratory ID:	08-148-03					
Arsenic	6.4	0.56	EPA 200.8	8-19-24	8-20-24	
Cadmium	ND	0.22	EPA 200.8	8-19-24	8-20-24	
Chromium	12	1.1	EPA 200.8	8-19-24	8-20-24	
Lead	1.3	0.44	EPA 200.8	8-19-24	8-20-24	
Mercury	ND	0.13	EPA 7470A	8-22-24	8-22-24	

Client ID:	MW-5					
Laboratory ID:	08-148-04					
Arsenic	0.68	0.56	EPA 200.8	8-19-24	8-20-24	
Cadmium	ND	0.22	EPA 200.8	8-19-24	8-20-24	
Chromium	ND	1.1	EPA 200.8	8-19-24	8-20-24	
Lead	ND	0.44	EPA 200.8	8-19-24	8-20-24	
Mercury	ND	0.13	EPA 7470A	8-22-24	8-22-24	



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**TOTAL METALS
 EPA 200.8/7470A**

Matrix: Water
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-1					
Laboratory ID:	08-148-05					
Arsenic	27	0.56	EPA 200.8	8-19-24	8-20-24	
Cadmium	0.60	0.22	EPA 200.8	8-19-24	8-20-24	
Chromium	12	1.1	EPA 200.8	8-19-24	8-20-24	
Lead	3.3	0.44	EPA 200.8	8-19-24	8-20-24	
Mercury	1.0	0.13	EPA 7470A	8-22-24	8-22-24	

Client ID:	MW-6					
Laboratory ID:	08-148-06					
Arsenic	2.6	0.56	EPA 200.8	8-19-24	8-20-24	
Cadmium	ND	0.22	EPA 200.8	8-19-24	8-20-24	
Chromium	ND	1.1	EPA 200.8	8-19-24	8-20-24	
Lead	ND	0.44	EPA 200.8	8-19-24	8-20-24	
Mercury	ND	0.13	EPA 7470A	8-22-24	8-22-24	



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**TOTAL METALS
 EPA 200.8/7470A
 QUALITY CONTROL**

Matrix: Water
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0819WM1					
Arsenic	ND	0.56	EPA 200.8	8-19-24	8-20-24	
Cadmium	ND	0.22	EPA 200.8	8-19-24	8-20-24	
Chromium	ND	1.1	EPA 200.8	8-19-24	8-20-24	
Lead	ND	0.44	EPA 200.8	8-19-24	8-20-24	
Laboratory ID:	MB0822W1					
Mercury	ND	0.13	EPA 7470A	8-22-24	8-22-24	



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**TOTAL METALS
 EPA 200.8/7470A
 QUALITY CONTROL**

Matrix: Water
 Units: ug/L (ppb)

Analyte	Result		Spike Level		Source	Percent	Recovery	RPD		Flags
					Result	Recovery	Limits	RPD	Limit	
DUPLICATE										
Laboratory ID:	08-028-05									
	ORIG	DUP								
Arsenic	ND	ND	NA	NA		NA	NA	NA	20	
Cadmium	ND	ND	NA	NA		NA	NA	NA	20	
Chromium	ND	ND	NA	NA		NA	NA	NA	20	
Lead	ND	ND	NA	NA		NA	NA	NA	20	
Laboratory ID:	08-140-02									
Mercury	ND	ND	NA	NA		NA	NA	NA	20	
MATRIX SPIKES										
Laboratory ID:	08-028-05									
	MS	MSD	MS	MSD		MS	MSD			
Arsenic	113	119	111	111	ND	102	108	75-125	5	20
Cadmium	115	121	111	111	ND	104	109	75-125	5	20
Chromium	114	121	111	111	ND	103	109	75-125	6	20
Lead	113	118	111	111	ND	102	107	75-125	4	20
Laboratory ID:	08-140-02									
Mercury	12.2	12.2	12.5	12.5	ND	97	97	75-125	0	20
SPIKE BLANK										
Laboratory ID:	SB0819WM1									
Arsenic	115		111		N/A	103		85-115		
Cadmium	117		111		N/A	105		85-115		
Chromium	115		111		N/A	103		85-115		
Lead	111		111		N/A	100		85-115		
Laboratory ID:	SB0822W1									
Mercury	12.3		12.5		N/A	99		80-120		



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DISSOLVED METALS
EPA 200.8/7470A

Matrix: Water
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-4					
Laboratory ID:	08-148-01					
Arsenic	2.4	0.50	EPA 200.8		8-20-24	
Cadmium	ND	0.20	EPA 200.8		8-20-24	
Chromium	ND	1.0	EPA 200.8		8-20-24	
Lead	ND	0.40	EPA 200.8		8-20-24	
Mercury	ND	0.13	EPA 7470A		8-22-24	

Client ID:	MW-3					
Laboratory ID:	08-148-02					
Arsenic	1.4	0.56	EPA 200.8	8-19-24	8-20-24	
Cadmium	ND	0.22	EPA 200.8	8-19-24	8-20-24	
Chromium	9.2	1.1	EPA 200.8	8-19-24	8-20-24	
Lead	ND	0.44	EPA 200.8	8-19-24	8-20-24	
Mercury	ND	0.13	EPA 7470A		8-22-24	

Client ID:	MW-2					
Laboratory ID:	08-148-03					
Arsenic	3.6	0.56	EPA 200.8	8-19-24	8-20-24	
Cadmium	ND	0.22	EPA 200.8	8-19-24	8-20-24	
Chromium	8.5	1.1	EPA 200.8	8-19-24	8-20-24	
Lead	ND	0.44	EPA 200.8	8-19-24	8-20-24	
Mercury	ND	0.13	EPA 7470A		8-22-24	

Client ID:	MW-5					
Laboratory ID:	08-148-04					
Arsenic	0.67	0.50	EPA 200.8		8-20-24	
Cadmium	ND	0.20	EPA 200.8		8-20-24	
Chromium	ND	1.0	EPA 200.8		8-20-24	
Lead	ND	0.40	EPA 200.8		8-20-24	
Mercury	ND	0.13	EPA 7470A		8-22-24	



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DISSOLVED METALS
EPA 200.8/7470A

Matrix: Water
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	MW-1					
Laboratory ID:	08-148-05					
Arsenic	27	0.56	EPA 200.8	8-19-24	8-20-24	
Cadmium	0.31	0.22	EPA 200.8	8-19-24	8-20-24	
Chromium	5.9	1.1	EPA 200.8	8-19-24	8-20-24	
Lead	0.87	0.44	EPA 200.8	8-19-24	8-20-24	
Mercury	0.16	0.13	EPA 7470A		8-22-24	

Client ID:	MW-6					
Laboratory ID:	08-148-06					
Arsenic	2.1	0.50	EPA 200.8		8-20-24	
Cadmium	ND	0.20	EPA 200.8		8-20-24	
Chromium	ND	1.0	EPA 200.8		8-20-24	
Lead	ND	0.40	EPA 200.8		8-20-24	
Mercury	ND	0.13	EPA 7470A		8-22-24	



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**DISSOLVED METALS
 EPA 200.8/7470A
 QUALITY CONTROL**

Matrix: Water
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0819D1					
Arsenic	ND	0.50	EPA 200.8		8-19-24	
Cadmium	ND	0.20	EPA 200.8		8-19-24	
Chromium	ND	1.0	EPA 200.8		8-19-24	
Lead	ND	0.40	EPA 200.8		8-19-24	

Laboratory ID:	MB0822D1					
Mercury	ND	0.13	EPA 7470A		8-22-24	

Laboratory ID:	MB0819WM1					
Arsenic	ND	0.56	EPA 200.8		8-19-24	
Cadmium	ND	0.22	EPA 200.8		8-19-24	
Chromium	ND	1.1	EPA 200.8		8-19-24	
Lead	ND	0.44	EPA 200.8		8-19-24	

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
DUPLICATE								
Laboratory ID:	08-140-02							
	ORIG	DUP						
Arsenic	ND	ND	NA	NA	NA	NA	NA	20
Cadmium	ND	ND	NA	NA	NA	NA	NA	20
Chromium	ND	ND	NA	NA	NA	NA	NA	20
Lead	ND	ND	NA	NA	NA	NA	NA	20

Laboratory ID:	08-140-02							
Mercury	ND	ND	NA	NA	NA	NA	NA	20

Laboratory ID:	08-028-05							
	ORIG	DUP						
Arsenic	ND	ND	NA	NA	NA	NA	NA	20
Cadmium	ND	ND	NA	NA	NA	NA	NA	20
Chromium	ND	ND	NA	NA	NA	NA	NA	20
Lead	ND	ND	NA	NA	NA	NA	NA	20



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**DISSOLVED METALS
 EPA 200.8/7470A
 QUALITY CONTROL**

Matrix: Water
 Units: ug/L (ppb)

Analyte	Result		Spike Level		Source	Percent	Recovery	RPD		Flags
					Result	Recovery	Limits	RPD	Limit	
MATRIX SPIKES										
Laboratory ID:	08-140-02									
	MS	MSD	MS	MSD		MS	MSD			
Arsenic	85.4	88.0	80.0	80.0	ND	107	110	75-125	3	20
Cadmium	75.8	78.8	80.0	80.0	ND	95	99	75-125	4	20
Chromium	73.8	75.0	80.0	80.0	ND	92	94	75-125	2	20
Lead	76.0	76.8	80.0	80.0	ND	95	96	75-125	1	20
Laboratory ID:	08-140-02									
Mercury	11.9	12.2	12.5	12.5	ND	95	97	75-125	2	20
Laboratory ID:	08-028-05									
	MS	MSD	MS	MSD		MS	MSD			
Arsenic	113	119	80.0	80.0	ND	142	149	75-125	5	20
Cadmium	115	121	80.0	80.0	ND	144	152	75-125	5	20
Chromium	114	121	80.0	80.0	ND	143	151	75-125	6	20
Lead	113	118	80.0	80.0	ND	141	148	75-125	4	20
SPIKE BLANK										
Laboratory ID:	SB0819D1									
Arsenic	79.4		80.0		N/A	99		85-115		
Cadmium	80.2		80.0		N/A	100		85-115		
Chromium	80.4		80.0		N/A	101		85-115		
Lead	81.6		80.0		N/A	102		85-115		
Laboratory ID:	SB0822D1									
Mercury	12.1		12.5		N/A	97		80-120		
Laboratory ID:	SB0819WM1									
Arsenic	115		80.0		N/A	143		85-115		
Cadmium	117		80.0		N/A	146		85-115		
Chromium	115		80.0		N/A	143		85-115		
Lead	111		80.0		N/A	138		85-115		





Data Qualifiers and Abbreviations

- A - Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
 - B - The analyte indicated was also found in the blank sample.
 - C - The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
 - E - The value reported exceeds the quantitation range and is an estimate.
 - F - Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
 - H - The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
 - I - Compound recovery is outside of the control limits.
 - J - The value reported was below the practical quantitation limit. The value is an estimate.
 - K - Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
 - L - The RPD is outside of the control limits.
 - M - Hydrocarbons in the gasoline range are impacting the diesel range result.
 - M1 - Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
 - N - Hydrocarbons in the lube oil range are impacting the diesel range result.
 - N1 - Hydrocarbons in diesel range are impacting lube oil range results.
 - O - Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
 - P - The RPD of the detected concentrations between the two columns is greater than 40.
 - Q - Surrogate recovery is outside of the control limits.
 - S - Surrogate recovery data is not available due to the necessary dilution of the sample.
 - T - The sample chromatogram is not similar to a typical _____.
 - U - The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
 - U1 - The practical quantitation limit is elevated due to interferences present in the sample.
 - V - Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
 - W - Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
 - X - Sample extract treated with a mercury cleanup procedure.
 - X1 - Sample extract treated with a sulfuric acid/silica gel cleanup procedure.
 - X2 - Sample extract treated with a silica gel cleanup procedure.
 - Y - The calibration verification for this analyte exceeded the 20% drift specified in methods 8260 & 8270, and therefore the reported result should be considered an estimate. The overall performance of the calibration verification standard met the acceptance criteria of the method.
 - Y1 - Negative effects of the matrix from this sample on the instrument caused values for this analyte in the bracketing continuing calibration verification standard (CCVs) to be outside of 20% acceptance criteria. Because of this, quantitation limits and sample concentrations should be considered estimates.
 - Z -
- ND - Not Detected at PQL
 PQL - Practical Quantitation Limit
 RPD - Relative Percent Difference





Onsite Environmental Inc.
Analytical Laboratory Testing Services

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Phone: (425) 883-3881 • www.onsite-env.com

Chain of Custody

Turnaround Request
(in working days)

(Check One)

- Same Day 1 Day
 2 Days 3 Days
 Standard (7 Days)

_____ (other)

Laboratory Number: **08-148**

Company: **MACEY 3 ADDRESS**
 Project Number: **0209325-000**
 Project Name: **CUSTOM (RYW057)**
 Project Manager: **A. KAPPAOS**
 Sampled by: **S. FISHER, M. BANU & ASSER**

Lab ID	Sample Identification	Signature
1	MW-4	<i>[Signature]</i>
2	MW-3	<i>[Signature]</i>
3	MW-2	<i>[Signature]</i>
4	MW-5	<i>[Signature]</i>
5	MW-1	<i>[Signature]</i>
6	MW-6	<i>[Signature]</i>

Date Sampled	Time Sampled	Matrix
8/12/24	10:40	GW
8/12/24	12:50	GW
8/12/24	13:00	GW
8/12/24	15:30	GW
8/12/24	15:40	GW
8/12/24	17:05	GW

Number of Containers	NWTPH-HCID	NWTPH-Gx/BTEX (8021 <input type="checkbox"/> 8260 <input checkbox="" type="checkbox/>)</th> <th>NWTPH-Gx</th> <th>NWTPH-Dx (SG Clean-up <input type="/>)	Volatiles 8260	Halogenated Volatiles 8260	EDB EPA 8011 (Waters Only)	Semivolatiles 8270/SIM (with low-level PAHs)	PAHs 8270/SIM (low-level)	PCBs 8082	Organochlorine Pesticides 8081	Organophosphorus Pesticides 8270/SIM	Chlorinated Acid Herbicides 8151	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664	% Moisture		
6			X	X				X	X									
6			X	X				X	X									
6			X	X				X	X									
6			X	X				X	X									
6			X	X				X	X									
6			X	X				X	X									

Company	Date	Time
MACEY 3 ADDRESS	8/13/24	16:02
<i>[Signature]</i>	8/13/24	16:02

Comments/Special Instructions
 * DISSOLVED METALS FIELD FILTERED
 - Please send results to
 kapaos@halogaldiv.com
 vphdivan@halogaldiv.com
 Data Package: Standard Level III Level IV
 Chromatograms with final report Electronic Data Deliverables (EDDs)

Data Usability Summary Report

Project Name: Custom Plywood

Project Description: Groundwater Samples

Sample Date(s): February and August 2024

Analytical Laboratory: OnSite Environmental, Inc. – Redmond, WA

Validation Performed by: Eric Hitchens

Validation Reviewed by: Gabrielle Davis

Validation Date: 3 September 2024

Haley & Aldrich, Inc. prepared this Data Usability Summary Report (DUSR) to summarize the review and validation of the analytical results for Sample Delivery Group(s) (SDGs) listed. This DUSR is organized into the following sections:

- 1. Sample Delivery Group Numbers**
 - 2. Precision and Accuracy [for SDG(s) above]**
 - 3. Explanations**
 - 4. Glossary**
 - 5. Abbreviations**
 - 6. Qualifiers**
- References**

This data validation and usability assessment was performed per the guidance and requirements established by the United States Environmental Protection Agency (USEPA) using the following reference materials:

- National Functional Guidelines (NFG) for Inorganic Data Review.
- National Functional Guidelines (NFG) for Organic Data Review.

Data reported in this sampling event were reported to the laboratory reporting limit (RL).

Sample data were qualified in accordance with the laboratory's standard operating procedures (SOPs). The results presented in each laboratory report were found to be compliant with the data quality objectives (DQOs) for the project and are therefore usable; any exceptions are noted in the following pages.

1. Sample Delivery Group Numbers

1.1 SAMPLE MANAGEMENT

This DUSR summarizes the review of SDG numbers:

- 2408-148; and
- 2402-047.

Samples were collected, preserved, and shipped following standard chain of custody (COC) protocols. Samples were also received appropriately, identified correctly, and analyzed according to the COC.

Issues noted with sample management are listed below:

- Both reports initially reported TPH analysis with silica gel treatment when it was not requested; revisions were issued to remove the unnecessary analysis from the final reports.

Analyses were performed on the following samples:

Sample ID	Sample Type	Lab ID	Sample Date	Matrix	Methods
MW-1-20240201	N	2402-047-01	02/01/2024	WG	A, B, C, D
MW-3-20240201	N	2402-047-02	02/01/2024	WG	A, B, C, D
MW-5-20240202	N	2402-047-03	02/02/2024	WG	A, B, C, D
MW-6-20240202	N	2402-047-04	02/02/2024	WG	A, B, C, D
MW-4-20240202	N	2402-047-05	02/02/2024	WG	A, B, C, D
MW-2-20240202	N	2402-047-06	02/02/2024	WG	A, B, C, D
MW-4-20240812	N	2408-148-01	08/12/2024	WG	A, B, C, D
MW-3-20240812	N	2408-148-02	08/12/2024	WG	A, B, C, D
MW-2-20240812	N	2408-148-03	08/12/2024	WG	A, B, C, D
MW-5-20240812	N	2408-148-04	08/12/2024	WG	A, B, C, D
MW-1-20240812	N	2408-148-05	08/12/2024	WG	A, B, C, D
MW-6-20240812	N	2408-148-06	08/12/2024	WG	A, B, C, D

Method Holding Times			
A.	E200.8	Metals (by Mass Spectrometer)	180 days for liquid, preserved
B.	NWTPH-DX	Total Petroleum Hydrocarbons (TPH) Diesel Range Organics	14 days for liquid, preserved 7 days for liquid, unpreserved
C.	SW7470A	Mercury (in Liquids)	28 days extraction for liquid, preserved
D.	SW8270ESIM	Polycyclic Aromatic Hydrocarbons (PAHs)	7 days extraction / 40 days analysis for liquid, unpreserved

1.2 HOLDING TIMES/PRESERVATION

The samples arrived at the laboratory at the proper temperature and were prepared and analyzed within the holding time and preservation criteria specified per method protocol, with the following exceptions:

- SDG 2402-047: Mercury was analyzed outside of analytical holding time; qualify data UJ.

1.3 REPORTING LIMITS AND SAMPLE DILUTIONS

All sample dilutions were reviewed and found to be justified.

1.4 SURROGATE RECOVERY COMPLIANCE

[Refer to Section E 1.2.](#) The percent recovery (%R) for each surrogate compound added to each project sample were determined to be within the laboratory-specified quality control (QC) limits, with the following exceptions:

SDG	Sample ID	Surrogate	Dilution	%R	Qualification
2408-148	MW-3	o-Terphenyl	1X	33	J- target compounds
2408-148	MW-1	Terphenyl-d14	1X	155	None, sample is ND

** o-Terphenyl targets the following compounds: Diesel-Range Organics, Lube Oil-Range Organics*

1.5 LABORATORY CONTROL SAMPLES

[Refer to Section E 1.3.](#) Compounds associated with the laboratory control sample/laboratory control sample duplicate (LCS/LCSD) analyses associated with client samples exhibited recoveries and relative percent differences (RPDs) within the specified limits.

1.6 MATRIX SPIKE SAMPLES

[Refer to Section E 1.4.](#) The sample(s) below were used for matrix spike/matrix spike duplicate (MS/MSD):

Lab Sample Number	Matrix Spike/Matrix Spike Duplicate Sample Client ID	Method(s)
MW-1-20240201MS/MSD	MW-1-20240201	USEPA 7470A

The MS/MSD recoveries and the relative percent difference (RPD) between the MS and MSD results were within the specified limits.

1.7 BLANK SAMPLE ANALYSIS

[Refer to Section E 1.5.](#) Method blank samples had no detections, indicating that no contamination from laboratory activities occurred.

1.8 DUPLICATE SAMPLE ANALYSIS

[Refer to Section E 1.6.](#) The following sample(s) were used for laboratory duplicate analysis and the RPDs were all below 20 percent (or the absolute difference rule was satisfied if detects were less than 5 times the RL):

Lab Sample Number	Laboratory Duplicate Sample Client ID	Method(s)
MW-1-20240201LR	MW-1-20240201	EPA 7470A

1.9 SYSTEM PERFORMANCE AND OVERALL ASSESSMENT

The results presented in this report were found to comply with the DQOs for the project and the guidelines specified by the analytical method. Based on the review of this report, the data are useable and acceptable as no data was rejected. A summary of qualifiers applied to this dataset is shown in Table 1.

2. Precision and Accuracy [for SDG(s) above]

[Refer to Section E 1.7.](#) Where required by the method, some measurement of analytical accuracy and precision was reported for each method with the site samples.

3. Explanations

The following explanations include more detailed information regarding each of the sections in the DUSR above. Not all sections in the Explanations are represented:

- E 1.2 Surrogate Recovery Compliance
 - Surrogates, also known as system monitoring compounds, are compounds added to each sample prior to sample preparation to determine the efficiency of the extraction procedure by evaluating the percent recovery (%R) of the compounds.
- E 1.3 Laboratory Control Samples
 - The laboratory control sample/laboratory control sample duplicate (LCS/LCSD) analyses are used to assess the precision and accuracy of the analytical method independent of matrix interferences.
- E 1.4 Matrix Spike Samples
 - Matrix spike/matrix spike duplicate (MS/MSD) data are used to assess the precision and accuracy of the analytical method and evaluate the effects of the sample matrix on the sample preparation procedures and measurement methodologies.
 - For inorganic methods, when a matrix spike recovery falls outside of the control limits and the sample result is less than four times the spike added, a post-digestion spike (PDS) is performed.
- E 1.5 Blank Sample Analysis
 - Method blanks are prepared by the analytical laboratory and analyzed concurrently with the project samples to assess possible laboratory contamination.
- E 1.6 Laboratory Duplicate Sample Analysis
 - The laboratory duplicate sample analysis is used by the laboratory at the time of the analysis to demonstrate acceptable method precision. The RPD or absolute difference was evaluated for each duplicate sample pair to monitor the reproducibility of the data.
- E 1.7 Precision and Accuracy
 - Precision measures the reproducibility of repetitive measurements. In a laboratory environment, this will be measured by determining the relative percent difference (RPD) found between a primary and a duplicate sample. This can be an LCS/LCSD pair, a MS/MSD pair, a laboratory duplicate performed on a site sample, or a field duplicate collected and analyzed concurrently with a site sample.
 - Accuracy is a statistical measurement of the correctness of a measured value and includes components of random error (variability caused by imprecision) and systematic error. In a laboratory environment, this will be measured by determining the percent recovery (%R) of certain spiked compounds. This can be assessed using LCS, blank spike (BS), MS, and/or surrogate recoveries.

4. Glossary

Not all of the following symbols, acronyms, or qualifiers occur in this document.

- Sample Types:
 - EB Equipment Blank Sample
 - FB Field Blank Sample
 - FD Field Duplicate Sample
 - N Primary Sample
 - TB Trip Blank Sample
- Units:
 - $\mu\text{g}/\text{kg}$ micrograms per kilogram
 - $\mu\text{g}/\text{L}$ micrograms per liter
 - $\mu\text{g}/\text{m}^3$ micrograms per cubic meter
 - mg/kg milligrams per kilogram
 - mg/L milligrams per liter
 - ppb v/v parts per billion volume/volume
 - pCi/L picocuries per liter
 - pg/g picograms per gram
 - pg/L picograms per liter
- Matrices:
 - AA Ambient Air
 - GS Soil Gas
 - GW/WG Groundwater
 - QW Water Quality
 - IA Indoor Air
 - SE Sediment
 - SO Soil
 - SSV Sub-slab Vapor
 - WQ Water Quality control matrix
 - WS Surface Water
- Table Footnotes:
 - NA Not applicable
 - ND Non-detect
 - NR Not reported
- Common Symbols:
 - % percent
 - < less than
 - \leq less than or equal to
 - > greater than
 - \geq greater than or equal to
 - = equal
 - $^{\circ}\text{C}$ degrees Celsius
 - \pm plus or minus
 - \sim approximately
 - x times (multiplier)

- Fractions:
 - N Normal (method cannot be filtered)
 - D Dissolved (filtered)
 - T Total (unfiltered)

5. Abbreviations

%D	Percent Difference	MDL	Laboratory Method Detection Limit
%R	Percent Recovery	MS/MSD	Matrix Spike/Matrix Spike Duplicate
%RSD	Percent Relative Standard Deviation	NA	not applicable
%v/v	Percent volume by volume	ND	Non-Detect
2s	2 sigma	NFG	National Functional Guidelines
4,4-DDT	4 4-dichlorodiphenyltrichloroethane	NH ₃	Ammonia
Abs Diff	Absolute Difference	NYSDEC	New York State Department of Environmental Conservation
amu	atomic mass unit	PAH	Polycyclic Aromatic Hydrocarbon
BPJ	Best Professional Judgement	PCB	Polychlorinated Biphenyl
BS	Blank Spike	PDS	Post-Digestion Spike
CCB	Continuing Calibration Blank	PEM	Performance Evaluation Mixture
CCV	Continuing Calibration Verification	PFAS	Per- and Polyfluoroalkyl Substances
CCVL	Continuing Calibration Verification Low	PFBA	Perfluorbutanoic Acid
COC	Chain of Custody	PFD	Perfluorodecalin
COM	Combined Isotope Calculation	PFOA	Perfluorooctanoic Acid
Cr (VI)	Hexavalent Chromium	PFOS	Perfluorooctane sulfonate
CRI	Collision Reaction Interface	PFPeA	Perfluoropentanoic Acid
DoD	Department of Defense	QAPP	Quality Assurance Project Plan
DQO	data quality objective	QC	Quality Control
DUSR	Data Usability Summary Report	QSM	Quality Systems Manual
EIS	Extraction Internal Standard	R ²	R-squared value
EMPC	Estimated Maximum Possible Concentration	Ra-226	Radium-226
FBK	Field Blank Contamination	Ra-228	Radium-228
FDP	Field Duplicate	RESC	Resolution Check Measure
GC	Gas Chromatograph	RL	Laboratory Reporting Limit
GC/MS	Gas Chromatography/Mass Spectrometry	RPD	Relative Percent Difference
GPC	Gel Permeation Chromatography	RRF	Relative Response Factor
H ₂	Hydrogen gas	RT	Retention Time
HCl	Hydrochloric Acid	SAP	Sampling Analysis Plan
ICAL	Initial Calibration	SDG	Sample Delivery Group
ICB	Initial Calibration Blank	SIM	Selected ion monitoring
ICP/MS	Inductively Coupled Plasma/Mass Spectrometry	SOP	Standard Operating Procedure
ICV	Initial Calibration Verification	SPE	Solid-Phase Extraction
ICVL	Initial Calibration Verification Low	SVOC	Semi-Volatile Organic Compound
IPA	Isopropyl Alcohol	TCLP	Toxicity Characteristic Leaching Procedure
LC	Laboratory Control	TIC	Tentatively Identified Compound
LCS/LCSD	Laboratory Control Sample/Laboratory Control Sample Duplicate	TKN	Total Kjeldahl Nitrogen
MBK	Method Blank Contamination	TPH	Total Petroleum Hydrocarbon
MDC	Minimum Detectable Concentration	TPU	Total Propagated Uncertainty
		USEPA	U.S. Environmental Protection Agency
		VOC	Volatile Organic Compound
		WP	Work Plan

6. Qualifiers

The qualifiers below are from the USEPA National Functional Guidelines and the data in the DUSR may contain these qualifiers:

- Concentration (C) Qualifiers:
 - U The compound was analyzed for but not detected. The associated value is either the compound quantitation limit if not detected by the analytical instrument or could be the reported or blank concentration if qualified by blank contamination. This can also be displayed as less than the associated compound quantitation limit (<RL or <MDL), or “ND”.
 - B The compound was found in the sample and its associated blank. Its presence in the sample may be suspect.
- Quantitation (Q) Qualifiers:
 - E The compound was quantitated above the calibration range.
 - D The concentration is based on a diluted sample analysis.
- Validation Qualifiers:
 - J The compound was positively identified; however, the associated numerical value is an estimated concentration only.
 - J+ The result is an estimated quantity, but the result may be biased high.
 - J- The result is an estimated quantity, but the result may be biased low.
 - J/UJ as listed in exception tables J applies to detected data and UJ applies to non-detected data as reported by the laboratory.
 - UJ The compound was not detected. The reported sample quantitation limit is approximate.
 - NJ The analysis indicated the presence of a compound for which there is presumptive evidence to make a tentative identification; the associated numerical value is an estimated concentration only.
 - R The sample results were rejected as unusable; the compound may or may not be present in the sample.
 - S Result is suspect. See DUSR for details.

References

1. United States Environmental Protection Agency, 2020a. National Functional Guidelines for Inorganic Superfund Methods Data Review. EPA-542-R-20-006. November.
2. United States Environmental Protection Agency, 2020b. National Functional Guidelines for Organic Superfund Methods Data Review. EPA-540-R-20-005. November.

Attachments:

Table 1 – System Performance Summary

TABLE

TABLE 1
SYSTEM PERFORMANCE SUMMARY
CUSTOM PLYWOOD

SDG	Method	Sample ID	Lab ID	Analyte	Fraction	Reportable Result	Reported Result	Validated Result	Reason for Qualifier
2408-148	NWTPH-DX	MW-3-20240812	2408-148-02	Diesel Range Organics	Not Applicable	Yes	0.15	0.15 J-	SUR
2408-148	NWTPH-DX	MW-3-20240812	2408-148-02	Petroleum Hydrocarbons as Lube Oil	Not Applicable	Yes	0.32	0.32 J-	SUR

Notes:

SUR = Surrogate percent recovery outside the specified limits.

J- = The result is an estimated quantity, but the result may be biased low.