

GALLOWAY ENVIRONMENTAL, INC 15600 NE 8th Street, Suite B1, 617 (425) 894-8607 Bellevue, WA 98008 Dylan@GallowayEnvironmental.com

June 25, 2021

Tamarah Hancock Scarsella Bros. Inc. PO Box 68697 Seattle, Washington 98168

Emailed to: Tamarah.k@scarsellabros.com and Jenifer.m@scarsellabros.com

SUBJECT: SUMMARY STATUS REPORT — GROUNDWATER MONITORING WELL SAMPLE RESULTS AT THE FIRWOOD PIT PROPERTY IN EDGEWOOD, WA SEPA PROJECT #1808 – CLEAR AND GRADE PERMIT #3492

Dear Ms. Hancock:

This letter report presents a summary of Galloway Environmental, Inc.'s (GEI's) findings from the groundwater monitoring event at the Firwood Pit property for May 2021.

The scope of work for this quarterly groundwater monitoring report is generally based on the City of Edgewood's Final Conditions for Firwood Pit Reclamation – Original SEPA project #1808, Clear and Grade Permit #3492 which was updated on November 26, 2018 and your direction. This report generally includes: 1) chemical analytical results of water samples, and 2) physical properties of groundwater in the monitoring wells.

INTRODUCTION

The Firwood Mine was a sand and gravel surface mine that was exhausted of its aggregate resource before March 1, 2000 when it was assigned and leased to Scarsella Bros., Inc. by the Tim Corliss and Son Company and is now in the reclamation process. The mine is located in the general area east of Freeman Road, adjacent to the west side of 90th Avenue East, south of 33rd Street East, and northeast of Simons Creek. The Site is in the City of Edgewood, Pierce County, Washington.

GROUNDWATER SAMPLE COLLECTION SUMMARY

On May 28, 2021, GEI collected groundwater samples from monitoring wells MW-1, MW-2, MW-3b, and MW-4. Prior to sample collection, GEI gauged and purged the wells to attain groundwater samples that were representative of the site.

During the purging process for the monitoring wells, water quality parameters were measured using a multiparameter water quality meter (model YSI 556 MPS) to measure pH, temperature, conductivity, and dissolved oxygen (DO). The YSI 556 MPS meter was calibrated prior to well purging using a 3-point pH calibration process (pH valued at 4.02, 7.02, and 10.04) and a 3-point conductivity process (conductivity valued at 84 micrograms per centimeter [μ g/cm], 1,413 μ g/cm, and 12,880 μ g/cm). Additionally, to quantify turbidity throughout the purging process, GEI utilized a turbidity meter (model Lamotte #2020T).

Sample collection was initiated upon attaining stabilized water quality parameters, including pH within 0.1 SU, conductivity within 5%, and temperature within 0.1 degrees Celsius (°C) for at least three consecutive readings.

The pH measured at the conclusion of the purging process ranged from 6.91 (MW-4) to 7.40 (MW-2). The conductivity measured at the conclusion of the purging process ranged from 0.400 milliSiemens per

Firwood Pit Groundwater Monitoring Report June 25, 2021 Page 2



centimeter (mS/cm) (MW-2) to 0.625 mS/cm (MW-3b). The DO measured at the conclusion of the purging process ranged from 2.62 milligrams per liter (mg/L) (MW-3b) to 6.82 mg/L (MW-2). Turbidity measured at the conclusion of the purging process ranged from 1.19 Nephelometric Turbidity Units (NTU) (MW-2) to 3.75 NTU at MW-4. The water color was observed to be clear in all wells at the conclusion of the purging process. GEI purged a minimum of three well volumes from each well, or until water quality parameters were stabilized.

GEI purged a total of 5.5 gallons from MW-1, 3.5 gallons from MW-2, 5.5 gallons from MW-3b, and 3.5 gallons from MW-4.

GROUNDWATER SAMPLE ANALYSES SUMMARY

All samples were submitted to OnSite Environmental, Inc., located at 14648 NE 95th St., Redmond, Washington (OnSite Environmental) for analyses of petroleum hydrocarbons using the Northwest Total Petroleum Hydrocarbons as Hydrocarbon Identification (NWTPH-HCID) method, polycyclic aromatic hydrocarbons (PAHs) using the United States Environmental Protection Agency (USEPA) method 8270E/SIM, and dissolved arsenic using the USEPA method 200.8.

Because the initial NWTPH-HCID laboratory analyses detected the presence of total petroleum hydrocarbons as lube oil (TPH-Oil) in the sample that was collected from MW-1, the sample was further analyzed to quantify concentrations of TPH-Oil in the sample using the Northwest Total Petroleum Hydrocarbons as Diesel Extended (NWTPH-Dx) method.

Below is a summary of the laboratory analytical results. Attachment 1 includes the laboratory analytical report for this monitoring event.

Laboratory analyses resulted in the detection of TPH-Oil (as lube oil range organics) in the sample that was collected from MW-1 at a concentration of 0.30 mg/L, below the Washington States Department of Ecology (Ecology) Model Toxics Control Act (MTCA) Method A Cleanup Level of 0.50 mg/L.

Laboratory analyses resulted in no other detections of analytes at concentrations equal to or exceeding their respective laboratory practical quantitation limits (PQLs) in any of the samples analyzed.

As per the City of Edgewood's Final Conditions for the Clear and Grade Permit (#3492), petroleum hydrocarbon compounds and dissolved arsenic will be tested quarterly; and PAHs will be tested annually.

Should you have any questions regarding this report or if you would like to discuss our findings, please contact us at any of the addresses listed on top of this letter.

Respectfully Submitted, GALLOWAY ENVIRONMENTAL, INC.

Dylan Galloway, REA President cc: Jenifer A. Morrison, SBI

Attachment 1 Laboratory Analytical Reports



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

June 9, 2021

Dylan Galloway Galloway Environmental, Inc. 15600 NE 8th Street Suite B1, 617 Bellevue WA 98008

Re: Analytical Data for Project 28027 Laboratory Reference No. 2105-282

Dear Dylan:

Enclosed are the analytical results and associated quality control data for samples submitted on May 28, 2021.

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,

David Baumeister Project Manager

Enclosures



Case Narrative

Samples were collected on May 28, 2021 and received by the laboratory on May 28, 2021. 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.

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.



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

HYDROCARBON IDENTIFICATION NWTPH-HCID

Matrix: Water Units: mg/L (ppm)

0 (11)				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW-1					
Laboratory ID:	05-282-01					
Gasoline Range Organics	ND	0.10	NWTPH-HCID	6-1-21	6-3-21	
Diesel Range Organics	ND	0.20	NWTPH-HCID	6-1-21	6-3-21	
Lube Oil Range Organics	Detected	0.20	NWTPH-HCID	6-1-21	6-3-21	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	90	50-150				
Client ID:	MW-2					
Laboratory ID:	05-282-02					
Gasoline Range Organics	ND	0.10	NWTPH-HCID	6-1-21	6-3-21	
Diesel Range Organics	ND	0.20	NWTPH-HCID	6-1-21	6-3-21	
Lube Oil Range Organics	ND	0.20	NWTPH-HCID	6-1-21	6-3-21	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	91	50-150				
Client ID:	MW-3b					
Laboratory ID:	05-282-03					
Gasoline Range Organics	ND	0.10	NWTPH-HCID	6-1-21	6-2-21	
Diesel Range Organics	ND	0.20	NWTPH-HCID	6-1-21	6-2-21	
Lube Oil Range Organics	ND	0.20	NWTPH-HCID	6-1-21	6-2-21	
Surrogate:	Percent Recovery	Control Limits		-	-	
o-Terphenyl	88	50-150				
Client ID:	MW-4					
Laboratory ID:	05-282-04					
Gasoline Range Organics	ND	0.10	NWTPH-HCID	6-1-21	6-3-21	
Diesel Range Organics	ND	0.20	NWTPH-HCID	6-1-21	6-3-21	
Lube Oil Range Organics	ND	0.20	NWTPH-HCID	6-1-21	6-3-21	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	85	50-150				



3

HYDROCARBON IDENTIFICATION NWTPH-HCID QUALITY CONTROL

Matrix: Water Units: mg/L (ppm)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0601W1					
Gasoline Range Organics	ND	0.10	NWTPH-HCID	6-1-21	6-3-21	
Diesel Range Organics	ND	0.20	NWTPH-HCID	6-1-21	6-3-21	
Lube Oil Range Organics	ND	0.20	NWTPH-HCID	6-1-21	6-3-21	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	95	50-150				



4

PAHs EPA 8270E/SIM

Matrix: Water Units: ug/L

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW-1					
Laboratory ID:	05-282-01					
Naphthalene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
2-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
1-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Acenaphthylene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Acenaphthene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Fluorene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Phenanthrene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Anthracene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Fluoranthene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Pyrene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Benzo[a]anthracene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Chrysene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Benzo[b]fluoranthene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Benzo[a]pyrene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	67	25 - 106				
Pyrene-d10	85	28 - 104				
Terphenyl-d14	88	40 - 139				



5

PAHs EPA 8270E/SIM

Matrix: Water Units: ug/L

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW-2					
Laboratory ID:	05-282-02					
Naphthalene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
2-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
1-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Acenaphthylene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Acenaphthene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Fluorene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Phenanthrene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Anthracene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Fluoranthene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Pyrene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Benzo[a]anthracene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Chrysene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Benzo[b]fluoranthene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Benzo[a]pyrene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	48	25 - 106				
Pyrene-d10	80	28 - 104				
Terphenyl-d14	82	40 - 139				



PAHs EPA 8270E/SIM

Matrix: Water Units: ug/L

5				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW-3b					
Laboratory ID:	05-282-03					
Naphthalene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
2-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
1-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
Acenaphthylene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
Acenaphthene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
Fluorene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
Phenanthrene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
Anthracene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
Fluoranthene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
Pyrene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
Benzo[a]anthracene	ND	0.010	EPA 8270E/SIM	6-2-21	6-3-21	
Chrysene	ND	0.010	EPA 8270E/SIM	6-2-21	6-3-21	
Benzo[b]fluoranthene	ND	0.010	EPA 8270E/SIM	6-2-21	6-3-21	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270E/SIM	6-2-21	6-3-21	
Benzo[a]pyrene	ND	0.010	EPA 8270E/SIM	6-2-21	6-3-21	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270E/SIM	6-2-21	6-3-21	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270E/SIM	6-2-21	6-3-21	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270E/SIM	6-2-21	6-3-21	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	40	25 - 106				
Pyrene-d10	68	28 - 104				
Terphenyl-d14	74	40 - 139				



PAHs EPA 8270E/SIM

Matrix: Water Units: ug/L

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW-4					
Laboratory ID:	05-282-04					
Naphthalene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
2-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
1-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
Acenaphthylene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
Acenaphthene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
Fluorene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
Phenanthrene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
Anthracene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
Fluoranthene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
Pyrene	ND	0.10	EPA 8270E/SIM	6-2-21	6-3-21	
Benzo[a]anthracene	ND	0.010	EPA 8270E/SIM	6-2-21	6-3-21	
Chrysene	ND	0.010	EPA 8270E/SIM	6-2-21	6-3-21	
Benzo[b]fluoranthene	ND	0.010	EPA 8270E/SIM	6-2-21	6-3-21	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270E/SIM	6-2-21	6-3-21	
Benzo[a]pyrene	ND	0.010	EPA 8270E/SIM	6-2-21	6-3-21	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270E/SIM	6-2-21	6-3-21	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270E/SIM	6-2-21	6-3-21	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270E/SIM	6-2-21	6-3-21	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	59	25 - 106				
Pyrene-d10	65	28 - 104				
Terphenyl-d14	82	40 - 139				



PAHs EPA 8270E/SIM QUALITY CONTROL

. . .

Matrix: Water Units: ug/L

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
METHOD BLANK						
Laboratory ID:	MB0602W1					
Naphthalene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
2-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
1-Methylnaphthalene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Acenaphthylene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Acenaphthene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Fluorene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Phenanthrene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Anthracene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Fluoranthene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Pyrene	ND	0.10	EPA 8270E/SIM	6-2-21	6-2-21	
Benzo[a]anthracene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Chrysene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Benzo[b]fluoranthene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Benzo[a]pyrene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270E/SIM	6-2-21	6-2-21	
Surrogate:	Percent Recovery	Control Limits				
2-Fluorobiphenyl	65	25 - 106				
Pyrene-d10	85	28 - 104				
Terphenyl-d14	109	40 - 139				



9

PAHs EPA 8270E/SIM QUALITY CONTROL

Matrix: Water Units: ug/L

AnalyteResultSpike LevelRecoveryLimitsRPDLimSPIKE BLANKSLaboratory ID:SB0602W1SBSBDSBSBDSBSBDNaphthalene0.2950.2900.5000.500595829 - 96238Acenaphthylene0.3610.3630.5000.500727342 - 101128Acenaphthene0.3380.3230.5000.500737848 - 101621Phenanthrene0.3650.3890.5000.500778452 - 104920Anthracene0.3620.4100.5000.500728250 - 1061220Fluoranthene0.4350.4500.5000.500728250 - 1061220Fluoranthene0.4350.4500.5000.500728250 - 1061220Fluoranthene0.4350.4500.5000.500879056 - 113320Pyrene0.5540.4410.5000.500889460 - 131620Chrysene0.4390.4870.5000.500889762 - 1201020Benzo[a]anthracene0.4420.4690.5000.500889963 - 1231020Benzo[b]fluoranthene0.4470.4930.5000.500889461 - 123620 <t< th=""><th></th></t<>	
Laboratory ID: SB0602W1 SB SBD SB SBD SB SBD Naphthalene 0.295 0.290 0.500 0.500 59 58 29 - 96 2 38 Acenaphthylene 0.361 0.363 0.500 0.500 72 73 42 - 101 1 28 Acenaphthene 0.338 0.323 0.500 0.500 68 65 37 - 104 5 31 Fluorene 0.365 0.389 0.500 0.500 73 78 48 - 101 6 21 Phenanthrene 0.385 0.421 0.500 0.500 77 84 52 - 104 9 20 Anthracene 0.362 0.410 0.500 0.500 72 82 50 - 106 12 20 Fluoranthene 0.435 0.450 0.500 0.500 87 90 56 - 113 3 20 Pyrene 0.554 0.441 0.500	Flags
SB SBD SB SBD SB SBD SB SBD Naphthalene 0.295 0.290 0.500 0.500 59 58 29 - 96 2 38 Acenaphthylene 0.361 0.363 0.500 0.500 72 73 42 - 101 1 28 Acenaphthene 0.338 0.323 0.500 0.500 68 65 37 - 104 5 31 Fluorene 0.365 0.389 0.500 0.500 73 78 48 - 101 6 21 Phenanthrene 0.362 0.410 0.500 0.500 77 84 52 - 104 9 20 Anthracene 0.362 0.410 0.500 0.500 72 82 50 - 106 12 20 Fluoranthene 0.435 0.450 0.500 0.500 87 90 56 - 113 3 20 Pyrene 0.554 0.441 0.500 0.500 88	
Naphthalene0.2950.2900.5000.500595829 - 96238Acenaphthylene0.3610.3630.5000.500727342 - 101128Acenaphthene0.3380.3230.5000.500686537 - 104531Fluorene0.3650.3890.5000.500737848 - 101621Phenanthrene0.3850.4210.5000.500778452 - 104920Anthracene0.3620.4100.5000.500728250 - 1061220Fluoranthene0.4350.4500.5000.500879056 - 113320Pyrene0.5540.4410.5000.5001118855 - 1232327Benzo[a]anthracene0.4420.4690.5000.500889460 - 131620Chrysene0.4390.4870.5000.500889762 - 1201020Benzo[b]fluoranthene0.4470.4930.5000.500899963 - 1231020Benzo[j,k)fluoranthene0.4420.4690.5000.500889461 - 123620Indeno(1,2,3-c,d)pyrene0.4320.4640.5000.500869360 - 125720	
Acenaphthylene0.3610.3630.5000.500727342 - 101128Acenaphthene0.3380.3230.5000.500686537 - 104531Fluorene0.3650.3890.5000.500737848 - 101621Phenanthrene0.3850.4210.5000.500778452 - 104920Anthracene0.3620.4100.5000.500728250 - 1061220Fluoranthene0.4350.4500.5000.500879056 - 113320Pyrene0.5540.4410.5000.5001118855 - 1232327Benzo[a]anthracene0.4420.4690.5000.500889460 - 131620Chrysene0.4390.4870.5000.500889762 - 1201020Benzo[b]fluoranthene0.4470.4930.5000.500899963 - 1231020Benzo[j,k)fluoranthene0.4420.4690.5000.500889461 - 123620Indeno(1,2,3-c,d)pyrene0.4320.4640.5000.500869360 - 125720	
Acenaphthene0.3380.3230.5000.500686537 - 104531Fluorene0.3650.3890.5000.500737848 - 101621Phenanthrene0.3850.4210.5000.500778452 - 104920Anthracene0.3620.4100.5000.500728250 - 1061220Fluoranthene0.4350.4500.5000.500879056 - 113320Pyrene0.5540.4410.5000.5001118855 - 1232327Benzo[a]anthracene0.4320.4690.5000.500889460 - 131620Chrysene0.4390.4870.5000.500889762 - 1201020Benzo[b]fluoranthene0.4470.4930.5000.500899963 - 1231020Benzo[j,k)fluoranthene0.4420.4690.5000.500889461 - 123620Indeno(1,2,3-c,d)pyrene0.4320.4640.5000.500869360 - 125720	
Fluorene0.3650.3890.5000.500737848 - 101621Phenanthrene0.3850.4210.5000.500778452 - 104920Anthracene0.3620.4100.5000.500728250 - 1061220Fluoranthene0.4350.4500.5000.500879056 - 113320Pyrene0.5540.4410.5000.5001118855 - 1232327Benzo[a]anthracene0.4420.4690.5000.500889460 - 131620Chrysene0.4390.4870.5000.500889762 - 1201020Benzo[b]fluoranthene0.4470.4930.5000.500899963 - 1231020Benzo[j,k)fluoranthene0.4420.4690.5000.500889461 - 123620Indeno(1,2,3-c,d)pyrene0.4320.4640.5000.500869360 - 125720	
Phenanthrene0.3850.4210.5000.500778452 - 104920Anthracene0.3620.4100.5000.500728250 - 1061220Fluoranthene0.4350.4500.5000.500879056 - 113320Pyrene0.5540.4410.5000.5001118855 - 1232327Benzo[a]anthracene0.4420.4690.5000.500889460 - 131620Chrysene0.4390.4870.5000.500889762 - 1201020Benzo[b]fluoranthene0.4470.4930.5000.500899963 - 1231020Benzo(j,k)fluoranthene0.4650.4970.5000.500889461 - 123620Indeno(1,2,3-c,d)pyrene0.4320.4640.5000.500869360 - 125720	
Anthracene0.3620.4100.5000.500728250 - 1061220Fluoranthene0.4350.4500.5000.500879056 - 113320Pyrene0.5540.4410.5000.5001118855 - 1232327Benzo[a]anthracene0.4420.4690.5000.500889460 - 131620Chrysene0.4390.4870.5000.500889762 - 1201020Benzo[b]fluoranthene0.4470.4930.5000.500899963 - 1231020Benzo(j,k)fluoranthene0.4650.4970.5000.500889461 - 123620Indeno(1,2,3-c,d)pyrene0.4320.4640.5000.500869360 - 125720	
Fluoranthene0.4350.4500.5000.500879056 - 113320Pyrene0.5540.4410.5000.5001118855 - 1232327Benzo[a]anthracene0.4420.4690.5000.500889460 - 131620Chrysene0.4390.4870.5000.500889762 - 1201020Benzo[b]fluoranthene0.4470.4930.5000.500899963 - 1231020Benzo(j,k)fluoranthene0.4650.4970.5000.500939960 - 127720Benzo[a]pyrene0.4420.4690.5000.500889461 - 123620Indeno(1,2,3-c,d)pyrene0.4640.5000.500869360 - 125720	
Pyrene0.5540.4410.5000.5001118855 - 1232327Benzo[a]anthracene0.4420.4690.5000.500889460 - 131620Chrysene0.4390.4870.5000.500889762 - 1201020Benzo[b]fluoranthene0.4470.4930.5000.500899963 - 1231020Benzo(j,k)fluoranthene0.4650.4970.5000.500939960 - 127720Benzo[a]pyrene0.4420.4690.5000.500889461 - 123620Indeno(1,2,3-c,d)pyrene0.4320.4640.5000.500869360 - 125720	
Benzo[a]anthracene0.4420.4690.5000.500889460 - 131620Chrysene0.4390.4870.5000.500889762 - 1201020Benzo[b]fluoranthene0.4470.4930.5000.500899963 - 1231020Benzo(j,k)fluoranthene0.4650.4970.5000.500939960 - 127720Benzo[a]pyrene0.4420.4690.5000.500889461 - 123620Indeno(1,2,3-c,d)pyrene0.4320.4640.5000.500869360 - 125720	
Chrysene0.4390.4870.5000.500889762 - 1201020Benzo[b]fluoranthene0.4470.4930.5000.500899963 - 1231020Benzo(j,k)fluoranthene0.4650.4970.5000.500939960 - 127720Benzo[a]pyrene0.4420.4690.5000.500889461 - 123620Indeno(1,2,3-c,d)pyrene0.4320.4640.5000.500869360 - 125720	
Benzo[b]fluoranthene0.4470.4930.5000.500899963 - 1231020Benzo(j,k)fluoranthene0.4650.4970.5000.500939960 - 127720Benzo[a]pyrene0.4420.4690.5000.500889461 - 123620Indeno(1,2,3-c,d)pyrene0.4320.4640.5000.500869360 - 125720	
Benzo(j,k)fluoranthene0.4650.4970.5000.500939960 - 127720Benzo[a]pyrene0.4420.4690.5000.500889461 - 123620Indeno(1,2,3-c,d)pyrene0.4320.4640.5000.500869360 - 125720	
Benzo[a]pyrene 0.442 0.469 0.500 0.500 88 94 61 - 123 6 20 Indeno(1,2,3-c,d)pyrene 0.432 0.464 0.500 0.500 86 93 60 - 125 7 20	
Indeno(1,2,3-c,d)pyrene 0.432 0.464 0.500 0.500 86 93 60 - 125 7 20	
Dibenz[a,h]anthracene 0.441 0.475 0.500 0.500 88 95 61 - 124 7 20	
Benzo[g,h,i]perylene 0.444 0.484 0.500 0.500 89 97 59 - 122 9 20	
Surrogate:	
2-Fluorobiphenyl 59 55 25 - 106	
Pyrene-d10 83 80 28 - 104	
Terphenyl-d14 89 84 40 - 139	



DISSOLVED ARSENIC EPA 200.8

Matrix: Water Units: ug/L (ppb)

				Date	Date	
Analyte	Result	PQL	Method	Prepared	Analyzed	Flags
Client ID:	MW-1					
Laboratory ID:	05-282-01					
Arsenic	ND	3.0	EPA 200.8	5-28-21	6-2-21	
Client ID:	MW-2					
Laboratory ID:	05-282-02					
Arsenic	ND	3.0	EPA 200.8	5-28-21	6-2-21	
Client ID:	MW-3b					
Laboratory ID:	05-282-03					
Arsenic	ND	3.0	EPA 200.8	5-28-21	6-2-21	
Client ID:	MW-4					
Laboratory ID:	05-282-04					
Arsenic	ND	3.0	EPA 200.8	5-28-21	6-2-21	



DISSOLVED ARSENIC EPA 200.8 QUALITY CONTROL

Matrix: Water Units: ug/L (ppb)

								Date	Date	e	
Analyte		Result		PQL N		Method Prepared		Analyzed		Flags	
METHOD BLANK											
Laboratory ID:		MB0528F1									
Arsenic		ND		3.0	EP	A 200.8	1	5-28-21	6-2-2	21	
					•	_		_			
					Source	Perc	ent	Recovery		RPD	
Analyte	Res	sult	Spike	Level	Result	Reco	very	Limits	RPD	Limit	Flags
DUPLICATE											
Laboratory ID:	05-28	32-01									
	ORIG	DUP									
Arsenic	ND	ND	NA	NA		N	4	NA	NA	20	
MATRIX SPIKES											
Laboratory ID:	05-28	32-01									
	MS	MSD	MS	MSD		MS	MSD				
Arsenic	83.4	84.2	80.0	80.0	ND	104	105	75-125	1	20	



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:	05-282-01					
Diesel Range Organics	ND	0.20	NWTPH-Dx	6-1-21	6-3-21	
Lube Oil Range Organics	0.30	0.20	NWTPH-Dx	6-1-21	6-3-21	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	90	50-150				



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

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:	MB0601W1					
Diesel Range Organics	ND	0.20	NWTPH-Dx	6-1-21	6-3-21	
Lube Oil Range Organics	ND	0.20	NWTPH-Dx	6-1-21	6-3-21	
Surrogate:	Percent Recovery	Control Limits				
o-Terphenyl	95	50-150				
		S	ource Percen	t Recovery	RPE)

Analyte	Re	sult	Spike Level		Result	Recovery		Limits	RPD	Limit	Flags	
DUPLICATE												
Laboratory ID:	SB06	01W1										
	ORIG	DUP										
Diesel Fuel #2	0.456	0.431	NA	NA		NA		NA	6	NA		
Surrogate:												
o-Terphenyl						105	98	50-150				





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.
- 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.

Ζ-

ND - Not Detected at PQL

- PQL Practical Quantitation Limit
- **RPD** Relative Percent Difference



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

Reviewed/Date	Received	Relinquished	Received	Relinquished	Received	Relinquished	Signature				4 MW-4	3 MW-36	2 mw-2	1 mw-1	Lab ID Sample Identification	Sampled by: D Galloning	D mailwary	Project Name: Project Manager	Project Namon	Brief Number	Analytical Laboratory Testing Services 14648 NE 95th Street • Redmond, WA 98052 Phone: (425) 883-3881 • www.onsite-env.com	Environmental Inc.
Reviewed/Date					086	681	Company				8 1 05.51 0	8 hhzl	1 1130 1 8	\$ 12 0440 W 8	Sampled Sampled Matrix	(other)		(TPH analysis 5 Days)	2 Days 3 Days	Same Day 1 Day	(in working days) (Check One)	Chain of
					1-121 12/202	5/13/21 151	Date Time				×	×	X	X	Number of Containers NWTPH-HCID NWTPH-Gx/BTEX NWTPH-Gx NWTPH-Dx (Acid / SG Clean-up) Volatiles 8260C Halogenated Volatiles 8260C EDB EPA 8011 (Waters Only)					Laboratory Number:	Chain of Custody	
Chromatograms with final report Electronic Data Deliverables (EDDs)	Data Package: Standard Level III Level IV	Addeci 6/3/21 7 Addeci 6/3/21 7 Mardana 74									×	X	×		Semiv (with I PAHs PCBs Organ Organ Chlori Total F Total N TCLP HEM (olatiles ow-leve 8270D/ 8082A ochlorin ophosp hated A RCRA M ATCA M Metals oil and	8270D Il PAHe SIM (Ic horus ccid He letals fletals	/SIM	081B es 8270 8151A		r: 05-282	Page of