



# ANALYTICAL SUMMARY REPORT

March 01, 2022

AECOM - Honolulu  
1001 Bishop Street, Suite 1600  
Honolulu HI, 96813-3698

Work Order: B22011131 Quote ID: 5912

Project Name: CV18F0126/60571032.02.46.01

Energy Laboratories Inc Billings MT received the following 5 samples from AECOM - Honolulu on 1/19/2022 for analysis.

Lab ID	Client Sample ID	Collect Date	Received Date	Matrix	Test
B22011131-001	ERH2460 (RHMW12A)	01/17/22 17:15	01/19/2022	Ground Water	Metals Digestion by SW3010A DRO-Liquid-Liquid Extraction SW3520C Low Level PAH by 8270C SIM SW8270CSIM Separatory Funnel SW3510C Liquid-Liquid Ext. Carbon, Total Organic SW9060A Metals by ICP-MS, Dissolved SW6020 Metals by ICP-MS, Total SW6020 8260-Volatile Organic Compounds-Short List SW8260B EDB in Water by ECD SW8011 Gasoline Range Organics SW8015C Diesel Range Organics SW8015C Headspace Gas Analysis SW8015M Semi-Volatile Organic Compounds, Extended List SW8270C SW8011 Microextraction
B22011131-002	ERH2459 (Trip Blank)- 14694	01/17/22 17:15	01/19/2022	Trip Blank	8260-Volatile Organic Compounds-Short List SW8260B
B22011131-003	ERH2459 (Trip Blank)- 14733	01/17/22 17:15	01/19/2022	Trip Blank	Gasoline Range Organics SW8015C
B22011131-004	ERH2459 (Trip Blank)- 14705	01/17/22 17:15	01/19/2022	Trip Blank	EDB in Water by ECD SW8011 SW8011 Microextraction



## ANALYTICAL SUMMARY REPORT

B22011131-005    ERH2459 (Trip Blank)-    01/17/22 17:15    01/19/2022    Trip Blank    Headspace Gas Analysis  
14732    SW8015M

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The analyses presented in this report were performed by Energy Laboratories, Inc., 1120 S 27th St., Billings, MT 59101, unless otherwise noted. Any exceptions or problems with the analyses are noted in the report package. Any issues encountered during sample receipt are documented in the Work Order Receipt Checklist.

The results as reported relate only to the item(s) submitted for testing. This report shall be used or copied only in its entirety. Energy Laboratories, Inc. is not responsible for the consequences arising from the use of a partial report.

If you have any questions regarding these test results, please contact your Project Manager.

Report Approved By:



**CLIENT:** AECOM - Honolulu  
**Project:** CV18F0126/60571032.02.46.01  
**Work Order:** B22011131

**Report Date:** 3/1/2022

## CASE NARRATIVE

### General Comments:

For any question please contact your Project Manager at (406) 252-6325 or [billingspm@energylab.com](mailto:billingspm@energylab.com).

All analyses have been performed in accordance with DOD QSM Version 5.3 unless otherwise noted below. The specific methodologies used in obtaining the enclosed analytical results are indicated on the Analytical Summary Report and the Laboratory Analytical Report. The samples were received at the appropriate temperature and in accordance with the chain of custody unless otherwise noted in the Work Order Receipt Checklist.

The tests listed below are accredited and meet the requirements of DoD QSM Version 5.3 as verified by ANSI-ASQ National Accreditation Board (ANAB) certificate number ADE-2588. Exceptions to this require client authorization and records documenting this approval are attached in the Sample Management Records. Accreditation may not be offered or required for all methods and analytes reported in this package. Refer to the certificate and scope of accreditation located at <https://www.energylab.com/whyus/certifications-quality-control/> or contact your project manager.

Tests for Total Organic Carbon by SW0060A associated with analyst identified as ELI-CA were subcontracted to Energy Laboratories, PO Box 247, Casper, WY, EPA Number WY00002.

Project specific matrix quality control samples may not be reported if site specific samples were not submitted. Matrix quality control samples were performed on project samples where adequate volume was available. All quality control measures met criteria unless otherwise noted in the Analytical QC Exceptions report and in the Analysis Specific Comments below. Where available, sample management records are attached.

The Stage 4 Validation Package includes data reports for all analyses associated with the instrument calibration, quality control (QC) sample analysis, and sample analysis. All analytical data is within method specifications except as noted in the Analytical QC Exceptions report or the Analysis Specific Comments below. The analytical report identifies preparation batch and analytical run IDs associated with each result for a sample. Only the raw data associated with the parameters listed on this report should be validated.

### Analysis Specific Comments:

An Analytical QC Exceptions Report has been attached, summarizing all qualified QC results. Where qualified, an analyte exceeded quality control limits, but was not detected in the associated sample(s).



Trust our People. Trust our Data.

# Chain of Custody & Analytical Request Record – DoD Project

www.energylab.com

COC#202201-77NOI Page 1 of 1

### Account Information (Billing information)

Company/Name	AECOM	
Contact	Alethea Ramos / Margie Pascua	
Phone	808-529-7283 / 808-356-5373	
Mailing Address	1001 Bishop St., Suite 1600	
City, State, Zip	Honolulu, HI 96813	
Email	alethea.ramos@aecom.com / margie.pascua@aecom.com	
Receive Invoice	<input type="checkbox"/> Hard Copy <input checked="" type="checkbox"/> Email	Receive Report <input type="checkbox"/> Hard Copy <input checked="" type="checkbox"/> Email
Purchase Order	Quote	Bottle Order
N/A	N/A	N/A

### Report Information (if different than Account Information)

Company/Name	AECOM	
Contact	see Account information	
Phone		
Mailing Address		
City, State, Zip		
Email	USAPimaging@aecom.com	
Receive Report	<input type="checkbox"/> Hard Copy <input type="checkbox"/> Email	
Special Report/Formats:	<input checked="" type="checkbox"/> LEVEL IV <input type="checkbox"/> NELAC <input checked="" type="checkbox"/> EDD/EDT (contact laboratory) <input type="checkbox"/> Other _____	

### Comments

- 1 Project performed under DoD QSM
- 2 TPH-d/o needs 3520 extraction
3. Preliminary data (or level 1) in 1-2 business days; Level IV report in 10 working days.
- 4 Note: NOI log is separate from other COC's
- 5 \*SVOC/VOC (full suite); PAH SIM (naphthalene, 1-methylnaphthalene, 2-methylnaphthalene)

### Project Information

Project Name, PWSID, Permit, etc.	CV18F0126, 60571032.02.46.01		
Sampler Name	ES, JV, MMC	Sampler Phone	808-593-6607
Sample Origin State	Hawaii	EPA/State Compliance	<input type="checkbox"/> Yes <input type="checkbox"/> No
The following tests will be subcontracted to other certified laboratories as shown. Signing this COC is authorization to subcontract the analyses as indicated.			
Analysis	Subcontract Lab		
TOC	Energy Laboratories Inc., Casper		

### Matrix Codes

- A - Air
- W - Water
- S - Soils/Solids
- V - Vegetation
- B - Bioassay
- O - Other
- DW - Drinking Water

### Analysis Requested

	8260 VOC's (Full Suite) + DCA* [40ml VOA w/HCL]	8015 TPH-g [40ml VOA w/HCL]	RSK175 Methane [40ml VOA w/H2SO4]	8011 EDB [40ml VOA w/HCL]	SVOCs (full suite+Nap, 1-2-Methylnap) by 8270DSIM*	EPA 3630/8015 TPH-d/o +SGC [1-L AG w/H2SO4]	EPA 9060 TOC [250ml AG w/H3PO4]	EPA 6020 Total Lead [250ml HDPE w/HNO3]	EPA 6020 Diss. Lead [250ml HDPE w/HNO3] (field Filtered)	See Attached
1	X	X	X	X	X	X	X	X	X	✓
2	X	X	X	X						✓
3										
4										
5										
6										
7										
8										
9										
10										

All turnaround times are standard unless marked as RUSH.  
Energy Laboratories MUST be contacted prior to RUSH sample submittal for charges and scheduling – See Instructions Page

Sample Identification (Name, Location, Interval, etc.)	Collection		Number of Containers	Matrix (See Codes Above)	Analysis Requested										See Attached	RUSH TAT	ELI LAB ID (Laboratory Use Only)
	Date	Time			8260 VOC's (Full Suite) + DCA* [40ml VOA w/HCL]	8015 TPH-g [40ml VOA w/HCL]	RSK175 Methane [40ml VOA w/H2SO4]	8011 EDB [40ml VOA w/HCL]	SVOCs (full suite+Nap, 1-2-Methylnap) by 8270DSIM*	EPA 3630/8015 TPH-d/o +SGC [1-L AG w/H2SO4]	EPA 9060 TOC [250ml AG w/H3PO4]	EPA 6020 Total Lead [250ml HDPE w/HNO3]	EPA 6020 Diss. Lead [250ml HDPE w/HNO3] (field Filtered)				
1 ERH2460 (RHMW12A)	1/17/22	1315	19	GW	X	X	X	X	X	X	X	X	X	X	✓	BZ201131	
2 ERH2459 (Trip Blank)	1/17/22	1230	8	WQ	X	X	X	X							✓		
3																	
4 TB - 8260 - 14694																	
5 TB - 6020 - 14733	1/17/22																
6 TB - 8011 - 14705																	
7 TB - methane - 14732																	
8																	
9																	
10																	

Custody Record MUST be signed	Relinquished by (print)	Date/Time	Signature	Received by (print)	Date/Time	Signature			
	Relinquished by (print)	Date/Time	Signature	Received by Laboratory (print)	Date/Time	Signature			
LABORATORY USE ONLY									
Shipped By	Cooler ID(s)	Custody Seals Y N C B	Intact Y N	Receipt Temp °C	Temp Blank Y N	On Ice Y N	Payment Type CC Cash Check	Amount \$	Receipt Number (cash/check only)



Work Order Receipt Checklist

AECOM - Honolulu

B22011131

Login completed by: Leslie S. Cadreau
Reviewed by: BL2000\gmccartney
Reviewed Date: 1/22/2022

Date Received: 1/19/2022
Received by: Isc
Carrier name: FedEx

- Shipping container/cooler in good condition? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on all shipping container(s)/cooler(s)? Yes [checked] No [ ] Not Present [ ]
Custody seals intact on all sample bottles? Yes [checked] No [ ] Not Present [ ]
Chain of custody present? Yes [checked] No [ ]
Chain of custody signed when relinquished and received? Yes [checked] No [ ]
Chain of custody agrees with sample labels? Yes [checked] No [ ]
Samples in proper container/bottle? Yes [checked] No [ ]
Sample containers intact? Yes [checked] No [ ]
Sufficient sample volume for indicated test? Yes [checked] No [ ]
All samples received within holding time? (Exclude analyses that are considered field parameters such as pH, DO, Res Cl, Sulfite, Ferrous Iron, etc.) Yes [checked] No [ ]
Temp Blank received in all shipping container(s)/cooler(s)? Yes [checked] No [ ] Not Applicable [ ]
Container/Temp Blank temperature: 0.6°C On Ice
Water - VOA vials have zero headspace? Yes [checked] No [ ] Not Applicable [ ]
Water - pH acceptable upon receipt? Yes [checked] No [ ] Not Applicable [ ]

Standard Reporting Procedures:

Lab measurement of analytes considered field parameters that require analysis within 15 minutes of sampling such as pH, Dissolved Oxygen and Residual Chlorine, are qualified as being analyzed outside of recommended holding time.

Solid/soil samples are reported on a wet weight basis (as received) unless specifically indicated. If moisture corrected, data units are typically noted as -dry. For agricultural and mining soil parameters/characteristics, all samples are dried and ground prior to sample analysis.

Contact and Corrective Action Comments:

The collection time indicated on the Chain of Custody for all samples is in Hawaii-Aleutian Standard Time. The collection time has been converted (+4 Hours) to Mountain Standard Time.

## Qualifiers and Abbreviations

Qualifier	Qualifier Description
##	Limit of Quantitation (LOQ) for this analyte exceeds the Maximum Contaminant Level (MCL)
*	Result exceeds the Maximum Contaminant Level (MCL)
A	The analyte level was greater than four times the spike level - in accordance with the method, percent recovery is not calculated
B	Analyte detected in the method blank
C	Continuing calibration verification was outside of the quality control advisory limits
D	Limit of Quantitation (LOQ) increased due to sample matrix
E	Estimated value - result exceeds the instrument upper quantitation limit
H	Analysis performed past the method holding time
J	The reported result is an estimated value
L	Lowest Limit of Quantitation (LOQ) available for the analytical method used
N	Analyte concentration was not sufficiently high to calculate a Relative Percent Difference (RPD) for the serial dilution test
O	Diluted out
P	Poor method performance - method validations have shown no recoveries at low concentrations or method performance was erratic
Q	Values reported below the Limit of Quantitation (LOQ) are statistically invalid
R	Relative Percent Difference (RPD) exceeds advisory limit
S	Spike recovery outside of advisory limits
T	Analyte detected in the associated trip blank
U	Not detected at the Limit of Detection (LOD)
V	The RPD value for this duplicate represents the RER value and the RPD limit of 2 is the RER upper limit.

## Qualifiers and Abbreviations

### Abbreviation

Reporting	Explanation of Abbreviation
DF	Dilution Factor
DL	Detection Limit
LOD	Limit of Detection
LOQ	Limit of Quantitation
MCL	Maximum Contaminant Level
MDC	Minimum Detectable Concentration
ND	Not detected at the Limit of Quantitation (LOQ)
RBSL	Risk-Based Screening Levels
REC	Recovery
RER	Relative Error Ratio
RPD	Relative Percent Difference
SPK	Spike

Sample Types	Explanation of Abbreviation
CCB	Continuing Calibration Blank
CCV	Continuing Calibration Verification Standard
DUP	Sample Duplicate
ICSA	Interference Check Sample A
ICSAB	Interference Check Sample AB
ICV	Initial Calibration Verification Standard
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
LFB	Laboratory Fortified Blank
LRB	Laboratory Reagent Blank
MBLK	Method Blank
MS	Sample Matrix Spike
MSD	Sample Matrix Spike Duplicate
PDS	Post Digestion/Distillation Spike
QCS	Quality Control Sample
SD	Serial Dilution
SRM	Standard Reference Material



**LABORATORY ANALYTICAL REPORT**

Prepared by Billings, MT Branch

**Lab ID:** B22011131-001  
**Collection Date:** 01/17/2022 17:15  
**Date Received:** 01/19/2022  
**Report Date:** 03/01/2022

**Client:** AECOM - Honolulu  
**Client Sample ID:** ERH2460 (RHMW12A)  
**Project:** CV18F0126/60571032.02.46.01  
**Matrix:** Ground Water

Analyses	Result	Units	DF	Qual	LOQ	LOD	DL	MCL	Method	Analysis Date / By	RunID : Run Order	BatchID
<b>LOW LEVEL PAH BY 8270C SIM</b>												
1-Methylnaphthalene	ND	ug/L	1	U	0.10	0.048	0.020		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
2-Methylnaphthalene	ND	ug/L	1	U	0.10	0.048	0.017		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
Acenaphthene	ND	ug/L	1	U	0.10	0.048	0.030		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
Acenaphthylene	ND	ug/L	1	U	0.10	0.048	0.024		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
Anthracene	ND	ug/L	1	U	0.10	0.048	0.027		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
Benzo(a)anthracene	ND	ug/L	1	U	0.10	0.048	0.026		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
Benzo(a)pyrene	ND	ug/L	1	U	0.10	0.048	0.033		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
Benzo(b)fluoranthene	ND	ug/L	1	U	0.10	0.048	0.022		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
Benzo(g,h,i)perylene	ND	ug/L	1	U	0.10	0.048	0.026		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
Benzo(k)fluoranthene	ND	ug/L	1	U	0.10	0.048	0.028		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
Chrysene	ND	ug/L	1	U	0.10	0.048	0.044		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
Dibenzo(a,h)anthracene	ND	ug/L	1	U	0.10	0.048	0.035		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
Fluoranthene	ND	ug/L	1	U	0.10	0.048	0.022		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
Fluorene	ND	ug/L	1	U	0.10	0.048	0.022		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
Indeno(1,2,3-cd)pyrene	ND	ug/L	1	U	0.10	0.048	0.047		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
Naphthalene	ND	ug/L	1	U	0.10	0.048	0.028		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
Phenanthrene	ND	ug/L	1	U	0.10	0.048	0.028		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
Pyrene	ND	ug/L	1	U	0.10	0.048	0.023		SW8270CSIM	01/27/2022 07:45/jph	SV5975.I_220126B : 5	163072
<b>AGGREGATE ORGANICS</b>												
Organic Carbon, Total (TOC) - TOC Range is 0.3 to 0.3	0.30	mg/L	1	J	0.50	0.50	0.17		SW9060A	01/21/2022 00:05/eli-ca	SUB-C278921 : 17	C_R278921
<b>METALS, DISSOLVED</b>												
Lead	ND	mg/L	1	U	0.001	0.0001	0.00006		SW6020	01/21/2022 23:19/car	ICPMS207-B_220121A : 70	R373694
<b>METALS, TOTAL</b>												
Lead	ND	mg/L	1	U	0.001	0.0001	0.00008		SW6020	01/21/2022 23:25/car	ICPMS207-B_220121A : 71	163063
<b>VOLATILE ORGANIC COMPOUNDS</b>												
Benzene	ND	ug/L	1	U	1.0	0.20	0.091		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Bromobenzene	ND	ug/L	1	U	1.0	0.20	0.083		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Bromochloromethane	ND	ug/L	1	U	1.0	0.50	0.14		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Bromodichloromethane	ND	ug/L	1	U	1.0	0.25	0.12		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Bromoform	ND	ug/L	1	U	1.0	0.25	0.12		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Carbon tetrachloride	ND	ug/L	1	U	1.0	0.50	0.14		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Chlorobenzene	ND	ug/L	1	U	1.0	0.20	0.091		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Chlorodibromomethane	ND	ug/L	1	U	1.0	0.20	0.084		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Chloroethane	ND	ug/L	1	U	1.0	0.50	0.17		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Chloroform	ND	ug/L	1	U	1.0	0.20	0.079		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Chloromethane	ND	ug/L	1	U	1.0	0.50	0.16		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695





### LABORATORY ANALYTICAL REPORT

Prepared by Billings, MT Branch

Lab ID: B22011131-001

Collection Date: 01/17/2022 17:15

Date Received: 01/19/2022

Report Date: 03/01/2022

**Client:** AECOM - Honolulu  
**Client Sample ID:** ERH2460 (RHMW12A)  
**Project:** CV18F0126/60571032.02.46.01  
**Matrix:** Ground Water

Analyses	Result	Units	DF	Qual	LOQ	LOD	DL	MCL	Method	Analysis Date / By	RunID : Run Order	BatchID
<b>VOLATILE ORGANIC COMPOUNDS</b>												
1,2-Dibromoethane	ND	ug/L	1	U	1.0	0.20	0.092		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
2-Chlorotoluene	ND	ug/L	1	U	1.0	0.20	0.088		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
4-Chlorotoluene	ND	ug/L	1	U	1.0	0.20	0.073		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Dibromomethane	ND	ug/L	1	U	1.0	0.50	0.15		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
1,2-Dichlorobenzene	ND	ug/L	1	U	1.0	0.20	0.075		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
1,3-Dichlorobenzene	ND	ug/L	1	U	1.0	0.20	0.080		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
1,4-Dichlorobenzene	ND	ug/L	1	U	1.0	0.20	0.086		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Dichlorodifluoromethane	ND	ug/L	1	U	1.0	0.50	0.18		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
1,1-Dichloroethane	ND	ug/L	1	U	1.0	0.50	0.14		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
1,2-Dichloroethane	ND	ug/L	1	U	1.0	0.25	0.12		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
1,1-Dichloroethene	ND	ug/L	1	U	1.0	0.50	0.14		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
cis-1,2-Dichloroethene	ND	ug/L	1	U	1.0	0.25	0.11		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
trans-1,2-Dichloroethene	ND	ug/L	1	U	1.0	0.25	0.12		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
1,2-Dichloropropane	ND	ug/L	1	U	1.0	0.20	0.085		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
1,3-Dichloropropane	ND	ug/L	1	U	1.0	0.20	0.079		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
2,2-Dichloropropane	ND	ug/L	1	U	1.0	0.50	0.19		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
1,1-Dichloropropene	ND	ug/L	1	U	1.0	0.20	0.083		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
cis-1,3-Dichloropropene	ND	ug/L	1	U	1.0	0.20	0.073		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
trans-1,3-Dichloropropene	ND	ug/L	1	U	1.0	0.20	0.085		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Ethylbenzene	ND	ug/L	1	U	1.0	0.20	0.084		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Methyl ethyl ketone	ND	ug/L	1	U	20	5.0	1.8		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Methyl tert-butyl ether (MTBE)	ND	ug/L	1	U	1.0	0.25	0.10		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Methylene chloride	ND	ug/L	1	U	1.0	0.50	0.34		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Styrene	ND	ug/L	1	U	1.0	0.20	0.067		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
1,1,1,2-Tetrachloroethane	ND	ug/L	1	U	1.0	0.25	0.10		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
1,1,2,2-Tetrachloroethane	ND	ug/L	1	U	1.0	0.20	0.087		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Tetrachloroethene	ND	ug/L	1	U	1.0	0.20	0.067		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Toluene	ND	ug/L	1	U	1.0	0.20	0.068		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
1,1,1-Trichloroethane	ND	ug/L	1	U	1.0	0.50	0.13		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
1,1,2-Trichloroethane	ND	ug/L	1	U	1.0	0.25	0.11		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Trichloroethene	ND	ug/L	1	U	1.0	0.20	0.099		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Trichlorofluoromethane	ND	ug/L	1	U	1.0	0.50	0.13		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
1,2,3-Trichloropropane	ND	ug/L	1	U	1.0	0.50	0.24		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Vinyl chloride	ND	ug/L	1	U	1.0	0.50	0.15		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
m+p-Xylenes	ND	ug/L	1	U	1.0	0.50	0.15		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
o-Xylene	ND	ug/L	1	U	1.0	0.20	0.060		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Xylenes, Total	ND	ug/L	1	U	1.0	0.20	0.060		SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Surr: Dibromofluoromethane	102.0	%REC	1			80-119			SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Surr: 1,2-Dichloroethane-d4	105.0	%REC	1			81-118			SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695



### LABORATORY ANALYTICAL REPORT

Prepared by Billings, MT Branch

**Lab ID:** B22011131-001  
**Collection Date:** 01/17/2022 17:15  
**Date Received:** 01/19/2022  
**Report Date:** 03/01/2022

**Client:** AECOM - Honolulu  
**Client Sample ID:** ERH2460 (RHMW12A)  
**Project:** CV18F0126/60571032.02.46.01  
**Matrix:** Ground Water

Analyses	Result	Units	DF	Qual	LOQ	LOD	DL	MCL	Method	Analysis Date / By	RunID : Run Order	BatchID
<b>VOLATILE ORGANIC COMPOUNDS</b>												
Surr: Toluene-d8	96.0	%REC	1		89-112				SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
Surr: p-Bromofluorobenzene	101.0	%REC	1		85-114				SW8260B	01/21/2022 15:42/msc	VOA5975C.I_220121A : 12	R373695
<b>VOCS BY MICROEXTRACTION-ECD</b>												
1,2-Dibromoethane	ND	ug/L	1	U	0.010	0.0048	0.0025		SW8011	01/22/2022 10:35/clt	GECD.I_220121A : 49	163129
Surr: 1,1,1,2-Tetrachloroethane	89.0	%REC	1		70-130				SW8011	01/22/2022 10:35/clt	GECD.I_220121A : 49	163129
<b>PETROLEUM HYDROCARBONS-VOLATILE</b>												
C6 to C10	ND	ug/L	1	U	20	8.7	2.3		SW8015C	01/21/2022 10:16/jp	PE 1_220120A : 33	R373498
Total Purgeable Hydrocarbons	ND	ug/L	1	U	20	10	3.6		SW8015C	01/21/2022 10:16/jp	PE 1_220120A : 33	R373498
Surr: Trifluorotoluene	80.0	%REC	1		70-130				SW8015C	01/21/2022 10:16/jp	PE 1_220120A : 33	R373498
- Note 1: C6 to C10 is defined as all hydrocarbons eluting between 2-Methylpentane and 1,2,4-Trimethylbenzene.												
- Note 2: Total Purgeable Hydrocarbons are defined as the total hydrocarbon response regardless of elution time.												
<b>PETROLEUM HYDROCARBONS-SEMI-VOLATILE</b>												
Diesel Range Organics (C10 to C24)	ND	mg/L	1	U	0.30	0.14	0.037		SW8015C	01/23/2022 01:56/amn	GCFID-HP5-B_220122A : 14	163074
Oil Range Hydrocarbons (C24 to C40)	ND	mg/L	1	U	0.30	0.14	0.084		SW8015C	01/23/2022 01:56/amn	GCFID-HP5-B_220122A : 14	163074
Total Extractable Hydrocarbons	ND	mg/L	1	U	0.30	0.14	0.071		SW8015C	01/23/2022 01:56/amn	GCFID-HP5-B_220122A : 14	163074
Surr: o-Terphenyl	96.0	%REC	1		56-125				SW8015C	01/23/2022 01:56/amn	GCFID-HP5-B_220122A : 14	163074
Surr: n-Triacontane	101.0	%REC	1		50-150				SW8015C	01/23/2022 01:56/amn	GCFID-HP5-B_220122A : 14	163074
- Note: Total Extractable Hydrocarbons are defined as the total hydrocarbon responses regardless of elution time.												
- Since there were no detectable hydrocarbons, Silica Gel Treatment (SGT) results are equivalent to non-SGT results.												
<b>ORGANIC CHARACTERISTICS</b>												
Methane	ND	mg/L	1	U	0.0020	0.0012	0.00070		SW8015M	01/20/2022 12:20/jdw	FID-HEADSPACE_220120A : 20	R373491
<b>SEMI-VOLATILE ORGANIC COMPOUNDS</b>												
1,2,4-Trichlorobenzene	ND	ug/L	1	U	10	4.8	1.8		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
1,2-Dichlorobenzene	ND	ug/L	1	U	10	4.8	1.9		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
1,3-Dichlorobenzene	ND	ug/L	1	U	10	4.8	2.0		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
1,4-Dichlorobenzene	ND	ug/L	1	U	10	4.8	1.9		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
2,4,5-Trichlorophenol	ND	ug/L	1	U	10	4.8	2.1		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
2,4,6-Trichlorophenol	ND	ug/L	1	U	10	4.8	2.5		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
2,4-Dichlorophenol	ND	ug/L	1	U	10	4.8	1.6		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
2,4-Dimethylphenol	ND	ug/L	1	U	10	4.8	1.6		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
2,4-Dinitrophenol	ND	ug/L	1	U	10	9.6	4.1		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
2,4-Dinitrotoluene	ND	ug/L	1	U	10	4.8	2.9		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
2,6-Dinitrotoluene	ND	ug/L	1	U	10	4.8	3.1		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
2-Chloronaphthalene	ND	ug/L	1	U	10	4.8	2.1		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
2-Chlorophenol	ND	ug/L	1	U	10	4.8	2.4		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
2-Nitrophenol	ND	ug/L	1	U	10	4.8	2.3		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
3,3'-Dichlorobenzidine	ND	ug/L	1	U	10	4.8	2.0		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
4,6-Dinitro-2-methylphenol	ND	ug/L	1	U	10	9.6	2.2		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072



### LABORATORY ANALYTICAL REPORT

Prepared by Billings, MT Branch

Lab ID: B22011131-001

Collection Date: 01/17/2022 17:15

Date Received: 01/19/2022

Report Date: 03/01/2022

**Client:** AECOM - Honolulu  
**Client Sample ID:** ERH2460 (RHMW12A)  
**Project:** CV18F0126/60571032.02.46.01  
**Matrix:** Ground Water

Analyses	Result	Units	DF	Qual	LOQ	LOD	DL	MCL	Method	Analysis Date / By	RunID : Run Order	BatchID
<b>SEMI-VOLATILE ORGANIC COMPOUNDS</b>												
4-Bromophenyl phenyl ether	ND	ug/L	1	U	10	4.8	1.7		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
4-Chloro-3-methylphenol	ND	ug/L	1	U	10	4.8	1.4		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
4-Chlorophenol	ND	ug/L	1	U	10	4.8	2.5		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
4-Chlorophenyl phenyl ether	ND	ug/L	1	U	10	4.8	2.0		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
4-Nitrophenol	ND	ug/L	1	U	10	9.6	2.4		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Azobenzene	ND	ug/L	1	U	10	4.8	1.0		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
bis(-2-chloroethoxy)Methane	ND	ug/L	1	U	10	4.8	1.3		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
bis(-2-chloroethyl)Ether	ND	ug/L	1	U	10	4.8	2.5		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
bis(2-chloroisopropyl)Ether	ND	ug/L	1	U	10	4.8	1.4		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
bis(2-ethylhexyl)Phthalate	ND	ug/L	1	U	10	4.8	1.8		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Butylbenzylphthalate	ND	ug/L	1	U	10	4.8	1.5		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Di-n-butyl phthalate	ND	ug/L	1	U	10	4.8	0.90		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Di-n-octyl phthalate	ND	ug/L	1	U	10	4.8	1.3		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Diethyl phthalate	ND	ug/L	1	U	10	4.8	2.1		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Dimethyl phthalate	ND	ug/L	1	U	10	4.8	1.7		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Hexachlorobenzene	ND	ug/L	1	U	10	4.8	1.3		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Hexachlorobutadiene	ND	ug/L	1	U	10	4.8	2.2		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Hexachlorocyclopentadiene	ND	ug/L	1	U	10	4.8	2.9		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Hexachloroethane	ND	ug/L	1	U	10	4.8	1.7		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Isophorone	ND	ug/L	1	U	10	4.8	1.6		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
m+p-Cresols	ND	ug/L	1	U	10	4.8	1.7		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
n-Nitroso-di-n-propylamine	ND	ug/L	1	U	10	4.8	1.5		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
n-Nitrosodimethylamine	ND	ug/L	1	U	10	4.8	1.5		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
n-Nitrosodiphenylamine	ND	ug/L	1	U	10	4.8	1.1		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Nitrobenzene	ND	ug/L	1	U	10	4.8	2.2		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
o-Cresol	ND	ug/L	1	U	10	4.8	1.8		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Pentachlorophenol	ND	ug/L	1	U	10	9.6	4.1		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Phenol	ND	ug/L	1	U	10	4.8	1.4		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Pyridine	ND	ug/L	1	U	10	4.8	3.1		SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Surr: 2,4,6-Tribromophenol	80.0	%REC	1		43-140				SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Surr: 2-Fluorobiphenyl	62.0	%REC	1		44-119				SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Surr: 2-Fluorophenol	31.0	%REC	1		19-119				SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Surr: Nitrobenzene-d5	60.0	%REC	1		44-120				SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Surr: Phenol-d5	30.0	%REC	1		10-65				SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072
Surr: Terphenyl-d14	97.0	%REC	1		50-134				SW8270C	02/2/2022 04:06/dsm	SV5973N.I_220201A : 22	163072



### LABORATORY ANALYTICAL REPORT

Prepared by Billings, MT Branch

Lab ID: B22011131-002

Collection Date: 01/17/2022 17:15

Date Received: 01/19/2022

Report Date: 03/01/2022

**Client:** AECOM - Honolulu  
**Client Sample ID:** ERH2459 (Trip Blank)-14694  
**Project:** CV18F0126/60571032.02.46.01  
**Matrix:** Trip Blank

Analyses	Result	Units	DF	Qual	LOQ	LOD	DL	MCL	Method	Analysis Date / By	RunID : Run Order	BatchID
<b>VOLATILE ORGANIC COMPOUNDS</b>												
Benzene	ND	ug/L	1	U	1.0	0.20	0.091		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Bromobenzene	ND	ug/L	1	U	1.0	0.20	0.083		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Bromochloromethane	ND	ug/L	1	U	1.0	0.50	0.14		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Bromodichloromethane	ND	ug/L	1	U	1.0	0.25	0.12		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Bromoform	ND	ug/L	1	U	1.0	0.25	0.12		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Carbon tetrachloride	ND	ug/L	1	U	1.0	0.50	0.14		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Chlorobenzene	ND	ug/L	1	U	1.0	0.20	0.091		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Chlorodibromomethane	ND	ug/L	1	U	1.0	0.20	0.084		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Chloroethane	ND	ug/L	1	U	1.0	0.50	0.17		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Chloroform	ND	ug/L	1	U	1.0	0.20	0.079		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Chloromethane	ND	ug/L	1	U	1.0	0.50	0.16		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
1,2-Dibromoethane	ND	ug/L	1	U	1.0	0.20	0.092		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
2-Chlorotoluene	ND	ug/L	1	U	1.0	0.20	0.088		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
4-Chlorotoluene	ND	ug/L	1	U	1.0	0.20	0.073		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Dibromomethane	ND	ug/L	1	U	1.0	0.50	0.15		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
1,2-Dichlorobenzene	ND	ug/L	1	U	1.0	0.20	0.075		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
1,3-Dichlorobenzene	ND	ug/L	1	U	1.0	0.20	0.080		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
1,4-Dichlorobenzene	ND	ug/L	1	U	1.0	0.20	0.086		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Dichlorodifluoromethane	ND	ug/L	1	U	1.0	0.50	0.18		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
1,1-Dichloroethane	ND	ug/L	1	U	1.0	0.50	0.14		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
1,2-Dichloroethane	ND	ug/L	1	U	1.0	0.25	0.12		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
1,1-Dichloroethene	ND	ug/L	1	U	1.0	0.50	0.14		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
cis-1,2-Dichloroethene	ND	ug/L	1	U	1.0	0.25	0.11		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
trans-1,2-Dichloroethene	ND	ug/L	1	U	1.0	0.25	0.12		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
1,2-Dichloropropane	ND	ug/L	1	U	1.0	0.20	0.085		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
1,3-Dichloropropane	ND	ug/L	1	U	1.0	0.20	0.079		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
2,2-Dichloropropane	ND	ug/L	1	U	1.0	0.50	0.19		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
1,1-Dichloropropene	ND	ug/L	1	U	1.0	0.20	0.083		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
cis-1,3-Dichloropropene	ND	ug/L	1	U	1.0	0.20	0.073		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
trans-1,3-Dichloropropene	ND	ug/L	1	U	1.0	0.20	0.085		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Ethylbenzene	ND	ug/L	1	U	1.0	0.20	0.084		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Methyl ethyl ketone	ND	ug/L	1	U	20	5.0	1.8		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Methyl tert-butyl ether (MTBE)	ND	ug/L	1	U	1.0	0.25	0.10		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Methylene chloride	ND	ug/L	1	U	1.0	0.50	0.34		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Styrene	ND	ug/L	1	U	1.0	0.20	0.067		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
1,1,1,2-Tetrachloroethane	ND	ug/L	1	U	1.0	0.25	0.10		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
1,1,2,2-Tetrachloroethane	ND	ug/L	1	U	1.0	0.20	0.087		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Tetrachloroethene	ND	ug/L	1	U	1.0	0.20	0.067		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Toluene	ND	ug/L	1	U	1.0	0.20	0.068		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695



**LABORATORY ANALYTICAL REPORT**

Prepared by Billings, MT Branch

Lab ID: B22011131-002

Collection Date: 01/17/2022 17:15

Date Received: 01/19/2022

Report Date: 03/01/2022

**Client:** AECOM - Honolulu  
**Client Sample ID:** ERH2459 (Trip Blank)-14694  
**Project:** CV18F0126/60571032.02.46.01  
**Matrix:** Trip Blank

Analyses	Result	Units	DF	Qual	LOQ	LOD	DL	MCL	Method	Analysis Date / By	RunID : Run Order	BatchID
<b>VOLATILE ORGANIC COMPOUNDS</b>												
1,1,1-Trichloroethane	ND	ug/L	1	U	1.0	0.50	0.13		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
1,1,2-Trichloroethane	ND	ug/L	1	U	1.0	0.25	0.11		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Trichloroethene	ND	ug/L	1	U	1.0	0.20	0.099		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Trichlorofluoromethane	ND	ug/L	1	U	1.0	0.50	0.13		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
1,2,3-Trichloropropane	ND	ug/L	1	U	1.0	0.50	0.24		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Vinyl chloride	ND	ug/L	1	U	1.0	0.50	0.15		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
m+p-Xylenes	ND	ug/L	1	U	1.0	0.50	0.15		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
o-Xylene	ND	ug/L	1	U	1.0	0.20	0.060		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Xylenes, Total	ND	ug/L	1	U	1.0	0.20	0.060		SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Surr: Dibromofluoromethane	109.0	%REC	1		80-119				SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Surr: 1,2-Dichloroethane-d4	109.0	%REC	1		81-118				SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Surr: Toluene-d8	100.0	%REC	1		89-112				SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695
Surr: p-Bromofluorobenzene	102.0	%REC	1		85-114				SW8260B	01/21/2022 19:20/msc	VOA5975C.I_220121A : 19	R373695





**LABORATORY ANALYTICAL REPORT**

Prepared by Billings, MT Branch

**Lab ID:** B22011131-003

**Collection Date:** 01/17/2022 17:15

**Date Received:** 01/19/2022

**Report Date:** 03/01/2022

**Client:** AECOM - Honolulu  
**Client Sample ID:** ERH2459 (Trip Blank)-14733  
**Project:** CV18F0126/60571032.02.46.01  
**Matrix:** Trip Blank

Analyses	Result	Units	DF	Qual	LOQ	LOD	DL	MCL	Method	Analysis Date / By	RunID : Run Order	BatchID
<b>PETROLEUM HYDROCARBONS-VOLATILE</b>												
C6 to C10	ND	ug/L	1	U	20	8.7	2.3		SW8015C	01/20/2022 17:07/jp	PE 1_220120A : 12	R373498
Total Purgeable Hydrocarbons	ND	ug/L	1	U	20	10	3.6		SW8015C	01/20/2022 17:07/jp	PE 1_220120A : 12	R373498
Surr: Trifluorotoluene	79.0	%REC	1		70-130				SW8015C	01/20/2022 17:07/jp	PE 1_220120A : 12	R373498

- Note 1: C6 to C10 is defined as all hydrocarbons eluting between 2-Methylpentane and 1,2,4-Trimethylbenzene.

- Note 2: Total Purgeable Hydrocarbons are defined as the total hydrocarbon response regardless of elution time.



### LABORATORY ANALYTICAL REPORT

Prepared by Billings, MT Branch

**Lab ID:** B22011131-004  
**Collection Date:** 01/17/2022 17:15  
**Date Received:** 01/19/2022  
**Report Date:** 03/01/2022

**Client:** AECOM - Honolulu  
**Client Sample ID:** ERH2459 (Trip Blank)-14705  
**Project:** CV18F0126/60571032.02.46.01  
**Matrix:** Trip Blank

Analyses	Result	Units	DF	Qual	LOQ	LOD	DL	MCL	Method	Analysis Date / By	RunID : Run Order	BatchID
<b>VOCS BY MICROEXTRACTION-ECD</b>												
1,2-Dibromoethane	ND	ug/L	1	U	0.010	0.0049	0.0025		SW8011	01/22/2022 08:15/ct	GECD.I_220121A : 42	163129
Surr: 1,1,1,2-Tetrachloroethane	93.0	%REC	1		70-130				SW8011	01/22/2022 08:15/ct	GECD.I_220121A : 42	163129



### LABORATORY ANALYTICAL REPORT

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Client Sample ID:** ERH2459 (Trip Blank)-14732  
**Project:** CV18F0126/60571032.02.46.01  
**Matrix:** Trip Blank

**Lab ID:** B22011131-005  
**Collection Date:** 01/17/2022 17:15  
**Date Received:** 01/19/2022  
**Report Date:** 03/01/2022

Analyses	Result	Units	DF	Qual	LOQ	LOD	DL	MCL	Method	Analysis Date / By	RunID : Run Order	BatchID
<b>ORGANIC CHARACTERISTICS</b>												
Methane	ND	mg/L	1	U	0.0020	0.0012	0.00070		SW8015M	01/20/2022 12:26/jdw	FID-HEADSPACE_220120A : 21	R373491





### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5975.I\_220126A: 16  
**Method:** SW8270CSIM  
**Lab ID:** MB-163072

**SampType:** Method Blank  
**Analysis Date:** 01/27/2022 00:20  
**Units:** ug/L

**Batch ID:** 163072  
**Prep Date:** 01/19/2022 15:57  
**Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1-Methylnaphthalene	ND	0.10									
2-Methylnaphthalene	ND	0.10									
Acenaphthene	ND	0.10									
Acenaphthylene	ND	0.10									
Anthracene	ND	0.10									
Benzo(a)anthracene	ND	0.10									
Benzo(a)pyrene	ND	0.10									
Benzo(b)fluoranthene	ND	0.10									
Benzo(g,h,i)perylene	ND	0.10									
Benzo(k)fluoranthene	ND	0.10									
Chrysene	ND	0.10									
Dibenzo(a,h)anthracene	ND	0.10									
Fluoranthene	ND	0.10									
Fluorene	ND	0.10									
Indeno(1,2,3-cd)pyrene	ND	0.10									
Naphthalene	ND	0.10									
Phenanthrene	ND	0.10									
Pyrene	ND	0.10									

Associated Samples: **B22011131-001C**

**Run ID: Run Order:** SV5975.I\_220126A: 17  
**Method:** SW8270CSIM  
**Lab ID:** LLCS-163072

**SampType:** Laboratory Control Sample  
**Analysis Date:** 01/27/2022 00:52  
**Units:** ug/L

**Batch ID:** 163072  
**Prep Date:** 01/19/2022 15:58  
**Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1-Methylnaphthalene	2.6	0.10	5.0		51.0	41	115				



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5975.I\_220126A: 17      **SampType:** Laboratory Control Sample      **Batch ID:** 163072  
**Method:** SW8270CSIM      **Analysis Date:** 01/27/2022 00:52      **Prep Date:** 01/19/2022 15:58  
**Lab ID:** LLCS-163072      **Units:** ug/L      **Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
2-Methylnaphthalene	2.6	0.10	5.0		51.0	39	114				
Acenaphthene	3.4	0.10	5.0		67.0	48	114				
Acenaphthylene	3.5	0.10	5.0		70.0	35	121				
Anthracene	5.1	0.10	5.0		102.0	53	119				
Benzo(a)anthracene	5.8	0.10	5.0		116.0	59	120				
Benzo(a)pyrene	5.3	0.10	5.0		106.0	53	120				
Benzo(b)fluoranthene	5.3	0.10	5.0		106.0	53	126				
Benzo(g,h,i)perylene	5.3	0.10	5.0		106.0	44	128				
Benzo(k)fluoranthene	4.9	0.10	5.0		97.0	54	125				
Chrysene	5.5	0.10	5.0		109.0	57	120				
Dibenzo(a,h)anthracene	5.7	0.10	5.0		114.0	44	141				
Fluoranthene	5.2	0.10	5.0		104.0	58	120				
Fluorene	3.5	0.10	5.0		69.0	50	118				
Indeno(1,2,3-cd)pyrene	5.4	0.10	5.0		109.0	48	130				
Naphthalene	2.5	0.10	5.0		51.0	43	114				
Phenanthrene	4.5	0.10	5.0		91.0	53	115				
Pyrene	5.2	0.10	5.0		103.0	53	121				

Associated Samples: **B22011131-001C**

**Run ID: Run Order:** SV5975.I\_220126A: 18      **SampType:** Laboratory Control Sample Duplicate      **Batch ID:** 163072  
**Method:** SW8270CSIM      **Analysis Date:** 01/27/2022 01:25      **Prep Date:** 01/19/2022 15:58  
**Lab ID:** LLCSD-163072      **Units:** ug/L      **Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1-Methylnaphthalene	2.7	0.10	5.0		55.0	41	115	2.6	6.4	40.0	
2-Methylnaphthalene	2.8	0.10	5.0		57.0	39	114	2.6	9.5	40.0	



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5975.I\_220126A: 18      **SampType:** Laboratory Control Sample Duplicate      **Batch ID:** 163072  
**Method:** SW8270CSIM      **Analysis Date:** 01/27/2022 01:25      **Prep Date:** 01/19/2022 15:58  
**Lab ID:** LLCSD-163072      **Units:** ug/L      **Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Acenaphthene	3.6	0.10	5.0		71.0	48	114	3.4	5.1	40.0	
Acenaphthylene	3.7	0.10	5.0		74.0	35	121	3.5	6.6	40.0	
Anthracene	5.3	0.10	5.0		107.0	53	119	5.1	4.4	40.0	
Benzo(a)anthracene	5.8	0.10	5.0		115.0	59	120	5.8	0.4	40.0	
Benzo(a)pyrene	5.0	0.10	5.0		100.0	53	120	5.3	5.7	40.0	
Benzo(b)fluoranthene	5.1	0.10	5.0		102.0	53	126	5.3	3.7	40.0	
Benzo(g,h,i)perylene	5.3	0.10	5.0		105.0	44	128	5.3	0.7	40.0	
Benzo(k)fluoranthene	4.7	0.10	5.0		93.0	54	125	4.9	3.9	40.0	
Chrysene	5.4	0.10	5.0		109.0	57	120	5.5	0.5	40.0	
Dibenzo(a,h)anthracene	5.5	0.10	5.0		110.0	44	141	5.7	3.5	40.0	
Fluoranthene	5.4	0.10	5.0		107.0	58	120	5.2	3.1	40.0	
Fluorene	3.8	0.10	5.0		76.0	50	118	3.5	9.6	40.0	
Indeno(1,2,3-cd)pyrene	5.2	0.10	5.0		104.0	48	130	5.4	4.5	40.0	
Naphthalene	2.5	0.10	5.0		51.0	43	114	2.5	0.4	40.0	
Phenanthrene	4.8	0.10	5.0		96.0	53	115	4.5	5.6	40.0	
Pyrene	5.2	0.10	5.0		104.0	53	121	5.2	1.2	40.0	

Associated Samples: **B22011131-001C**

**Run ID: Run Order:** SV5975.I\_220126B: 12      **SampType:** Sample Matrix Spike      **Batch ID:** 163072  
**Method:** SW8270CSIM      **Analysis Date:** 01/27/2022 11:31      **Prep Date:** 01/20/2022 13:07  
**Lab ID:** B22011136-001CLMS      **Units:** ug/L      **Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1-Methylnaphthalene	3.4	0.10	5.1	0.0	66.0	18	117				
2-Methylnaphthalene	3.1	0.10	5.1	0.0	61.0	17	118				
Acenaphthene	4.2	0.10	5.1	0.0	82.0	40	92				



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5975.I\_220126B: 12      **SampType:** Sample Matrix Spike      **Batch ID:** 163072  
**Method:** SW8270CSIM      **Analysis Date:** 01/27/2022 11:31      **Prep Date:** 01/20/2022 13:07  
**Lab ID:** B22011136-001CLMS      **Units:** ug/L      **Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Acenaphthylene	4.0	0.10	5.1	0.0	78.0	37	96				
Anthracene	5.2	0.10	5.1	0.0	102.0	46	108				
Benzo(a)anthracene	5.7	0.10	5.1	0.0	112.0	41	105				S
Benzo(a)pyrene	4.9	0.10	5.1	0.0	96.0	42	110				
Benzo(b)fluoranthene	5.0	0.10	5.1	0.0	99.0	27	121				
Benzo(g,h,i)perylene	5.0	0.10	5.1	0.0	98.0	44	108				
Benzo(k)fluoranthene	4.7	0.10	5.1	0.0	91.0	44	111				
Chrysene	5.3	0.10	5.1	0.0	104.0	50	106				
Dibenzo(a,h)anthracene	5.4	0.10	5.1	0.0	106.0	47	111				
Fluoranthene	5.0	0.10	5.1	0.0	98.0	44	111				
Fluorene	4.4	0.10	5.1	0.0	86.0	42	99				
Indeno(1,2,3-cd)pyrene	5.1	0.10	5.1	0.0	101.0	33	112				
Naphthalene	3.7	0.10	5.1	0.0	73.0	22	108				
Phenanthrene	5.1	0.10	5.1	0.0	100.0	43	106				
Pyrene	5.3	0.10	5.1	0.0	103.0	41	106				

Associated Samples: **B22011131-001C**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5975.I\_220126B: 13      **SampType:** Sample Matrix Spike Duplicate      **Batch ID:** 163072  
**Method:** SW8270CSIM      **Analysis Date:** 01/27/2022 12:04      **Prep Date:** 01/20/2022 13:07  
**Lab ID:** B22011136-001CLMSD      **Units:** ug/L      **Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1-Methylnaphthalene	3.9	0.10	5.0	0.0	78.0	18	117	3.4	15.0	40.0	
2-Methylnaphthalene	4.4	0.10	5.0	0.0	86.0	18	117	3.1	34.0	40.0	
Acenaphthene	4.1	0.10	5.0	0.0	81.0	40	92	4.2	2.3	40.0	
Acenaphthylene	4.1	0.10	5.0	0.0	81.0	37	96	4.0	2.7	40.0	
Anthracene	4.7	0.10	5.0	0.0	93.0	46	108	5.2	10.0	40.0	
Benzo(a)anthracene	4.7	0.10	5.0	0.0	92.0	41	105	5.7	20.0	40.0	
Benzo(a)pyrene	4.0	0.10	5.0	0.0	79.0	42	110	4.9	21.0	40.0	
Benzo(b)fluoranthene	4.0	0.10	5.0	0.0	80.0	27	121	5.0	22.0	40.0	
Benzo(g,h,i)perylene	4.0	0.10	5.0	0.0	80.0	44	108	5.0	21.0	40.0	
Benzo(k)fluoranthene	3.7	0.10	5.0	0.0	74.0	44	111	4.7	22.0	40.0	
Chrysene	4.4	0.10	5.0	0.0	87.0	50	106	5.3	18.0	40.0	
Dibenzo(a,h)anthracene	4.4	0.10	5.0	0.0	87.0	47	111	5.4	21.0	40.0	
Fluoranthene	4.3	0.10	5.0	0.0	86.0	44	111	5.0	15.0	40.0	
Fluorene	4.2	0.10	5.0	0.0	84.0	42	99	4.4	2.7	40.0	
Indeno(1,2,3-cd)pyrene	4.2	0.10	5.0	0.0	82.0	33	112	5.1	21.0	40.0	
Naphthalene	4.1	0.10	5.0	0.0	82.0	22	108	3.7	9.9	40.0	
Phenanthrene	4.5	0.10	5.0	0.0	90.0	43	106	5.1	12.0	40.0	
Pyrene	4.6	0.10	5.0	0.0	91.0	41	106	5.3	14.0	40.0	

Associated Samples: **B22011131-001C**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5975.I\_220126B: 2      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R373846  
**Method:** SW8270CSIM      **Analysis Date:** 01/27/2022 06:08      **Prep Date:**  
**Lab ID:** 26-Jan-22\_CCV\_26      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1-Methylnaphthalene	2.1	0.10	2.0		103.0	80	120				
2-Methylnaphthalene	2.2	0.10	2.0		110.0	80	120				
Acenaphthene	2.0	0.10	2.0		101.0	80	120				
Acenaphthylene	2.0	0.10	2.0		101.0	80	120				
Anthracene	2.1	0.10	2.0		103.0	80	120				
Benzo(a)anthracene	2.1	0.10	2.0		105.0	80	120				
Benzo(a)pyrene	2.1	0.10	2.0		106.0	80	120				
Benzo(b)fluoranthene	2.3	0.10	2.0		115.0	80	120				
Benzo(g,h,i)perylene	2.2	0.10	2.0		110.0	80	120				
Benzo(k)fluoranthene	2.0	0.10	2.0		102.0	80	120				
Chrysene	2.0	0.10	2.0		100.0	80	120				
Dibenzo(a,h)anthracene	2.1	0.10	2.0		107.0	80	120				
Fluoranthene	2.1	0.10	2.0		106.0	80	120				
Fluorene	2.0	0.10	2.0		100.0	80	120				
Indeno(1,2,3-cd)pyrene	2.3	0.10	2.0		113.0	80	120				
Naphthalene	2.1	0.10	2.0		105.0	80	120				
Phenanthrene	1.9	0.10	2.0		93.0	80	120				
Pyrene	2.1	0.10	2.0		103.0	80	120				

Associated Samples: **B22011131-001C**

**Run ID: Run Order:** SV5975.I\_220126B: 15      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R373846  
**Method:** SW8270CSIM      **Analysis Date:** 01/28/2022 08:43      **Prep Date:**  
**Lab ID:** 26-Jan-22\_CCV\_39      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1-Methylnaphthalene	2.3	0.10	2.0		113.0	50	150				



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5975.I\_220126B: 15      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R373846  
**Method:** SW8270CSIM      **Analysis Date:** 01/28/2022 08:43      **Prep Date:**  
**Lab ID:** 26-Jan-22\_CCV\_39      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
2-Methylnaphthalene	2.5	0.10	2.0		124.0	50	150				
Acenaphthene	1.9	0.10	2.0		95.0	50	150				
Acenaphthylene	2.0	0.10	2.0		99.0	50	150				
Anthracene	2.2	0.10	2.0		111.0	50	150				
Benzo(a)anthracene	2.3	0.10	2.0		116.0	50	150				
Benzo(a)pyrene	2.3	0.10	2.0		116.0	50	150				
Benzo(b)fluoranthene	2.2	0.10	2.0		109.0	50	150				
Benzo(g,h,i)perylene	2.3	0.10	2.0		113.0	50	150				
Benzo(k)fluoranthene	2.0	0.10	2.0		102.0	50	150				
Chrysene	2.1	0.10	2.0		106.0	50	150				
Dibenzo(a,h)anthracene	2.3	0.10	2.0		114.0	50	150				
Fluoranthene	2.1	0.10	2.0		107.0	50	150				
Fluorene	2.1	0.10	2.0		106.0	50	150				
Indeno(1,2,3-cd)pyrene	2.4	0.10	2.0		121.0	50	150				
Naphthalene	2.5	0.10	2.0		123.0	50	150				
Phenanthrene	2.3	0.10	2.0		113.0	50	150				
Pyrene	2.2	0.10	2.0		109.0	50	150				

Associated Samples: **B22011131-001C**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SUB-C278921: 2      **SampType:** Method Blank      **Batch ID:** C\_R278921  
**Method:** SW9060A      **Analysis Date:** 01/20/2022 16:27      **Prep Date:**  
**Lab ID:** MBLK      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Organic Carbon, Total (TOC)	ND	0.20									

Associated Samples: **B22011131-001E**  
- TOC Range is 0.0 to 0.1

**Run ID: Run Order:** SUB-C278921: 1      **SampType:** Laboratory Control Sample      **Batch ID:** C\_R278921  
**Method:** SW9060A      **Analysis Date:** 01/20/2022 15:46      **Prep Date:**  
**Lab ID:** LCS      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Organic Carbon, Total (TOC)	5.2	0.50	5.0		104.0	91	111				

Associated Samples: **B22011131-001E**  
- TOC Range is 5.2 to 5.3

**Run ID: Run Order:** SUB-C278921: 8      **SampType:** Sample Matrix Spike      **Batch ID:** C\_R278921  
**Method:** SW9060A      **Analysis Date:** 01/21/2022 04:10      **Prep Date:**  
**Lab ID:** C22010624-001EMS      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Organic Carbon, Total (TOC)	5.9	0.50	5.0	0.68	105.0	91	111				

Associated Samples: **B22011131-001E**  
- TOC Range is 5.9 to 6.0





### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SUB-C278921: 9      **SampType:** Sample Matrix Spike Duplicate      **Batch ID:** C\_R278921  
**Method:** SW9060A      **Analysis Date:** 01/21/2022 04:51      **Prep Date:**  
**Lab ID:** C22010624-001EMSD      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Organic Carbon, Total (TOC)	6.1	0.50	5.0	0.68	109.0	91	111	5.9	3.0	10.0	

Associated Samples: **B22011131-001E**  
- TOC Range is 6.0 to 6.2

**Run ID: Run Order:** SUB-C278921: 3      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** C\_R278921  
**Method:** SW9060A      **Analysis Date:** 01/20/2022 17:05      **Prep Date:**  
**Lab ID:** CCV      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Organic Carbon, Total (TOC)	5.0	0.50	5.0		100.0	90	110				

Associated Samples: **B22011131-001E**  
- TOC Range is 5.0 to 5.1

**Run ID: Run Order:** SUB-C278921: 7      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** C\_R278921  
**Method:** SW9060A      **Analysis Date:** 01/21/2022 02:10      **Prep Date:**  
**Lab ID:** CCV      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Organic Carbon, Total (TOC)	5.1	0.50	5.0		103.0	90	110				

Associated Samples: **B22011131-001E**  
- TOC Range is 5.1 to 5.2



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** ICPMS207-B\_220121A: 29      **SampType:** Method Blank      **Batch ID:** R373694  
**Method:** SW6020      **Analysis Date:** 01/21/2022 19:03      **Prep Date:**  
**Lab ID:** LRB      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	ND	0.0005									

Associated Samples: **B22011131-001A**

**Run ID: Run Order:** ICPMS207-B\_220121A: 30      **SampType:** Laboratory Fortified Blank      **Batch ID:** R373694  
**Method:** SW6020      **Analysis Date:** 01/21/2022 19:09      **Prep Date:**  
**Lab ID:** LFB      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	0.050	0.001	0.050		100.0	88	115				

Associated Samples: **B22011131-001A**

**Run ID: Run Order:** ICPMS207-B\_220121A: 45      **SampType:** Sample Matrix Spike      **Batch ID:** R373694  
**Method:** SW6020      **Analysis Date:** 01/21/2022 20:43      **Prep Date:**  
**Lab ID:** B22011124-001AMS      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	0.049	0.001	0.050	0.00	99.0	88	115				

Associated Samples: **B22011131-001A**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** ICPMS207-B\_220121A: 46      **SampType:** Sample Matrix Spike Duplicate      **Batch ID:** R373694  
**Method:** SW6020      **Analysis Date:** 01/21/2022 20:49      **Prep Date:**  
**Lab ID:** B22011124-001AMSD      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	0.050	0.001	0.050	0.00	99.0	88	115	0.049	0.3	20.0	

Associated Samples: **B22011131-001A**

**Run ID: Run Order:** ICPMS207-B\_220121A: 44      **SampType:** Serial Dilution      **Batch ID:** R373694  
**Method:** SW6020      **Analysis Date:** 01/21/2022 20:36      **Prep Date:**  
**Lab ID:** B22011124-001ADIL      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	ND	0.001						0.00		10.0	

Associated Samples: **B22011131-001A**

**Run ID: Run Order:** ICPMS207-B\_220121A: 38      **SampType:** Method Blank      **Batch ID:** 163063  
**Method:** SW6020      **Analysis Date:** 01/21/2022 19:59      **Prep Date:** 01/19/2022 13:08  
**Lab ID:** MB-163063      **Units:** mg/L      **Prep Method:** SW3010A

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	ND	0.0005									

Associated Samples: **B22011131-001B**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** ICPMS207-B\_220121A: 40      **SampType:** Laboratory Control Sample      **Batch ID:** 163063  
**Method:** SW6020      **Analysis Date:** 01/21/2022 20:11      **Prep Date:** 01/19/2022 13:08  
**Lab ID:** LCS4-163063      **Units:** mg/L      **Prep Method:** SW3010A

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	0.101	0.001	0.100		101.0	88	115				

Associated Samples: **B22011131-001B**

**Run ID: Run Order:** ICPMS207-B\_220121A: 53      **SampType:** Sample Matrix Spike      **Batch ID:** 163063  
**Method:** SW6020      **Analysis Date:** 01/21/2022 21:32      **Prep Date:** 01/19/2022 14:41  
**Lab ID:** B22011124-001BMS4      **Units:** mg/L      **Prep Method:** SW3010A

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	0.103	0.001	0.100	0.00	103.0	88	115				

Associated Samples: **B22011131-001B**

**Run ID: Run Order:** ICPMS207-B\_220121A: 54      **SampType:** Sample Matrix Spike Duplicate      **Batch ID:** 163063  
**Method:** SW6020      **Analysis Date:** 01/21/2022 21:39      **Prep Date:** 01/19/2022 14:41  
**Lab ID:** B22011124-001BMSD4      **Units:** mg/L      **Prep Method:** SW3010A

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	0.103	0.001	0.100	0.00	103.0	88	115	0.103	0.3	20.0	

Associated Samples: **B22011131-001B**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** ICPMS207-B\_220121A: 82      **SampType:** Sample Matrix Spike      **Batch ID:** 163063  
**Method:** SW6020      **Analysis Date:** 01/22/2022 00:34      **Prep Date:** 01/19/2022 15:36  
**Lab ID:** B22011134-001BMS4      **Units:** mg/L      **Prep Method:** SW3010A

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	0.102	0.001	0.100	0.00	102.0	88	115				

Associated Samples: **B22011131-001B**

**Run ID: Run Order:** ICPMS207-B\_220121A: 83      **SampType:** Sample Matrix Spike Duplicate      **Batch ID:** 163063  
**Method:** SW6020      **Analysis Date:** 01/22/2022 00:40      **Prep Date:** 01/19/2022 15:36  
**Lab ID:** B22011134-001BMSD4      **Units:** mg/L      **Prep Method:** SW3010A

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	0.102	0.001	0.100	0.00	102.0	88	115	0.102	0.2	20.0	

Associated Samples: **B22011131-001B**

**Run ID: Run Order:** ICPMS207-B\_220121A: 52      **SampType:** Post Digestion/Distillation Spike      **Batch ID:** 163063  
**Method:** SW6020      **Analysis Date:** 01/21/2022 21:26      **Prep Date:** 01/19/2022 14:41  
**Lab ID:** B22011124-001BPDS1      **Units:** mg/L      **Prep Method:** SW3010A

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	0.052	0.001	0.052	0.00	101.0	80	120				

Associated Samples: **B22011131-001B**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** ICPMS207-B\_220121A: 81      **SampType:** Post Digestion/Distillation Spike      **Batch ID:** 163063  
**Method:** SW6020      **Analysis Date:** 01/22/2022 00:27      **Prep Date:** 01/19/2022 15:36  
**Lab ID:** B22011134-001BPDS1      **Units:** mg/L      **Prep Method:** SW3010A

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	0.051	0.001	0.052	0.00	100.0	80	120				

Associated Samples: **B22011131-001B**

**Run ID: Run Order:** ICPMS207-B\_220121A: 49      **SampType:** Serial Dilution      **Batch ID:** 163063  
**Method:** SW6020      **Analysis Date:** 01/21/2022 21:08      **Prep Date:** 01/19/2022 14:41  
**Lab ID:** B22011124-001BDIL      **Units:** mg/L      **Prep Method:** SW3010A

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	ND	0.001						0.00		10.0	

Associated Samples: **B22011131-001B**

**Run ID: Run Order:** ICPMS207-B\_220121A: 80      **SampType:** Serial Dilution      **Batch ID:** 163063  
**Method:** SW6020      **Analysis Date:** 01/22/2022 00:21      **Prep Date:** 01/19/2022 15:36  
**Lab ID:** B22011134-001BDIL      **Units:** mg/L      **Prep Method:** SW3010A

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	0.004	0.001						0.00		10.0	N

Associated Samples: **B22011131-001B**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** ICPMS207-B\_220121A: 131      **SampType:** Serial Dilution      **Batch ID:** 163063  
**Method:** SW6020      **Analysis Date:** 01/22/2022 11:36      **Prep Date:** 01/19/2022 15:36  
**Lab ID:** B22011134-001BDIL      **Units:** mg/L      **Prep Method:** SW3010A

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	0.004	0.001						0.00		10.0	N

Associated Samples: **B22011131-001B**

**Run ID: Run Order:** ICPMS207-B\_220121A: 132      **SampType:** Serial Dilution      **Batch ID:** 163063  
**Method:** SW6020      **Analysis Date:** 01/22/2022 11:42      **Prep Date:** 01/19/2022 14:41  
**Lab ID:** B22011124-001BDIL      **Units:** mg/L      **Prep Method:** SW3010A

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	ND	0.001						0.00		10.0	

Associated Samples: **B22011131-001B**

**Run ID: Run Order:** ICPMS207-B\_220121A: 63      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R373694  
**Method:** SW6020      **Analysis Date:** 01/21/2022 22:35      **Prep Date:**  
**Lab ID:** CCV      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	0.050	0.001	0.050		100.0	90	110				

Associated Samples: **B22011131-001A, B22011131-001B**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** ICPMS207-B\_220121A: 75      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R373694  
**Method:** SW6020      **Analysis Date:** 01/21/2022 23:50      **Prep Date:**  
**Lab ID:** CCV      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Lead	0.051	0.001	0.050		102.0	90	110				

Associated Samples: **B22011131-001A, B22011131-001B**





### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** VOA5975C.I\_220121A: 4  
**Method:** SW8260B  
**Lab ID:** MBLK012122\_

**SampType:** Method Blank  
**Analysis Date:** 01/21/2022 11:47  
**Units:** ug/L

**Batch ID:** R373695  
**Prep Date:**  
**Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Benzene	ND	0.50									
Bromobenzene	ND	0.50									
Bromochloromethane	ND	0.50									
Bromodichloromethane	ND	0.50									
Bromoform	ND	0.50									
Carbon tetrachloride	ND	0.50									
Chlorobenzene	ND	0.50									
Chlorodibromomethane	ND	0.50									
Chloroethane	ND	0.50									
Chloroform	ND	0.50									
Chloromethane	ND	0.50									
1,2-Dibromoethane	ND	0.50									
2-Chlorotoluene	ND	0.50									
Dibromomethane	ND	0.50									
1,2-Dichlorobenzene	ND	0.50									
4-Chlorotoluene	ND	0.50									
1,3-Dichlorobenzene	ND	0.50									
1,4-Dichlorobenzene	ND	0.50									
Dichlorodifluoromethane	ND	0.50									
1,1-Dichloroethane	ND	0.50									
1,2-Dichloroethane	ND	0.50									
1,1-Dichloroethene	ND	0.50									
cis-1,2-Dichloroethene	ND	0.50									
trans-1,2-Dichloroethene	ND	0.50									
1,2-Dichloropropane	ND	0.50									
1,3-Dichloropropane	ND	0.50									
2,2-Dichloropropane	ND	0.50									



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** VOA5975C.I.\_220121A: 4  
**Method:** SW8260B  
**Lab ID:** MBLK012122\_

**SampType:** Method Blank  
**Analysis Date:** 01/21/2022 11:47  
**Units:** ug/L

**Batch ID:** R373695  
**Prep Date:**  
**Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,1-Dichloropropene	ND	0.50									
cis-1,3-Dichloropropene	ND	0.50									
trans-1,3-Dichloropropene	ND	0.50									
Ethylbenzene	ND	0.50									
Methyl tert-butyl ether (MTBE)	ND	0.50									
Methyl ethyl ketone	ND	10									
Methylene chloride	ND	0.50									
Styrene	ND	0.50									
1,1,1,2-Tetrachloroethane	ND	0.50									
1,1,2,2-Tetrachloroethane	ND	0.50									
Tetrachloroethene	ND	0.50									
Toluene	ND	0.50									
1,1,1-Trichloroethane	ND	0.50									
1,1,2-Trichloroethane	ND	0.50									
Trichloroethene	ND	0.50									
Trichlorofluoromethane	ND	0.50									
1,2,3-Trichloropropane	ND	0.50									
Vinyl chloride	ND	0.50									
m+p-Xylenes	ND	0.50									
o-Xylene	ND	0.50									
Xylenes, Total	ND	0.50									
Surr: 1,2-Dichloroethane-d4	11	0.50	10		112.0	81	118				
Surr: Dibromofluoromethane	11	0.50	10		109.0	80	119				
Surr: p-Bromofluorobenzene	10	0.50	10		104.0	85	114				
Surr: Toluene-d8	10	0.50	10		104.0	89	112				

Associated Samples: **B22011131-001F, B22011131-002A**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** VOA5975C.I\_220121A: 3      **SampType:** Laboratory Control Sample      **Batch ID:** R373695  
**Method:** SW8260B      **Analysis Date:** 01/21/2022 10:53      **Prep Date:**  
**Lab ID:** LCS012122\_      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Benzene	5.1	0.50	5.0		101.0	79	120				
Bromobenzene	5.3	0.50	5.0		105.0	80	120				
Bromochloromethane	4.9	0.50	5.0		98.0	78	123				
Bromodichloromethane	5.0	0.50	5.0		99.0	79	125				
Bromoform	5.0	0.50	5.0		100.0	66	130				
Carbon tetrachloride	5.0	0.50	5.0		100.0	72	136				
Chlorobenzene	5.2	0.50	5.0		104.0	82	118				
Chlorodibromomethane	4.9	0.50	5.0		97.0	74	126				
Chloroethane	5.5	0.50	5.0		110.0	60	138				
Chloroform	4.7	0.50	5.0		95.0	79	124				
Chloromethane	4.5	0.50	5.0		91.0	50	139				
1,2-Dibromoethane	5.0	0.50	5.0		100.0	78	122				
2-Chlorotoluene	5.2	0.50	5.0		105.0	79	122				
Dibromomethane	4.9	0.50	5.0		99.0	79	123				
1,2-Dichlorobenzene	5.2	0.50	5.0		103.0	80	119				
4-Chlorotoluene	5.4	0.50	5.0		107.0	78	122				
1,3-Dichlorobenzene	5.3	0.50	5.0		106.0	80	119				
1,4-Dichlorobenzene	5.2	0.50	5.0		103.0	79	118				
Dichlorodifluoromethane	4.4	0.50	5.0		87.0	32	152				
1,1-Dichloroethane	5.2	0.50	5.0		104.0	77	125				
1,2-Dichloroethane	4.8	0.50	5.0		97.0	73	128				
1,1-Dichloroethene	5.1	0.50	5.0		103.0	71	131				
cis-1,2-Dichloroethene	5.2	0.50	5.0		104.0	78	123				
trans-1,2-Dichloroethene	5.1	0.50	5.0		101.0	75	124				
1,2-Dichloropropane	5.0	0.50	5.0		100.0	78	122				
1,3-Dichloropropane	4.9	0.50	5.0		98.0	80	119				
2,2-Dichloropropane	5.0	0.50	5.0		101.0	60	139				



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** VOA5975C.I\_220121A: 3      **SampType:** Laboratory Control Sample      **Batch ID:** R373695  
**Method:** SW8260B      **Analysis Date:** 01/21/2022 10:53      **Prep Date:**  
**Lab ID:** LCS012122\_      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,1-Dichloropropene	4.9	0.50	5.0		97.0	79	125				
cis-1,3-Dichloropropene	4.8	0.50	5.0		95.0	75	124				
trans-1,3-Dichloropropene	5.2	0.50	5.0		104.0	73	127				
Ethylbenzene	5.0	0.50	5.0		100.0	79	121				
Methyl tert-butyl ether (MTBE)	5.1	0.50	5.0		102.0	71	124				
Methyl ethyl ketone	49	10	50		98.0	56	143				
Methylene chloride	4.8	0.50	5.0		96.0	74	124				
Styrene	5.1	0.50	5.0		102.0	78	123				
1,1,1,2-Tetrachloroethane	4.9	0.50	5.0		98.0	78	124				
1,1,2,2-Tetrachloroethane	5.0	0.50	5.0		99.0	71	121				
Tetrachloroethene	4.9	0.50	5.0		99.0	74	129				
Toluene	5.1	0.50	5.0		102.0	80	121				
1,1,1-Trichloroethane	5.0	0.50	5.0		99.0	74	131				
1,1,2-Trichloroethane	5.1	0.50	5.0		102.0	80	119				
Trichloroethene	5.0	0.50	5.0		100.0	79	123				
Trichlorofluoromethane	4.5	0.50	5.0		90.0	65	141				
1,2,3-Trichloropropane	5.1	0.50	5.0		101.0	73	125				
Vinyl chloride	4.8	0.50	5.0		95.0	58	137				
m+p-Xylenes	9.9	0.50	10		99.0	80	121				
o-Xylene	5.1	0.50	5.0		101.0	78	122				
Xylenes, Total	15	0.50	15		100.0	79	121				
Surr: 1,2-Dichloroethane-d4	11	0.50	10		109.0	81	118				
Surr: Dibromofluoromethane	11	0.50	10		108.0	80	119				
Surr: p-Bromofluorobenzene	10	0.50	10		103.0	85	114				
Surr: Toluene-d8	11	0.50	10		108.0	89	112				

Associated Samples: **B22011131-001F, B22011131-002A**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** VOA5975C.I\_220121A: 22

**SampType:** Sample Matrix Spike

**Batch ID:** R373695

**Method:** SW8260B

**Analysis Date:** 01/21/2022 20:15

**Prep Date:**

**Lab ID:** B22011125-001FMS

**Units:** ug/L

**Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Benzene	5.2	0.50	5.0	0.0	104.0	79	120				
Bromobenzene	5.3	0.50	5.0	0.0	106.0	80	120				
Bromochloromethane	4.8	0.50	5.0	0.0	97.0	78	123				
Bromodichloromethane	5.3	0.50	5.0	0.0	105.0	79	125				
Bromoform	5.1	0.50	5.0	0.0	102.0	66	130				
Carbon tetrachloride	5.2	0.50	5.0	0.0	104.0	72	136				
Chlorobenzene	5.4	0.50	5.0	0.0	107.0	82	118				
Chlorodibromomethane	5.2	0.50	5.0	0.0	104.0	74	126				
Chloroethane	5.6	0.50	5.0	0.0	112.0	60	138				
Chloroform	4.8	0.50	5.0	0.0	97.0	79	124				
Chloromethane	4.7	0.50	5.0	0.0	94.0	50	139				
1,2-Dibromoethane	5.2	0.50	5.0	0.0	105.0	78	122				
2-Chlorotoluene	5.4	0.50	5.0	0.0	107.0	79	122				
Dibromomethane	5.1	0.50	5.0	0.0	103.0	79	123				
1,2-Dichlorobenzene	5.2	0.50	5.0	0.0	104.0	80	119				
4-Chlorotoluene	5.5	0.50	5.0	0.0	111.0	78	122				
1,3-Dichlorobenzene	5.4	0.50	5.0	0.0	107.0	80	119				
1,4-Dichlorobenzene	5.3	0.50	5.0	0.0	106.0	79	118				
Dichlorodifluoromethane	4.6	0.50	5.0	0.0	92.0	32	152				
1,1-Dichloroethane	5.4	0.50	5.0	0.0	108.0	77	125				
1,2-Dichloroethane	4.8	0.50	5.0	0.0	96.0	73	128				
1,1-Dichloroethene	5.3	0.50	5.0	0.0	107.0	71	131				
cis-1,2-Dichloroethene	5.2	0.50	5.0	0.0	103.0	78	123				
trans-1,2-Dichloroethene	5.0	0.50	5.0	0.0	101.0	75	124				
1,2-Dichloropropane	5.3	0.50	5.0	0.0	105.0	78	122				
1,3-Dichloropropane	5.0	0.50	5.0	0.0	99.0	80	119				
2,2-Dichloropropane	5.1	0.50	5.0	0.0	102.0	60	139				



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** VOA5975C.I\_220121A: 22      **SampType:** Sample Matrix Spike      **Batch ID:** R373695  
**Method:** SW8260B      **Analysis Date:** 01/21/2022 20:15      **Prep Date:**  
**Lab ID:** B22011125-001FMS      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,1-Dichloropropene	5.0	0.50	5.0	0.0	101.0	79	125				
cis-1,3-Dichloropropene	4.9	0.50	5.0	0.0	99.0	75	124				
trans-1,3-Dichloropropene	5.2	0.50	5.0	0.0	105.0	73	127				
Ethylbenzene	5.3	0.50	5.0	0.0	106.0	79	121				
Methyl tert-butyl ether (MTBE)	5.2	0.50	5.0	0.0	105.0	71	124				
Methyl ethyl ketone	46	10	50	0.0	92.0	56	143				
Methylene chloride	4.9	0.50	5.0	0.0	99.0	74	124				
Styrene	5.3	0.50	5.0	0.0	107.0	78	123				
1,1,1,2-Tetrachloroethane	5.2	0.50	5.0	0.0	105.0	78	124				
1,1,2,2-Tetrachloroethane	5.3	0.50	5.0	0.0	105.0	71	121				
Tetrachloroethene	5.3	0.50	5.0	0.0	106.0	74	129				
Toluene	5.4	0.50	5.0	0.0	109.0	80	121				
1,1,1-Trichloroethane	5.2	0.50	5.0	0.0	104.0	74	131				
1,1,2-Trichloroethane	5.2	0.50	5.0	0.0	103.0	80	119				
Trichloroethene	5.3	0.50	5.0	0.0	106.0	79	123				
Trichlorofluoromethane	5.1	0.50	5.0	0.0	101.0	65	141				
1,2,3-Trichloropropane	4.8	0.50	5.0	0.0	97.0	73	125				
Vinyl chloride	4.8	0.50	5.0	0.0	97.0	58	137				
m+p-Xylenes	10	0.50	10	0.0	104.0	80	121				
o-Xylene	5.3	0.50	5.0	0.0	106.0	78	122				
Xylenes, Total	16	0.50	15	0.0	105.0	79	121				
Surr: 1,2-Dichloroethane-d4	10	0.50	10	0.0	100.0	81	118				
Surr: Dibromofluoromethane	9.7	0.50	10	0.0	97.0	80	119				
Surr: p-Bromofluorobenzene	9.6	0.50	10	0.0	96.0	85	114				
Surr: Toluene-d8	10	0.50	10	0.0	102.0	89	112				

Associated Samples: **B22011131-001F, B22011131-002A**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** VOA5975C.I\_220121A: 23      **SampType:** Sample Matrix Spike Duplicate      **Batch ID:** R373695  
**Method:** SW8260B      **Analysis Date:** 01/21/2022 20:42      **Prep Date:**  
**Lab ID:** B22011125-001FMSD      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Benzene	5.3	0.50	5.0	0.0	107.0	79	120	5.2	2.9	20.0	
Bromobenzene	5.4	0.50	5.0	0.0	108.0	80	120	5.3	1.8	20.0	
Bromochloromethane	5.1	0.50	5.0	0.0	102.0	78	123	4.8	5.4	20.0	
Bromodichloromethane	5.3	0.50	5.0	0.0	106.0	79	125	5.3	0.2	20.0	
Bromoform	5.0	0.50	5.0	0.0	101.0	66	130	5.1	1.7	20.0	
Carbon tetrachloride	5.4	0.50	5.0	0.0	108.0	72	136	5.2	3.7	20.0	
Chlorobenzene	5.4	0.50	5.0	0.0	108.0	82	118	5.4	0.6	20.0	
Chlorodibromomethane	5.1	0.50	5.0	0.0	102.0	74	126	5.2	1.2	20.0	
Chloroethane	5.4	0.50	5.0	0.0	107.0	60	138	5.6	4.4	20.0	
Chloroform	4.8	0.50	5.0	0.0	96.0	79	124	4.8	0.7	20.0	
Chloromethane	4.8	0.50	5.0	0.0	96.0	50	139	4.7	2.1	20.0	
1,2-Dibromoethane	5.0	0.50	5.0	0.0	101.0	78	122	5.2	3.5	20.0	
2-Chlorotoluene	5.5	0.50	5.0	0.0	110.0	79	122	5.4	2.8	20.0	
Dibromomethane	5.1	0.50	5.0	0.0	103.0	79	123	5.1	0.1	20.0	
1,2-Dichlorobenzene	5.4	0.50	5.0	0.0	108.0	80	119	5.2	2.9	20.0	
4-Chlorotoluene	5.6	0.50	5.0	0.0	112.0	78	122	5.5	1.3	20.0	
1,3-Dichlorobenzene	5.5	0.50	5.0	0.0	110.0	80	119	5.4	2.3	20.0	
1,4-Dichlorobenzene	5.3	0.50	5.0	0.0	107.0	79	118	5.3	0.9	20.0	
Dichlorodifluoromethane	4.7	0.50	5.0	0.0	95.0	32	152	4.6	3.4	20.0	
1,1-Dichloroethane	5.4	0.50	5.0	0.0	108.0	77	125	5.4	0.4	20.0	
1,2-Dichloroethane	4.9	0.50	5.0	0.0	98.0	73	128	4.8	2.0	20.0	
1,1-Dichloroethene	5.4	0.50	5.0	0.0	107.0	71	131	5.3	0.3	20.0	
cis-1,2-Dichloroethene	5.2	0.50	5.0	0.0	105.0	78	123	5.2	1.6	20.0	
trans-1,2-Dichloroethene	5.3	0.50	5.0	0.0	106.0	75	124	5.0	5.5	20.0	
1,2-Dichloropropane	5.1	0.50	5.0	0.0	102.0	78	122	5.3	3.4	20.0	
1,3-Dichloropropane	4.9	0.50	5.0	0.0	98.0	80	119	5.0	1.2	20.0	
2,2-Dichloropropane	5.1	0.50	5.0	0.0	101.0	60	139	5.1	0.4	20.0	



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** VOA5975C.I\_220121A: 23      **SampType:** Sample Matrix Spike Duplicate      **Batch ID:** R373695  
**Method:** SW8260B      **Analysis Date:** 01/21/2022 20:42      **Prep Date:**  
**Lab ID:** B22011125-001FMSD      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,1-Dichloropropene	5.2	0.50	5.0	0.0	104.0	79	125	5.0	3.2	20.0	
cis-1,3-Dichloropropene	4.7	0.50	5.0	0.0	93.0	75	124	4.9	5.7	20.0	
trans-1,3-Dichloropropene	5.2	0.50	5.0	0.0	104.0	73	127	5.2	1.0	20.0	
Ethylbenzene	5.3	0.50	5.0	0.0	105.0	79	121	5.3	0.6	20.0	
Methyl tert-butyl ether (MTBE)	5.1	0.50	5.0	0.0	103.0	71	124	5.2	2.0	20.0	
Methyl ethyl ketone	51	10	50	0.0	102.0	56	143	46	11.0	20.0	
Methylene chloride	5.0	0.50	5.0	0.0	99.0	74	124	4.9	0.5	20.0	
Styrene	5.3	0.50	5.0	0.0	106.0	78	123	5.3	0.8	20.0	
1,1,1,2-Tetrachloroethane	5.1	0.50	5.0	0.0	102.0	78	124	5.2	2.2	20.0	
1,1,2,2-Tetrachloroethane	5.3	0.50	5.0	0.0	106.0	71	121	5.3	1.0	20.0	
Tetrachloroethene	5.3	0.50	5.0	0.0	107.0	74	129	5.3	1.0	20.0	
Toluene	5.4	0.50	5.0	0.0	108.0	80	121	5.4	0.4	20.0	
1,1,1-Trichloroethane	5.3	0.50	5.0	0.0	106.0	74	131	5.2	1.6	20.0	
1,1,2-Trichloroethane	5.1	0.50	5.0	0.0	102.0	80	119	5.2	0.9	20.0	
Trichloroethene	5.3	0.50	5.0	0.0	106.0	79	123	5.3	0.5	20.0	
Trichlorofluoromethane	5.0	0.50	5.0	0.0	101.0	65	141	5.1	0.5	20.0	
1,2,3-Trichloropropane	5.3	0.50	5.0	0.0	106.0	73	125	4.8	9.6	20.0	
Vinyl chloride	5.2	0.50	5.0	0.0	103.0	58	137	4.8	6.7	20.0	
m+p-Xylenes	10	0.50	10	0.0	104.0	80	121	10	0.0	20.0	
o-Xylene	5.3	0.50	5.0	0.0	106.0	78	122	5.3	0.0	20.0	
Xylenes, Total	16	0.50	15	0.0	105.0	79	121	16	0.0	20.0	
Surr: 1,2-Dichloroethane-d4	10	0.50	10	0.0	105.0	81	118	0.0			
Surr: Dibromofluoromethane	10	0.50	10	0.0	103.0	80	119	0.0			
Surr: p-Bromofluorobenzene	10	0.50	10	0.0	104.0	85	114	0.0			
Surr: Toluene-d8	11	0.50	10	0.0	106.0	89	112	0.0			

Associated Samples: B22011131-001F, B22011131-002A





### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** VOA5975C.I\_220121A: 2      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R373695  
**Method:** SW8260B      **Analysis Date:** 01/21/2022 10:17      **Prep Date:**  
**Lab ID:** CCV012122\_      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Benzene	5.1	0.50	5.0		102.0	80	120				
Bromobenzene	5.1	0.50	5.0		102.0	80	120				
Bromochloromethane	5.0	0.50	5.0		100.0	80	120				
Bromodichloromethane	4.9	0.50	5.0		98.0	80	120				
Bromoform	4.9	0.50	5.0		99.0	80	120				
Carbon tetrachloride	5.0	0.50	5.0		100.0	80	120				
Chlorobenzene	5.0	0.50	5.0		100.0	80	120				
Chlorodibromomethane	4.9	0.50	5.0		99.0	80	120				
Chloroethane	5.7	0.50	5.0		115.0	80	120				
Chloroform	4.8	0.50	5.0		96.0	80	120				
Chloromethane	4.9	0.50	5.0		99.0	80	120				
1,2-Dibromoethane	5.0	0.50	5.0		100.0	80	120				
2-Chlorotoluene	5.0	0.50	5.0		99.0	80	120				
Dibromomethane	5.0	0.50	5.0		101.0	80	120				
1,2-Dichlorobenzene	4.9	0.50	5.0		98.0	80	120				
4-Chlorotoluene	5.1	0.50	5.0		101.0	80	120				
1,3-Dichlorobenzene	4.9	0.50	5.0		97.0	80	120				
1,4-Dichlorobenzene	4.9	0.50	5.0		98.0	80	120				
Dichlorodifluoromethane	4.7	0.50	5.0		94.0	80	120				
1,1-Dichloroethane	5.0	0.50	5.0		100.0	80	120				
1,2-Dichloroethane	4.9	0.50	5.0		98.0	80	120				
1,1-Dichloroethene	4.9	0.50	5.0		99.0	80	120				
cis-1,2-Dichloroethene	5.1	0.50	5.0		102.0	80	120				
trans-1,2-Dichloroethene	4.9	0.50	5.0		98.0	80	120				
1,2-Dichloropropane	5.1	0.50	5.0		101.0	80	120				
1,3-Dichloropropane	5.0	0.50	5.0		100.0	80	120				
2,2-Dichloropropane	5.2	0.50	5.0		103.0	80	120				



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** VOA5975C.I\_220121A: 2      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R373695  
**Method:** SW8260B      **Analysis Date:** 01/21/2022 10:17      **Prep Date:**  
**Lab ID:** CCV012122\_      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,1-Dichloropropene	5.1	0.50	5.0		101.0	80	120				
cis-1,3-Dichloropropene	5.0	0.50	5.0		100.0	80	120				
trans-1,3-Dichloropropene	5.2	0.50	5.0		104.0	80	120				
Ethylbenzene	4.9	0.50	5.0		97.0	80	120				
Methyl tert-butyl ether (MTBE)	4.9	0.50	5.0		98.0	80	120				
Methyl ethyl ketone	48	10	50		95.0	80	120				
Methylene chloride	4.9	0.50	5.0		97.0	80	120				
Styrene	5.0	0.50	5.0		99.0	80	120				
1,1,1,2-Tetrachloroethane	4.9	0.50	5.0		98.0	80	120				
1,1,2,2-Tetrachloroethane	5.0	0.50	5.0		100.0	80	120				
Tetrachloroethene	4.8	0.50	5.0		95.0	80	120				
Toluene	5.1	0.50	5.0		102.0	80	120				
1,1,1-Trichloroethane	5.0	0.50	5.0		99.0	80	120				
1,1,2-Trichloroethane	5.1	0.50	5.0		103.0	80	120				
Trichloroethene	5.0	0.50	5.0		100.0	80	120				
Trichlorofluoromethane	4.8	0.50	5.0		96.0	80	120				
1,2,3-Trichloropropane	4.9	0.50	5.0		99.0	80	120				
Vinyl chloride	4.8	0.50	5.0		97.0	80	120				
m+p-Xylenes	10	0.50	10		100.0	80	120				
o-Xylene	4.9	0.50	5.0		98.0	80	120				
Xylenes, Total	15	0.50	15		99.0	80	120				
Surr: 1,2-Dichloroethane-d4	11	0.50	10		109.0	80	120				
Surr: Dibromofluoromethane	10	0.50	10		102.0	80	120				
Surr: p-Bromofluorobenzene	10	0.50	10		100.0	80	120				
Surr: Toluene-d8	10	0.50	10		103.0	80	120				

Associated Samples: B22011131-001F, B22011131-002A



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** VOA5975C.I\_220121A: 24      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R373695  
**Method:** SW8260B      **Analysis Date:** 01/21/2022 21:37      **Prep Date:**  
**Lab ID:** CCV012122\_Closing      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Benzene	5.3	0.50	5.0		106.0	50	150				
Bromobenzene	5.3	0.50	5.0		106.0	50	150				
Bromochloromethane	5.3	0.50	5.0		106.0	50	150				
Bromodichloromethane	5.3	0.50	5.0		106.0	50	150				
Bromoform	5.1	0.50	5.0		101.0	50	150				
Carbon tetrachloride	5.4	0.50	5.0		108.0	50	150				
Chlorobenzene	5.4	0.50	5.0		107.0	50	150				
Chlorodibromomethane	5.2	0.50	5.0		104.0	50	150				
Chloroethane	4.3	0.50	5.0		85.0	50	150				
Chloroform	5.1	0.50	5.0		102.0	50	150				
Chloromethane	4.8	0.50	5.0		95.0	50	150				
1,2-Dibromoethane	5.1	0.50	5.0		102.0	50	150				
2-Chlorotoluene	5.4	0.50	5.0		108.0	50	150				
Dibromomethane	5.2	0.50	5.0		104.0	50	150				
1,2-Dichlorobenzene	5.3	0.50	5.0		106.0	50	150				
4-Chlorotoluene	5.5	0.50	5.0		110.0	50	150				
1,3-Dichlorobenzene	5.2	0.50	5.0		105.0	50	150				
1,4-Dichlorobenzene	5.3	0.50	5.0		106.0	50	150				
Dichlorodifluoromethane	4.9	0.50	5.0		97.0	50	150				
1,1-Dichloroethane	5.3	0.50	5.0		106.0	50	150				
1,2-Dichloroethane	4.9	0.50	5.0		98.0	50	150				
1,1-Dichloroethene	5.2	0.50	5.0		104.0	50	150				
cis-1,2-Dichloroethene	5.2	0.50	5.0		103.0	50	150				
trans-1,2-Dichloroethene	5.1	0.50	5.0		103.0	50	150				
1,2-Dichloropropane	5.3	0.50	5.0		105.0	50	150				
1,3-Dichloropropane	5.3	0.50	5.0		106.0	50	150				
2,2-Dichloropropane	5.0	0.50	5.0		99.0	50	150				



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** VOA5975C.I\_220121A: 24      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R373695  
**Method:** SW8260B      **Analysis Date:** 01/21/2022 21:37      **Prep Date:**  
**Lab ID:** CCV012122\_Closing      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,1-Dichloropropene	5.4	0.50	5.0		108.0	50	150				
cis-1,3-Dichloropropene	5.1	0.50	5.0		102.0	50	150				
trans-1,3-Dichloropropene	5.3	0.50	5.0		105.0	50	150				
Ethylbenzene	5.3	0.50	5.0		107.0	50	150				
Methyl tert-butyl ether (MTBE)	4.7	0.50	5.0		93.0	50	150				
Methyl ethyl ketone	49	10	50		98.0	50	150				
Methylene chloride	5.0	0.50	5.0		101.0	50	150				
Styrene	5.4	0.50	5.0		108.0	50	150				
1,1,1,2-Tetrachloroethane	5.3	0.50	5.0		106.0	50	150				
1,1,2,2-Tetrachloroethane	5.1	0.50	5.0		102.0	50	150				
Tetrachloroethene	5.4	0.50	5.0		107.0	50	150				
Toluene	5.5	0.50	5.0		110.0	50	150				
1,1,1-Trichloroethane	5.3	0.50	5.0		106.0	50	150				
1,1,2-Trichloroethane	5.0	0.50	5.0		101.0	50	150				
Trichloroethene	5.5	0.50	5.0		109.0	50	150				
Trichlorofluoromethane	5.0	0.50	5.0		101.0	50	150				
1,2,3-Trichloropropane	5.0	0.50	5.0		100.0	50	150				
Vinyl chloride	4.9	0.50	5.0		98.0	50	150				
m+p-Xylenes	11	0.50	10		105.0	50	150				
o-Xylene	5.4	0.50	5.0		108.0	50	150				
Xylenes, Total	16	0.50	15		106.0	50	150				
Surr: 1,2-Dichloroethane-d4	11	0.50	10		106.0	50	150				
Surr: Dibromofluoromethane	10	0.50	10		104.0	50	150				
Surr: p-Bromofluorobenzene	10	0.50	10		102.0	50	150				
Surr: Toluene-d8	11	0.50	10		107.0	50	150				

Associated Samples: B22011131-001F, B22011131-002A



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** GECD.I\_220121A: 35      **SampType:** Method Blank      **Batch ID:** 163129  
**Method:** SW8011      **Analysis Date:** 01/22/2022 05:36      **Prep Date:** 01/21/2022 07:48  
**Lab ID:** MB-163129      **Units:** ug/L      **Prep Method:** SW8011

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,2-Dibromoethane	ND	0.0050									
Surr: 1,1,1,2-Tetrachloroethane	0.094	0.020	0.10		94.0	70	130				

Associated Samples: **B22011131-001H, B22011131-004A**

**Run ID: Run Order:** GECD.I\_220121A: 36      **SampType:** Laboratory Control Sample      **Batch ID:** 163129  
**Method:** SW8011      **Analysis Date:** 01/22/2022 05:56      **Prep Date:** 01/21/2022 07:48  
**Lab ID:** LCS-163129      **Units:** ug/L      **Prep Method:** SW8011

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,2-Dibromoethane	0.24	0.010	0.25		94.0	60	140				
Surr: 1,1,1,2-Tetrachloroethane	0.093	0.020	0.10		93.0	70	130				

Associated Samples: **B22011131-001H, B22011131-004A**

**Run ID: Run Order:** GECD.I\_220121A: 37      **SampType:** Laboratory Control Sample      **Batch ID:** 163129  
**Method:** SW8011      **Analysis Date:** 01/22/2022 06:16      **Prep Date:** 01/21/2022 07:48  
**Lab ID:** LCS1-163129      **Units:** ug/L      **Prep Method:** SW8011

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,2-Dibromoethane	0.10	0.010	0.10		100.0	60	140				
Surr: 1,1,1,2-Tetrachloroethane	0.093	0.020	0.10		93.0	70	130				

Associated Samples: **B22011131-001H, B22011131-004A**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** GECD.I\_220121A: 50      **SampType:** Sample Matrix Spike      **Batch ID:** 163129  
**Method:** SW8011      **Analysis Date:** 01/22/2022 10:55      **Prep Date:** 01/21/2022 07:48  
**Lab ID:** B22011131-001HMS      **Units:** ug/L      **Prep Method:** SW8011

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,2-Dibromoethane	0.23	0.010	0.24	0.0	93.0	60	140				
Surr: 1,1,1,2-Tetrachloroethane	0.085	0.020	0.097	0.0	88.0	70	130				

Associated Samples: **B22011131-001H, B22011131-004A**

**Run ID: Run Order:** GECD.I\_220121A: 51      **SampType:** Sample Matrix Spike Duplicate      **Batch ID:** 163129  
**Method:** SW8011      **Analysis Date:** 01/22/2022 11:15      **Prep Date:** 01/21/2022 07:48  
**Lab ID:** B22011131-001HMSD      **Units:** ug/L      **Prep Method:** SW8011

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,2-Dibromoethane	0.22	0.010	0.24	0.0	92.0	60	140	0.23	1.6	20.0	
Surr: 1,1,1,2-Tetrachloroethane	0.088	0.020	0.097	0.0	91.0	70	130	0.0			

Associated Samples: **B22011131-001H, B22011131-004A**

**Run ID: Run Order:** GECD.I\_220121A: 34      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** 163128  
**Method:** SW8011      **Analysis Date:** 01/22/2022 05:16      **Prep Date:** 01/21/2022 07:45  
**Lab ID:** CAL3-163128      **Units:** ug/L      **Prep Method:** SW8011

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,2-Dibromoethane	0.098	0.010	0.10		98.0	80	120				
Surr: 1,1,1,2-Tetrachloroethane	0.091	0.020	0.10		91.0	80	120				

Associated Samples: **B22011131-001H, B22011131-004A**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** GECD.I\_220121A: 52      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** 163129  
**Method:** SW8011      **Analysis Date:** 01/22/2022 11:55      **Prep Date:** 01/21/2022 07:48  
**Lab ID:** CK5-163129      **Units:** ug/L      **Prep Method:** SW8011

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,2-Dibromoethane	0.39	0.010	0.40		96.0	80	120				
Surr: 1,1,1,2-Tetrachloroethane	0.41	0.020	0.40		104.0	80	120				

Associated Samples: **B22011131-001H, B22011131-004A**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** PE 1\_220120A: 4      **SampType:** Method Blank      **Batch ID:** R373498  
**Method:** SW8015C      **Analysis Date:** 01/20/2022 11:24      **Prep Date:**  
**Lab ID:** MBLK\_0120PE106r      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
C6 to C10	ND	10									
Total Purgeable Hydrocarbons	ND	10									
Surr: Trifluorotoluene	20	1.0	25		80.0	70	130				

Associated Samples: **B22011131-001G, B22011131-003A**

**Run ID: Run Order:** PE 1\_220120A: 19      **SampType:** Method Blank      **Batch ID:** R373498  
**Method:** SW8015C      **Analysis Date:** 01/20/2022 22:16      **Prep Date:**  
**Lab ID:** MBLK\_0120PE125r      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
C6 to C10	ND	10									
Total Purgeable Hydrocarbons	ND	10									
Surr: Trifluorotoluene	20	1.0	25		81.0	70	130				

Associated Samples: **B22011131-001G, B22011131-003A**

**Run ID: Run Order:** PE 1\_220120A: 31      **SampType:** Method Blank      **Batch ID:** R373498  
**Method:** SW8015C      **Analysis Date:** 01/21/2022 09:08      **Prep Date:**  
**Lab ID:** MBLK\_0120PE144r      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
C6 to C10	ND	10									
Total Purgeable Hydrocarbons	ND	10									
Surr: Trifluorotoluene	20	1.0	25		81.0	70	130				

Associated Samples: **B22011131-001G, B22011131-003A**





### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** PE 1\_220120A: 3      **SampType:** Laboratory Control Sample      **Batch ID:** R373498  
**Method:** SW8015C      **Analysis Date:** 01/20/2022 10:49      **Prep Date:**  
**Lab ID:** LCS\_0120PE105r      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
C6 to C10	160	20	170		94.0	78	122				
Total Purgeable Hydrocarbons	189	20	200		95.0	70	130				
Surr: Trifluorotoluene	23	1.0	25		91.0	70	130				

Associated Samples: **B22011131-001G, B22011131-003A**

**Run ID: Run Order:** PE 1\_220120A: 18      **SampType:** Laboratory Control Sample      **Batch ID:** R373498  
**Method:** SW8015C      **Analysis Date:** 01/20/2022 21:42      **Prep Date:**  
**Lab ID:** LCS\_0120PE124r      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
C6 to C10	151	20	170		89.0	78	122				
Total Purgeable Hydrocarbons	178	20	200		89.0	70	130				
Surr: Trifluorotoluene	22	1.0	25		88.0	70	130				

Associated Samples: **B22011131-001G, B22011131-003A**

**Run ID: Run Order:** PE 1\_220120A: 30      **SampType:** Laboratory Control Sample      **Batch ID:** R373498  
**Method:** SW8015C      **Analysis Date:** 01/21/2022 08:34      **Prep Date:**  
**Lab ID:** LCS\_0120PE143r      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
C6 to C10	156	20	170		92.0	78	122				
Total Purgeable Hydrocarbons	184	20	200		92.0	70	130				
Surr: Trifluorotoluene	22	1.0	25		89.0	70	130				

Associated Samples: **B22011131-001G, B22011131-003A**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** PE 1\_220120A: 14      **SampType:** Sample Matrix Spike      **Batch ID:** R373498  
**Method:** SW8015C      **Analysis Date:** 01/20/2022 18:50      **Prep Date:**  
**Lab ID:** B22011124-001GMS      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
C6 to C10	157	20	170	0.0	92.0	78	122				
Total Purgeable Hydrocarbons	186	20	200	0.0	93.0	70	130				
Surr: Trifluorotoluene	23	1.0	25	0.0	91.0	70	130				

Associated Samples: **B22011131-001G, B22011131-003A**

**Run ID: Run Order:** PE 1\_220120A: 15      **SampType:** Sample Matrix Spike Duplicate      **Batch ID:** R373498  
**Method:** SW8015C      **Analysis Date:** 01/20/2022 19:24      **Prep Date:**  
**Lab ID:** B22011124-001GMSD      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
C6 to C10	157	20	170	0.0	92.0	78	122	157	0.0	20.0	
Total Purgeable Hydrocarbons	185	20	200	0.0	93.0	70	130	186	0.2	20.0	
Surr: Trifluorotoluene	23	1.0	25	0.0	91.0	70	130	0.0			

Associated Samples: **B22011131-001G, B22011131-003A**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** GCFID-HP5-B\_220122A: 4      **SampType:** Method Blank      **Batch ID:** 163074  
**Method:** SW8015C      **Analysis Date:** 01/22/2022 15:56      **Prep Date:** 01/19/2022 16:29  
**Lab ID:** MB-163074      **Units:** mg/L      **Prep Method:** SW3520C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Diesel Range Organics (C10 to C24)	ND	0.15									
Oil Range Hydrocarbons (C24 to C40)	ND	0.15									
Total Extractable Hydrocarbons	ND	0.15									
Surr: o-Terphenyl	0.19	0.0020	0.20		95.0	56	125				
Surr: n-Triacontane	0.099	0.0020	0.10		99.0	50	150				

Associated Samples: **B22011131-001D**

**Run ID: Run Order:** GCFID-HP5-B\_220122A: 3      **SampType:** Laboratory Control Sample      **Batch ID:** 163074  
**Method:** SW8015C      **Analysis Date:** 01/22/2022 15:13      **Prep Date:** 01/19/2022 16:29  
**Lab ID:** LCS-163074      **Units:** mg/L      **Prep Method:** SW3520C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Diesel Range Organics (C10 to C24)	12	0.30	15		83.0	36	132				
Total Extractable Hydrocarbons	13	0.30	15		89.0	60	132				
Surr: o-Terphenyl	0.19	0.0020	0.20		97.0	56	125				

Associated Samples: **B22011131-001D**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** GCFID-HP5-B\_220122A: 24      **SampType:** Laboratory Control Sample      **Batch ID:** 163074  
**Method:** SW8015C      **Analysis Date:** 01/23/2022 16:52      **Prep Date:** 01/19/2022 16:30  
**Lab ID:** LCS-163074-RRO      **Units:** mg/L      **Prep Method:** SW3520C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
TEH(Oil Range)	5.0	0.30	5.0		100.0	41	113				
Surr: n-Triacontane	0.098	0.0020	0.10		98.0	50	150				

Associated Samples: **B22011131-001D**

**Run ID: Run Order:** GCFID-HP5-B\_220122A: 10      **SampType:** Sample Matrix Spike      **Batch ID:** 163074  
**Method:** SW8015C      **Analysis Date:** 01/22/2022 20:56      **Prep Date:** 01/19/2022 16:30  
**Lab ID:** B22011136-001DMS      **Units:** mg/L      **Prep Method:** SW3520C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Diesel Range Organics (C10 to C24)	12	0.30	15	0.041	81.0	36	132				
Total Extractable Hydrocarbons	13	0.30	15	0.13	86.0	60	132				
Surr: o-Terphenyl	0.18	0.0020	0.20	0.0	91.0	56	125				

Associated Samples: **B22011131-001D**

**Run ID: Run Order:** GCFID-HP5-B\_220122A: 11      **SampType:** Sample Matrix Spike Duplicate      **Batch ID:** 163074  
**Method:** SW8015C      **Analysis Date:** 01/22/2022 21:39      **Prep Date:** 01/19/2022 16:30  
**Lab ID:** B22011136-001DMSD      **Units:** mg/L      **Prep Method:** SW3520C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Diesel Range Organics (C10 to C24)	13	0.30	15	0.041	86.0	36	132	12	6.9	20.0	
Total Extractable Hydrocarbons	14	0.30	15	0.13	92.0	60	132	13	7.2	20.0	
Surr: o-Terphenyl	0.19	0.0020	0.20	0.0	97.0	56	125	0.0			

Associated Samples: **B22011131-001D**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** GCFID-HP5-B\_220122A: 25      **SampType:** Sample Matrix Spike      **Batch ID:** 163074  
**Method:** SW8015C      **Analysis Date:** 01/23/2022 18:18      **Prep Date:** 01/19/2022 16:30  
**Lab ID:** B22011136-001DMS-RRO      **Units:** mg/L      **Prep Method:** SW3520C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
TEH(Oil Range)	4.8	0.30	5.0	0.13	93.0	41	113				
Surr: n-Triacontane	0.094	0.0020	0.10	0.0	94.0	50	150				

Associated Samples: **B22011131-001D**

**Run ID: Run Order:** GCFID-HP5-B\_220122A: 26      **SampType:** Sample Matrix Spike Duplicate      **Batch ID:** 163074  
**Method:** SW8015C      **Analysis Date:** 01/23/2022 19:01      **Prep Date:** 01/19/2022 16:30  
**Lab ID:** B22011136-001DMSD-RRO      **Units:** mg/L      **Prep Method:** SW3520C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
TEH(Oil Range)	4.9	0.30	4.9	0.13	98.0	41	113	4.8	1.9	20.0	
Surr: n-Triacontane	0.088	0.0020	0.097	0.0	91.0	50	150	0.0			

Associated Samples: **B22011131-001D**

**Run ID: Run Order:** PE 1\_220120A: 2      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R373498  
**Method:** SW8015C      **Analysis Date:** 01/20/2022 10:15      **Prep Date:**  
**Lab ID:** CCV\_0120PE104r      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
C6 to C10	184	20	168		110.0	80	120				
Total Purgeable Hydrocarbons	221	20	200		110.0	80	120				
Surr: Trifluorotoluene	24	1.0	25		96.0	80	120				

Associated Samples: **B22011131-001G, B22011131-003A**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** PE 1\_220120A: 17      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R373498  
**Method:** SW8015C      **Analysis Date:** 01/20/2022 21:07      **Prep Date:**  
**Lab ID:** CCV\_0120PE123r      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
C6 to C10	180	20	168		107.0	80	120				
Total Purgeable Hydrocarbons	216	20	200		108.0	80	120				
Surr: Trifluorotoluene	23	1.0	25		94.0	80	120				

Associated Samples: **B22011131-001G, B22011131-003A**

**Run ID: Run Order:** PE 1\_220120A: 29      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R373498  
**Method:** SW8015C      **Analysis Date:** 01/21/2022 07:59      **Prep Date:**  
**Lab ID:** CCV\_0120PE142r      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
C6 to C10	169	20	168		101.0	80	120				
Total Purgeable Hydrocarbons	203	20	200		101.0	80	120				
Surr: Trifluorotoluene	22	1.0	25		90.0	80	120				

Associated Samples: **B22011131-001G, B22011131-003A**

**Run ID: Run Order:** PE 1\_220120A: 45      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R373498  
**Method:** SW8015C      **Analysis Date:** 01/21/2022 18:51      **Prep Date:**  
**Lab ID:** CCV\_0120PE161r      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
C6 to C10	181	20	168		107.0	80	120				
Total Purgeable Hydrocarbons	218	20	200		109.0	80	120				
Surr: Trifluorotoluene	22	1.0	25		89.0	80	120				

Associated Samples: **B22011131-001G, B22011131-003A**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** GCFID-HP5-B\_220122A: 12      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R373590  
**Method:** SW8015C      **Analysis Date:** 01/22/2022 23:05      **Prep Date:**  
**Lab ID:** CCV\_0122HP519r-W      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
TEH(Oil Range)	4.6	0.30	5.0		93.0	80	120				
Surr: n-Triacontane	0.21	0.0020	0.20		105.0	80	120				

Associated Samples: **B22011131-001D**

**Run ID: Run Order:** GCFID-HP5-B\_220122A: 13      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R373590  
**Method:** SW8015C      **Analysis Date:** 01/22/2022 23:47      **Prep Date:**  
**Lab ID:** CCV\_0122HP520r      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Diesel Range Organics (C10 to C24)	15	0.30	15		101.0	80	120				
Total Extractable Hydrocarbons	16	0.30	15		104.0	80	120				
Surr: o-Terphenyl	0.22	0.0020	0.20		109.0	80	120				

Associated Samples: **B22011131-001D**

**Run ID: Run Order:** GCFID-HP5-B\_220122A: 22      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R373590  
**Method:** SW8015C      **Analysis Date:** 01/23/2022 09:46      **Prep Date:**  
**Lab ID:** CCV\_0122HP534r-W      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
TEH(Oil Range)	4.4	0.30	5.0		89.0	80	120				
Surr: n-Triacontane	0.20	0.0020	0.20		101.0	80	120				

Associated Samples: **B22011131-001D**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** GCFID-HP5-B\_220122A: 23      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R373590  
**Method:** SW8015C      **Analysis Date:** 01/23/2022 10:29      **Prep Date:**  
**Lab ID:** CCV\_0122HP535r      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Diesel Range Organics (C10 to C24)	14	0.30	15		92.0	80	120				
Total Extractable Hydrocarbons	14	0.30	15		95.0	80	120				
Surr: o-Terphenyl	0.20	0.0020	0.20		100.0	80	120				

Associated Samples: **B22011131-001D**





### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** FID-HEADSPACE\_220120A: 4      **SampType:** Method Blank      **Batch ID:** R373491  
**Method:** SW8015M      **Analysis Date:** 01/20/2022 10:29      **Prep Date:**  
**Lab ID:** MBLK      **Units:** mg/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Methane	ND	0.0010			0.0						

Associated Samples: B22011131-001I, B22011131-005A

**Run ID: Run Order:** FID-HEADSPACE\_220120A: 2      **SampType:** Laboratory Control Sample      **Batch ID:** R373491  
**Method:** SW8015M      **Analysis Date:** 01/20/2022 09:21      **Prep Date:**  
**Lab ID:** LCS      **Units:** ppm      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Methane	99	2.0	100		99.0	85	115				

Associated Samples: B22011131-001I, B22011131-005A

**Run ID: Run Order:** FID-HEADSPACE\_220120A: 3      **SampType:** Laboratory Control Sample Duplicate      **Batch ID:** R373491  
**Method:** SW8015M      **Analysis Date:** 01/20/2022 09:25      **Prep Date:**  
**Lab ID:** LCSD      **Units:** ppm      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Methane	99	2.0	100		99.0	85	115	99	0.1	20.0	

Associated Samples: B22011131-001I, B22011131-005A



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** FID-HEADSPACE\_220120A: 10  
**Method:** SW8015M  
**Lab ID:** B22011126-001IDUP  
**SampType:** Sample Duplicate  
**Analysis Date:** 01/20/2022 11:12  
**Units:** mg/L

**Batch ID:** R373491  
**Prep Date:**  
**Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Methane	0.010	0.0020			0.0			0.010	1.1	20.0	

Associated Samples: B22011131-001I, B22011131-005A

**Run ID: Run Order:** FID-HEADSPACE\_220120A: 1  
**Method:** SW8015M  
**Lab ID:** CCV  
**SampType:** Continuing Calibration Verification Standard  
**Analysis Date:** 01/20/2022 09:16  
**Units:** ppm

**Batch ID:** R373491  
**Prep Date:**  
**Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Methane	98	2.0	100		98.0	85	115				

Associated Samples: B22011131-001I, B22011131-005A

**Run ID: Run Order:** FID-HEADSPACE\_220120A: 26  
**Method:** SW8015M  
**Lab ID:** CCV  
**SampType:** Continuing Calibration Verification Standard  
**Analysis Date:** 01/20/2022 12:56  
**Units:** ppm

**Batch ID:** R373491  
**Prep Date:**  
**Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Methane	96	2.0	100		96.0	85	115				

Associated Samples: B22011131-001I, B22011131-005A



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5973N.I\_220201A: 12  
**Method:** SW8270C  
**Lab ID:** MB-163072

**SampType:** Method Blank  
**Analysis Date:** 02/01/2022 22:46  
**Units:** ug/L

**Batch ID:** 163072  
**Prep Date:** 01/19/2022 15:57  
**Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,2,4-Trichlorobenzene	ND	5.0									
1,2-Dichlorobenzene	ND	5.0									
1,3-Dichlorobenzene	ND	5.0									
1,4-Dichlorobenzene	ND	5.0									
2,4,5-Trichlorophenol	ND	5.0									
2,4,6-Trichlorophenol	ND	5.0									
2,4-Dichlorophenol	ND	5.0									
2,4-Dimethylphenol	ND	5.0									
2,4-Dinitrophenol	ND	10									
2,4-Dinitrotoluene	ND	5.0									
2,6-Dinitrotoluene	ND	5.0									
2-Chloronaphthalene	ND	5.0									
2-Chlorophenol	ND	5.0									
2-Nitrophenol	ND	5.0									
3,3'-Dichlorobenzidine	ND	10									
4,6-Dinitro-2-methylphenol	ND	10									
4-Bromophenyl phenyl ether	ND	5.0									
4-Chloro-3-methylphenol	ND	5.0									
4-Chlorophenol	ND	5.0									
4-Chlorophenyl phenyl ether	ND	5.0									
4-Nitrophenol	ND	10									
Azobenzene	ND	5.0									
bis(-2-chloroethoxy)Methane	ND	5.0									
bis(-2-chloroethyl)Ether	ND	5.0									
bis(2-chloroisopropyl)Ether	ND	5.0									
bis(2-ethylhexyl)Phthalate	ND	5.0									
Butylbenzylphthalate	ND	5.0									



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5973N.I\_220201A: 12      **SampType:** Method Blank      **Batch ID:** 163072  
**Method:** SW8270C      **Analysis Date:** 02/01/2022 22:46      **Prep Date:** 01/19/2022 15:57  
**Lab ID:** MB-163072      **Units:** ug/L      **Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Diethyl phthalate	ND	5.0									
Dimethyl phthalate	ND	5.0									
Di-n-butyl phthalate	ND	5.0									
Di-n-octyl phthalate	ND	5.0									
Hexachlorobenzene	ND	5.0									
Hexachlorobutadiene	ND	5.0									
Hexachlorocyclopentadiene	ND	5.0									
Hexachloroethane	ND	5.0									
Isophorone	ND	5.0									
m+p-Cresols	ND	5.0									
Nitrobenzene	ND	5.0									
n-Nitrosodimethylamine	ND	5.0									
n-Nitroso-di-n-propylamine	ND	5.0									
n-Nitrosodiphenylamine	ND	10									
o-Cresol	ND	5.0									
Pentachlorophenol	ND	10									
Phenol	ND	5.0									
Pyridine	ND	5.0									
Surr: 2,4,6-Tribromophenol	164	5.0	200		82.0	43	140				
Surr: 2-Fluorobiphenyl	62	5.0	100		62.0	44	119				
Surr: 2-Fluorophenol	74	5.0	200		37.0	19	119				
Surr: Nitrobenzene-d5	63	5.0	100		63.0	44	120				
Surr: Phenol-d5	74	5.0	200		37.0	10	65				
Surr: Terphenyl-d14	93	5.0	100		93.0	50	134				

Associated Samples: **B22011131-001C**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5973N.I\_220201A: 13      **SampType:** Laboratory Control Sample      **Batch ID:** 163072  
**Method:** SW8270C      **Analysis Date:** 02/01/2022 23:18      **Prep Date:** 01/19/2022 15:57  
**Lab ID:** LCS-163072      **Units:** ug/L      **Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,2,4-Trichlorobenzene	64	10	100		64.0	29	116				
1,2-Dichlorobenzene	57	10	100		57.0	32	111				
1,3-Dichlorobenzene	60	10	100		60.0	28	110				
1,4-Dichlorobenzene	54	10	100		54.0	29	112				
2,4,5-Trichlorophenol	77	10	100		77.0	53	123				
2,4,6-Trichlorophenol	82	10	100		82.0	50	125				
2,4-Dichlorophenol	64	10	100		64.0	47	121				
2,4-Dimethylphenol	43	10	100		43.0	31	124				
2,4-Dinitrophenol	63	10	100		63.0	23	142				
2,4-Dinitrotoluene	73	10	100		73.0	57	128				
2,6-Dinitrotoluene	89	10	100		89.0	50	118				
2-Chloronaphthalene	80	10	100		80.0	40	116				
2-Chlorophenol	60	10	100		60.0	38	117				
2-Nitrophenol	72	10	100		72.0	47	123				
3,3'-Dichlorobenzidine	65	10	100		65.0	27	129				
4,6-Dinitro-2-methylphenol	67	10	100		67.0	44	137				
4-Bromophenyl phenyl ether	81	10	100		81.0	55	124				
4-Chloro-3-methylphenol	81	10	100		81.0	52	119				
4-Chlorophenol	65	10	100		65.0	41	81				
4-Chlorophenyl phenyl ether	86	10	100		86.0	53	121				
4-Nitrophenol	41	10	100		41.0	15	36				S
Azobenzene	73	10	100		73.0	61	116				
bis(-2-chloroethoxy)Methane	83	10	100		83.0	48	120				
bis(-2-chloroethyl)Ether	74	10	100		74.0	43	118				
bis(2-chloroisopropyl)Ether	60	10	100		60.0	37	130				
bis(2-ethylhexyl)Phthalate	86	10	100		86.0	55	135				
Butylbenzylphthalate	84	10	100		84.0	53	134				



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5973N.I\_220201A: 13      **SampType:** Laboratory Control Sample      **Batch ID:** 163072  
**Method:** SW8270C      **Analysis Date:** 02/01/2022 23:18      **Prep Date:** 01/19/2022 15:57  
**Lab ID:** LCS-163072      **Units:** ug/L      **Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Diethyl phthalate	93	10	100		93.0	56	125				
Dimethyl phthalate	85	10	100		85.0	45	127				
Di-n-butyl phthalate	89	10	100		89.0	59	127				
Di-n-octyl phthalate	87	10	100		87.0	51	140				
Hexachlorobenzene	68	10	100		68.0	53	125				
Hexachlorobutadiene	54	10	100		54.0	22	124				
Hexachlorocyclopentadiene	52	10	100		52.0	39	91				
Hexachloroethane	54	10	100		54.0	21	115				
Isophorone	76	10	100		76.0	42	124				
m+p-Cresols	61	10	100		61.0	29	110				
Nitrobenzene	77	10	100		77.0	45	121				
n-Nitrosodimethylamine	46	10	100		46.0	20	45				S
n-Nitroso-di-n-propylamine	79	10	100		79.0	49	119				
n-Nitrosodiphenylamine	79	10	100		79.0	51	123				
o-Cresol	68	10	100		68.0	30	117				
Pentachlorophenol	78	10	100		78.0	35	138				
Phenol	43	10	100		43.0	37	75				
Pyridine	32	10	100		32.0	16	45				
Surr: 2,4,6-Tribromophenol	161	10	200		80.0	43	140				
Surr: 2-Fluorobiphenyl	70	10	100		70.0	44	119				
Surr: 2-Fluorophenol	74	10	200		37.0	19	119				
Surr: Nitrobenzene-d5	69	10	100		69.0	44	120				
Surr: Phenol-d5	81	10	200		41.0	10	65				
Surr: Terphenyl-d14	84	10	100		84.0	50	134				

Associated Samples: **B22011131-001C**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5973N.I\_220201A: 14      **SampType:** Laboratory Control Sample Duplicate      **Batch ID:** 163072  
**Method:** SW8270C      **Analysis Date:** 02/01/2022 23:50      **Prep Date:** 01/19/2022 15:58  
**Lab ID:** LCSD-163072      **Units:** ug/L      **Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,2,4-Trichlorobenzene	72	10	100		72.0	29	116	64	12.0	20.0	
1,2-Dichlorobenzene	67	10	100		67.0	32	111	57	16.0	20.0	
1,3-Dichlorobenzene	67	10	100		67.0	28	110	60	11.0	20.0	
1,4-Dichlorobenzene	62	10	100		62.0	29	112	54	13.0	20.0	
2,4,5-Trichlorophenol	79	10	100		79.0	53	123	77	2.6	20.0	
2,4,6-Trichlorophenol	89	10	100		89.0	50	125	82	8.3	20.0	
2,4-Dichlorophenol	72	10	100		72.0	47	121	64	12.0	20.0	
2,4-Dimethylphenol	57	10	100		57.0	31	124	43	29.0	20.0	R
2,4-Dinitrophenol	56	10	100		56.0	23	142	63	12.0	20.0	
2,4-Dinitrotoluene	83	10	100		83.0	57	128	73	12.0	20.0	
2,6-Dinitrotoluene	93	10	100		93.0	50	118	89	4.4	20.0	
2-Chloronaphthalene	90	10	100		90.0	40	116	80	12.0	20.0	
2-Chlorophenol	68	10	100		68.0	38	117	60	11.0	20.0	
2-Nitrophenol	79	10	100		79.0	47	123	72	10.0	20.0	
3,3'-Dichlorobenzidine	74	10	100		74.0	27	129	65	12.0	20.0	
4,6-Dinitro-2-methylphenol	75	10	100		75.0	44	137	67	12.0	20.0	
4-Bromophenyl phenyl ether	94	10	100		94.0	55	124	81	16.0	20.0	
4-Chloro-3-methylphenol	89	10	100		89.0	52	119	81	9.9	20.0	
4-Chlorophenol	69	10	100		69.0	41	81	65	6.6	20.0	
4-Chlorophenyl phenyl ether	97	10	100		97.0	53	121	86	13.0	20.0	
4-Nitrophenol	41	10	100		41.0	15	36	41	1.0	20.0	S
Azobenzene	80	10	100		80.0	61	116	73	10.0	20.0	
bis(-2-chloroethoxy)Methane	97	10	100		97.0	48	120	83	16.0	20.0	
bis(-2-chloroethyl)Ether	87	10	100		87.0	43	118	74	16.0	20.0	
bis(2-chloroisopropyl)Ether	69	10	100		69.0	37	130	60	13.0	20.0	
bis(2-ethylhexyl)Phthalate	94	10	100		94.0	55	135	86	9.7	20.0	
Butylbenzylphthalate	98	10	100		98.0	53	134	84	15.0	20.0	



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5973N.I\_220201A: 14      **SampType:** Laboratory Control Sample Duplicate      **Batch ID:** 163072  
**Method:** SW8270C      **Analysis Date:** 02/01/2022 23:50      **Prep Date:** 01/19/2022 15:58  
**Lab ID:** LCSD-163072      **Units:** ug/L      **Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Diethyl phthalate	105	10	100		105.0	56	125	93	12.0	20.0	
Dimethyl phthalate	95	10	100		95.0	45	127	85	11.0	20.0	
Di-n-butyl phthalate	100	10	100		100.0	59	127	89	12.0	20.0	
Di-n-octyl phthalate	95	10	100		95.0	51	140	87	8.4	20.0	
Hexachlorobenzene	85	10	100		85.0	53	125	68	22.0	20.0	R
Hexachlorobutadiene	62	10	100		62.0	22	124	54	15.0	20.0	
Hexachlorocyclopentadiene	62	10	100		62.0	39	91	52	17.0	20.0	
Hexachloroethane	61	10	100		61.0	21	115	54	12.0	20.0	
Isophorone	87	10	100		87.0	42	124	76	13.0	20.0	
m+p-Cresols	67	10	100		67.0	29	110	61	10.0	20.0	
Nitrobenzene	89	10	100		89.0	45	121	77	14.0	20.0	
n-Nitrosodimethylamine	56	10	100		56.0	20	45	46	19.0	20.0	S
n-Nitroso-di-n-propylamine	97	10	100		97.0	49	119	79	20.0	20.0	R
n-Nitrosodiphenylamine	96	10	100		96.0	51	123	79	20.0	20.0	
o-Cresol	77	10	100		77.0	30	117	68	13.0	20.0	
Pentachlorophenol	99	10	100		99.0	35	138	78	23.0	20.0	R
Phenol	47	10	100		47.0	37	75	43	9.0	20.0	
Pyridine	41	10	100		41.0	16	45	32	24.0	20.0	R
Surr: 2,4,6-Tribromophenol	191	10	200		95.0	43	140	0.0	0.0		
Surr: 2-Fluorobiphenyl	75	10	100		75.0	44	119	0.0	0.0		
Surr: 2-Fluorophenol	83	10	200		42.0	19	119	0.0	0.0		
Surr: Nitrobenzene-d5	79	10	100		79.0	44	120	0.0	0.0		
Surr: Phenol-d5	89	10	200		44.0	10	65	0.0	0.0		
Surr: Terphenyl-d14	95	10	100		95.0	50	134	0.0	0.0		

Associated Samples: **B22011131-001C**





### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5973N.I\_220201B: 8  
**Method:** SW8270C  
**Lab ID:** B22011136-001CMS

**SampType:** Sample Matrix Spike  
**Analysis Date:** 02/02/2022 09:48  
**Units:** ug/L

**Batch ID:** 163072  
**Prep Date:** 01/20/2022 07:52  
**Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,2,4-Trichlorobenzene	66	10	99	0.0	66.0	29	116				
1,2-Dichlorobenzene	62	10	99	0.0	63.0	32	111				
1,3-Dichlorobenzene	61	10	99	0.0	62.0	28	110				
1,4-Dichlorobenzene	58	10	99	0.0	58.0	29	112				
2,4,5-Trichlorophenol	75	10	99	0.0	76.0	53	123				
2,4,6-Trichlorophenol	83	10	99	0.0	84.0	50	125				
2,4-Dichlorophenol	68	10	99	0.0	68.0	47	121				
2,4-Dimethylphenol	60	10	99	0.0	61.0	31	124				
2,4-Dinitrophenol	63	10	99	0.0	63.0	23	142				
2,4-Dinitrotoluene	83	10	99	0.0	83.0	57	128				
2,6-Dinitrotoluene	91	10	99	0.0	92.0	50	118				
2-Chloronaphthalene	89	10	99	0.0	90.0	40	116				
2-Chlorophenol	65	10	99	0.0	66.0	38	117				
2-Nitrophenol	74	10	99	0.0	74.0	47	123				
3,3'-Dichlorobenzidine	59	10	99	0.0	60.0	27	129				
4,6-Dinitro-2-methylphenol	66	10	99	0.0	67.0	44	137				
4-Bromophenyl phenyl ether	82	10	99	0.0	83.0	55	124				
4-Chloro-3-methylphenol	82	10	99	0.0	83.0	52	119				
4-Chlorophenol	59	10	99	0.0	60.0	41	81				
4-Chlorophenyl phenyl ether	91	10	99	0.0	92.0	53	121				
4-Nitrophenol	37	10	99	0.0	38.0	15	36				S
Azobenzene	76	10	99	0.0	76.0	61	116				
bis(-2-chloroethoxy)Methane	87	10	99	0.0	87.0	48	120				
bis(-2-chloroethyl)Ether	78	10	99	0.0	79.0	43	118				
bis(2-chloroisopropyl)Ether	79	10	99	0.0	80.0	37	130				
bis(2-ethylhexyl)Phthalate	68	10	99	0.0	68.0	55	135				
Butylbenzylphthalate	82	10	99	0.0	83.0	53	134				



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5973N.I\_220201B: 8      **SampType:** Sample Matrix Spike      **Batch ID:** 163072  
**Method:** SW8270C      **Analysis Date:** 02/02/2022 09:48      **Prep Date:** 01/20/2022 07:52  
**Lab ID:** B22011136-001CMS      **Units:** ug/L      **Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Diethyl phthalate	92	10	99	0.0	93.0	56	125				
Dimethyl phthalate	91	10	99	0.0	92.0	45	127				
Di-n-butyl phthalate	89	10	99	0.0	90.0	59	127				
Di-n-octyl phthalate	69	10	99	0.0	70.0	51	140				
Hexachlorobenzene	69	10	99	0.0	70.0	53	125				
Hexachlorobutadiene	55	10	99	0.0	55.0	22	124				
Hexachlorocyclopentadiene	54	10	99	0.0	54.0	39	91				
Hexachloroethane	56	10	99	0.0	56.0	21	115				
Isophorone	73	10	99	0.0	74.0	42	124				
m+p-Cresols	63	10	99	0.0	63.0	29	110				
Nitrobenzene	82	10	99	0.0	82.0	45	121				
n-Nitrosodimethylamine	50	10	99	0.0	50.0	20	45				S
n-Nitroso-di-n-propylamine	86	10	99	0.0	87.0	49	119				
n-Nitrosodiphenylamine	82	10	99	0.0	83.0	51	123				
o-Cresol	71	10	99	0.0	72.0	30	117				
Pentachlorophenol	71	10	99	0.0	71.0	35	138				
Phenol	43	10	99	0.0	43.0	37	75				
Pyridine	31	10	99	0.0	32.0	16	45				
Surr: 2,4,6-Tribromophenol	149	10	198	0.0	75.0	43	140				
Surr: 2-Fluorobiphenyl	76	10	99	0.0	77.0	44	119				
Surr: 2-Fluorophenol	76	10	198	0.0	38.0	19	119				
Surr: Nitrobenzene-d5	72	10	99	0.0	73.0	44	120				
Surr: Phenol-d5	84	10	198	0.0	42.0	10	65				
Surr: Terphenyl-d14	84	10	99	0.0	85.0	50	134				

Associated Samples: **B22011131-001C**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5973N.I\_220201B: 9      **SampType:** Sample Matrix Spike Duplicate      **Batch ID:** 163072  
**Method:** SW8270C      **Analysis Date:** 02/02/2022 10:20      **Prep Date:** 01/20/2022 07:52  
**Lab ID:** B22011136-001CMSD      **Units:** ug/L      **Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,2,4-Trichlorobenzene	59	10	99	0.0	59.0	29	116	66	11.0	20.0	
1,2-Dichlorobenzene	46	10	99	0.0	47.0	32	111	62	30.0	20.0	R
1,3-Dichlorobenzene	46	10	99	0.0	46.0	28	110	61	29.0	20.0	R
1,4-Dichlorobenzene	42	10	99	0.0	43.0	29	112	58	31.0	20.0	R
2,4,5-Trichlorophenol	69	10	99	0.0	69.0	53	123	75	9.4	20.0	
2,4,6-Trichlorophenol	73	10	99	0.0	74.0	50	125	83	12.0	20.0	
2,4-Dichlorophenol	55	10	99	0.0	55.0	47	121	68	21.0	20.0	R
2,4-Dimethylphenol	49	10	99	0.0	49.0	31	124	60	21.0	20.0	R
2,4-Dinitrophenol	55	10	99	0.0	55.0	23	142	63	14.0	20.0	
2,4-Dinitrotoluene	75	10	99	0.0	76.0	57	128	83	9.7	20.0	
2,6-Dinitrotoluene	84	10	99	0.0	85.0	50	118	91	8.0	20.0	
2-Chloronaphthalene	81	10	99	0.0	82.0	40	116	89	9.4	20.0	
2-Chlorophenol	52	10	99	0.0	52.0	38	117	65	23.0	20.0	R
2-Nitrophenol	63	10	99	0.0	63.0	47	123	74	16.0	20.0	
3,3'-Dichlorobenzidine	60	10	99	0.0	61.0	27	129	59	1.3	20.0	
4,6-Dinitro-2-methylphenol	64	10	99	0.0	65.0	44	137	66	3.2	20.0	
4-Bromophenyl phenyl ether	79	10	99	0.0	80.0	55	124	82	4.0	20.0	
4-Chloro-3-methylphenol	74	10	99	0.0	74.0	52	119	82	11.0	20.0	
4-Chlorophenol	52	10	99	0.0	52.0	41	81	59	14.0	20.0	
4-Chlorophenyl phenyl ether	84	10	99	0.0	85.0	53	121	91	8.4	20.0	
4-Nitrophenol	36	10	99	0.0	36.0	15	36	37	5.0	20.0	
Azobenzene	70	10	99	0.0	71.0	61	116	76	7.2	20.0	
bis(-2-chloroethoxy)Methane	75	10	99	0.0	75.0	48	120	87	15.0	20.0	
bis(-2-chloroethyl)Ether	69	10	99	0.0	70.0	43	118	78	12.0	20.0	
bis(2-chloroisopropyl)Ether	64	10	99	0.0	64.0	37	130	79	22.0	20.0	R
bis(2-ethylhexyl)Phthalate	84	10	99	0.0	85.0	55	135	68	22.0	20.0	R
Butylbenzylphthalate	89	10	99	0.0	90.0	53	134	82	8.5	20.0	



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5973N.I\_220201B: 9      **SampType:** Sample Matrix Spike Duplicate      **Batch ID:** 163072  
**Method:** SW8270C      **Analysis Date:** 02/02/2022 10:20      **Prep Date:** 01/20/2022 07:52  
**Lab ID:** B22011136-001CMSD      **Units:** ug/L      **Prep Method:** SW3510C

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Diethyl phthalate	88	10	99	0.0	88.0	56	125	92	4.9	20.0	
Dimethyl phthalate	82	10	99	0.0	82.0	45	127	91	11.0	20.0	
Di-n-butyl phthalate	91	10	99	0.0	92.0	59	127	89	1.7	20.0	
Di-n-octyl phthalate	84	10	99	0.0	85.0	51	140	69	19.0	20.0	
Hexachlorobenzene	68	10	99	0.0	68.0	53	125	69	2.1	20.0	
Hexachlorobutadiene	49	10	99	0.0	50.0	22	124	55	9.9	20.0	
Hexachlorocyclopentadiene	57	10	99	0.0	58.0	39	91	54	6.0	20.0	
Hexachloroethane	42	10	99	0.0	42.0	21	115	56	29.0	20.0	R
Isophorone	64	10	99	0.0	64.0	42	124	73	14.0	20.0	
m+p-Cresols	55	10	99	0.0	56.0	29	110	63	13.0	20.0	
Nitrobenzene	74	10	99	0.0	74.0	45	121	82	10.0	20.0	
n-Nitrosodimethylamine	42	10	99	0.0	43.0	20	45	50	16.0	20.0	
n-Nitroso-di-n-propylamine	69	10	99	0.0	69.0	49	119	86	23.0	20.0	R
n-Nitrosodiphenylamine	78	10	99	0.0	79.0	51	123	82	5.3	20.0	
o-Cresol	57	10	99	0.0	58.0	30	117	71	21.0	20.0	R
Pentachlorophenol	74	10	99	0.0	75.0	35	138	71	4.6	20.0	
Phenol	36	10	99	0.0	36.0	37	75	43	17.0	20.0	S
Pyridine	27	10	99	0.0	27.0	16	45	31	16.0	20.0	
Surr: 2,4,6-Tribromophenol	141	10	198	0.0	71.0	43	140	0.0	0.0		
Surr: 2-Fluorobiphenyl	61	10	99	0.0	62.0	44	119	0.0	0.0		
Surr: 2-Fluorophenol	55	10	198	0.0	28.0	19	119	0.0	0.0		
Surr: Nitrobenzene-d5	60	10	99	0.0	60.0	44	120	0.0	0.0		
Surr: Phenol-d5	62	10	198	0.0	31.0	10	65	0.0	0.0		
Surr: Terphenyl-d14	78	10	99	0.0	79.0	50	134	0.0	0.0		

Associated Samples: **B22011131-001C**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5973N.I\_220201A: 29

**SampType:** Continuing Calibration Verification Standard

**Batch ID:** R374090

**Method:** SW8270C

**Analysis Date:** 02/01/2022 21:09

**Prep Date:**

**Lab ID:** 01-Feb-22\_CCV\_9

**Units:** ug/L

**Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,2,4-Trichlorobenzene	78	10	75		104.0	80	120				
1,2-Dichlorobenzene	82	10	75		109.0	80	120				
1,3-Dichlorobenzene	82	10	75		110.0	80	120				
1,4-Dichlorobenzene	81	10	75		108.0	80	120				
2,4,5-Trichlorophenol	84	10	75		112.0	80	120				
2,4,6-Trichlorophenol	84	10	75		112.0	80	120				
2,4-Dichlorophenol	79	10	75		105.0	80	120				
2,4-Dimethylphenol	72	10	75		96.0	80	120				
2,4-Dinitrophenol	67	10	75		89.0	80	120				
2,4-Dinitrotoluene	84	10	75		112.0	80	120				
2,6-Dinitrotoluene	89	10	75		118.0	80	120				
2-Chloronaphthalene	86	10	75		115.0	80	120				
2-Chlorophenol	81	10	75		108.0	80	120				
2-Nitrophenol	74	10	75		98.0	80	120				
3,3'-Dichlorobenzidine	69	10	75		92.0	80	120				
4,6-Dinitro-2-methylphenol	63	10	75		84.0	80	120				
4-Bromophenyl phenyl ether	76	10	75		101.0	80	120				
4-Chloro-3-methylphenol	75	10	75		100.0	80	120				
4-Chlorophenol	75	10	75		100.0	80	120				
4-Chlorophenyl phenyl ether	82	10	75		109.0	80	120				
4-Nitrophenol	90	10	75		119.0	80	120				
Azobenzene	75	10	75		100.0	80	120				
bis(-2-chloroethoxy)Methane	79	10	75		105.0	80	120				
bis(-2-chloroethyl)Ether	84	10	75		112.0	80	120				
bis(2-chloroisopropyl)Ether	66	10	75		88.0	80	120				
bis(2-ethylhexyl)Phthalate	83	10	75		110.0	80	120				
Butylbenzylphthalate	81	10	75		109.0	80	120				



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5973N.I\_220201A: 29      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R374090  
**Method:** SW8270C      **Analysis Date:** 02/01/2022 21:09      **Prep Date:**  
**Lab ID:** 01-Feb-22\_CCV\_9      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Diethyl phthalate	88	10	75		117.0	80	120				
Dimethyl phthalate	84	10	75		112.0	80	120				
Di-n-butyl phthalate	84	10	75		112.0	80	120				
Di-n-octyl phthalate	81	10	75		108.0	80	120				
Hexachlorobenzene	74	10	75		98.0	80	120				
Hexachlorobutadiene	76	10	75		101.0	80	120				
Hexachlorocyclopentadiene	75	10	75		100.0	80	120				
Hexachloroethane	82	10	75		110.0	80	120				
Isophorone	69	10	75		92.0	80	120				
m+p-Cresols	75	10	75		100.0	80	120				
Nitrobenzene	81	10	75		108.0	80	120				
n-Nitrosodimethylamine	85	10	75		113.0	80	120				
n-Nitroso-di-n-propylamine	86	10	75		114.0	80	120				
n-Nitrosodiphenylamine	75	10	75		100.0	80	120				
o-Cresol	84	10	75		112.0	80	120				
Pentachlorophenol	79	10	75		106.0	80	120				
Phenol	75	10	75		101.0	80	120				
Pyridine	84	10	75		112.0	80	120				
Surr: 2,4,6-Tribromophenol	80	10	75		107.0	80	120				
Surr: 2-Fluorobiphenyl	72	10	75		96.0	80	120				
Surr: 2-Fluorophenol	84	10	75		112.0	80	120				
Surr: Nitrobenzene-d5	75	10	75		101.0	80	120				
Surr: Phenol-d5	84	10	75		112.0	80	120				
Surr: Terphenyl-d14	76	10	75		102.0	80	120				

Associated Samples: **B22011131-001C**



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5973N.I\_220201A: 25      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R374090  
**Method:** SW8270C      **Analysis Date:** 02/02/2022 05:42      **Prep Date:**  
**Lab ID:** 01-Feb-22\_CCV\_25      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
1,2,4-Trichlorobenzene	69	10	75		92.0	50	150				
1,2-Dichlorobenzene	74	10	75		98.0	50	150				
1,3-Dichlorobenzene	75	10	75		100.0	50	150				
1,4-Dichlorobenzene	74	10	75		99.0	50	150				
2,4,5-Trichlorophenol	81	10	75		108.0	50	150				
2,4,6-Trichlorophenol	82	10	75		109.0	50	150				
2,4-Dichlorophenol	76	10	75		101.0	50	150				
2,4-Dimethylphenol	69	10	75		92.0	50	150				
2,4-Dinitrophenol	63	10	75		84.0	50	150				
2,4-Dinitrotoluene	71	10	75		95.0	50	150				
2,6-Dinitrotoluene	81	10	75		108.0	50	150				
2-Chloronaphthalene	81	10	75		108.0	50	150				
2-Chlorophenol	77	10	75		102.0	50	150				
2-Nitrophenol	71	10	75		94.0	50	150				
3,3'-Dichlorobenzidine	77	10	75		103.0	50	150				
4,6-Dinitro-2-methylphenol	61	10	75		82.0	50	150				
4-Bromophenyl phenyl ether	74	10	75		98.0	50	150				
4-Chloro-3-methylphenol	77	10	75		103.0	50	150				
4-Chlorophenol	72	10	75		96.0	50	150				
4-Chlorophenyl phenyl ether	80	10	75		107.0	50	150				
4-Nitrophenol	81	10	75		108.0	50	150				
Azobenzene	74	10	75		98.0	50	150				
bis(-2-chloroethoxy)Methane	79	10	75		105.0	50	150				
bis(-2-chloroethyl)Ether	78	10	75		103.0	50	150				
bis(2-chloroisopropyl)Ether	72	10	75		96.0	50	150				
bis(2-ethylhexyl)Phthalate	74	10	75		98.0	50	150				
Butylbenzylphthalate	75	10	75		100.0	50	150				



### Analytical QC Summary Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

**Run ID: Run Order:** SV5973N.I\_220201A: 25      **SampType:** Continuing Calibration Verification Standard      **Batch ID:** R374090  
**Method:** SW8270C      **Analysis Date:** 02/02/2022 05:42      **Prep Date:**  
**Lab ID:** 01-Feb-22\_CCV\_25      **Units:** ug/L      **Prep Method:**

Analytes	Result	LOQ	Spk value	Spk RefVal	%REC	LowLimit	HighLimit	RPD RefVal	%RPD	RPDLimit	Qual
Diethyl phthalate	81	10	75		108.0	50	150				
Dimethyl phthalate	77	10	75		103.0	50	150				
Di-n-butyl phthalate	77	10	75		102.0	50	150				
Di-n-octyl phthalate	76	10	75		101.0	50	150				
Hexachlorobenzene	72	10	75		96.0	50	150				
Hexachlorobutadiene	70	10	75		94.0	50	150				
Hexachlorocyclopentadiene	71	10	75		95.0	50	150				
Hexachloroethane	77	10	75		103.0	50	150				
Isophorone	72	10	75		96.0	50	150				
m+p-Cresols	77	10	75		102.0	50	150				
Nitrobenzene	70	10	75		94.0	50	150				
n-Nitrosodimethylamine	78	10	75		104.0	50	150				
n-Nitroso-di-n-propylamine	78	10	75		104.0	50	150				
n-Nitrosodiphenylamine	68	10	75		90.0	50	150				
o-Cresol	73	10	75		98.0	50	150				
Pentachlorophenol	71	10	75		94.0	50	150				
Phenol	78	10	75		104.0	50	150				
Pyridine	71	10	75		95.0	50	150				
Surr: 2,4,6-Tribromophenol	73	10	75		98.0	50	150				
Surr: 2-Fluorobiphenyl	74	10	75		98.0	50	150				
Surr: 2-Fluorophenol	77	10	75		103.0	50	150				
Surr: Nitrobenzene-d5	72	10	75		96.0	50	150				
Surr: Phenol-d5	80	10	75		107.0	50	150				
Surr: Terphenyl-d14	73	10	75		97.0	50	150				

Associated Samples: **B22011131-001C**



### Analytical QC Exceptions Report

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu  
**Workorder:** B22011131  
**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

Analysis Method	Analysis	Batch ID	Associated Samples	Sample Type	Lab ID	Analysis Date	Analysis Time	Analyte	%REC	Low Limit	High Limit	% RPD	RPD Limit	Qual		
SW6020	Metals by ICP-MS, Total	163063	001B	SD	B22011134-001BDIL	1/22/2022	00:21	Lead					10.0	N		
						1/22/2022	11:36	Lead					10.0	N		
SW8270C	Semi-Volatile Organic Compounds, Extended List	163072	001C	LCS-DOD	LCS-163072	2/1/2022	23:18	4-Nitrophenol	41.0	15	36			S		
										n-Nitrosodimethylamine	46.0	20	45			S
				LCSD-DOD	LCSD-163072	2/1/2022	23:50	2,4-Dimethylphenol	57.0	31	124	29	20.0	R		
								4-Nitrophenol	41.0	15	36	1.0	20.0	S		
								Hexachlorobenzene	85.0	53	125	22	20.0	R		
								n-Nitrosodimethylamine	56.0	20	45	19	20.0	S		
								Pentachlorophenol	99.0	35	138	23	20.0	R		
				MS-DOD	B22011136-001CMS	2/2/2022	09:48	4-Nitrophenol	38.0	15	36			S		
								n-Nitrosodimethylamine	50.0	20	45			S		
				MSD-DOD	B22011136-001CMSD	2/2/2022	10:20	1,2-Dichlorobenzene	47.0	32	111	30	20.0	R		
								1,3-Dichlorobenzene	46.0	28	110	29	20.0	R		
								1,4-Dichlorobenzene	43.0	29	112	31	20.0	R		
								2,4-Dichlorophenol	55.0	47	121	21	20.0	R		
								2,4-Dimethylphenol	49.0	31	124	21	20.0	R		
								2-Chlorophenol	52.0	38	117	23	20.0	R		
bis(2-chloroisopropyl)Ether	64.0	37	130					22	20.0	R						
bis(2-ethylhexyl)Phthalate	85.0	55	135					22	20.0	R						
Hexachloroethane	42.0	21	115					29	20.0	R						
n-Nitroso-di-n-propylamine	69.0	49	119	23	20.0	R										
o-Cresol	58.0	30	117	21	20.0	R										
Phenol	36.0	37	75	17	20.0	S										
SW8270C SIM	Low Level PAH by 8270C SIM	163072	001C	MS	B22011136-001CLMS	1/27/2022	11:31	Benzo(a)anthracene	112.0	41	105			S		



## Preparation and Analysis Dates Report

Work Order: B22011131

Client: AECOM - Honolulu

Project Name: CV18F0126/60571032.02.46.01

Report Date: 3/01/2022

Lab ID	Client Sample ID	Collection Date	Matrix	Test Name	TCLP Date	Prep Method	Prep Date	Prep Batch	Analysis Method	Analysis Date
001B	ERH2460 (RHMW12A)	01/17/2022 17:15	Ground Water	Metals by ICP-MS, Total		SW3010A	01/19/2022 14:41	163063	SW6020	01/21/2022 23:25
001C	ERH2460 (RHMW12A)	01/17/2022 17:15	Ground Water	Low Level PAH by 8270C SIM		SW3510C	01/19/2022 15:58	163072	SW8270CSIM	01/27/2022 07:45
				Semi-Volatile Organic Compounds, Extended List		SW3510C	01/19/2022 15:58	163072	SW8270C	02/02/2022 04:06
001D	ERH2460 (RHMW12A)	01/17/2022 17:15	Ground Water	Diesel Range Organics		SW3520C	01/19/2022 16:30	163074	SW8015C	01/23/2022 01:56
001H	ERH2460 (RHMW12A)	01/17/2022 17:15	Ground Water	EDB in Water by ECD		SW8011	01/21/2022 07:48	163129	SW8011	01/22/2022 10:35
004A	ERH2459 (Trip Blank)-14705	01/17/2022 17:15	Trip Blank	EDB in Water by ECD		SW8011	01/21/2022 07:49	163129	SW8011	01/22/2022 08:15



## Chemical Abstracts Service (CAS) Registry Numbers

Prepared by Billings, MT Branch

**Client:** AECOM - Honolulu

**Workorder:** B22011131

**Project:** CV18F0126/60571032.02.46.01

**Report Date:** 03/01/2022

Analyses	CAS No
<b>LOW LEVEL PAH BY 8270C SIM</b>	
1-Methylnaphthalene	90-12-0
2-Methylnaphthalene	91-57-6
Acenaphthene	83-32-9
Acenaphthylene	208-96-8
Anthracene	120-12-7
Benzo(a)anthracene	56-55-3
Benzo(a)pyrene	50-32-8
Benzo(b)fluoranthene	205-99-2
Benzo(g,h,i)perylene	191-24-2
Benzo(k)fluoranthene	207-08-9
Chrysene	218-01-9
Dibenzo(a,h)anthracene	53-70-3
Fluoranthene	206-44-0
Fluorene	86-73-7
Indeno(1,2,3-cd)pyrene	193-39-5
Naphthalene	91-20-3
Phenanthrene	85-01-8
Pyrene	129-00-0
<b>AGGREGATE ORGANICS</b>	
Organic Carbon, Total (TOC)	7440-44-0
<b>METALS, TOTAL</b>	
Lead	7439-92-1
<b>METALS, DISSOLVED</b>	
Lead	7439-92-1
<b>VOLATILE ORGANIC COMPOUNDS</b>	
Benzene	71-43-2
Bromobenzene	108-86-1
Bromochloromethane	74-97-5
Bromodichloromethane	75-27-4
Bromoform	75-25-2
Carbon tetrachloride	56-23-5
Chlorobenzene	108-90-7
Chlorodibromomethane	124-48-1
Chloroethane	75-00-3
Chloroform	67-66-3
Chloromethane	74-87-3
1,2-Dibromoethane	106-93-4

2-Chlorotoluene	95-49-8
4-Chlorotoluene	106-43-4
Dibromomethane	74-95-3
1,2-Dichlorobenzene	95-50-1
1,3-Dichlorobenzene	541-73-1
1,4-Dichlorobenzene	106-46-7
Dichlorodifluoromethane	75-71-8
1,1-Dichloroethane	75-34-3
1,2-Dichloroethane	107-06-2
1,1-Dichloroethene	75-35-4
cis-1,2-Dichloroethene	156-59-2
trans-1,2-Dichloroethene	156-60-5
1,2-Dichloropropane	78-87-5
1,3-Dichloropropane	142-28-9
2,2-Dichloropropane	594-20-7
1,1-Dichloropropene	563-58-6
cis-1,3-Dichloropropene	10061-01-5
trans-1,3-Dichloropropene	10061-02-6
Ethylbenzene	100-41-4
Methyl ethyl ketone	78-93-3
Methyl tert-butyl ether (MTBE)	1634-04-4
Methylene chloride	75-09-2
Styrene	100-42-5
1,1,1,2-Tetrachloroethane	630-20-6
1,1,2,2-Tetrachloroethane	79-34-5
Tetrachloroethene	127-18-4
Toluene	108-88-3
1,1,1-Trichloroethane	71-55-6
1,1,2-Trichloroethane	79-00-5
Trichloroethene	79-01-6
Trichlorofluoromethane	75-69-4
1,2,3-Trichloropropane	96-18-4
Vinyl chloride	75-01-4
m+p-Xylenes	179601-23-1
o-Xylene	95-47-6
Xylenes, Total	1330-20-7

#### **VOCS BY MICROEXTRACTION-ECD**

1,2-Dibromoethane	106-93-4
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#### **PETROLEUM HYDROCARBONS-VOLATILE**

C6 to C10  
Total Purgeable Hydrocarbons

#### **PETROLEUM HYDROCARBONS-SEMI-VOLATILE**

Diesel Range Organics (C10 to C24)  
Oil Range Hydrocarbons (C24 to C40)  
Total Extractable Hydrocarbons

#### **ORGANIC CHARACTERISTICS**

Methane	74-82-8
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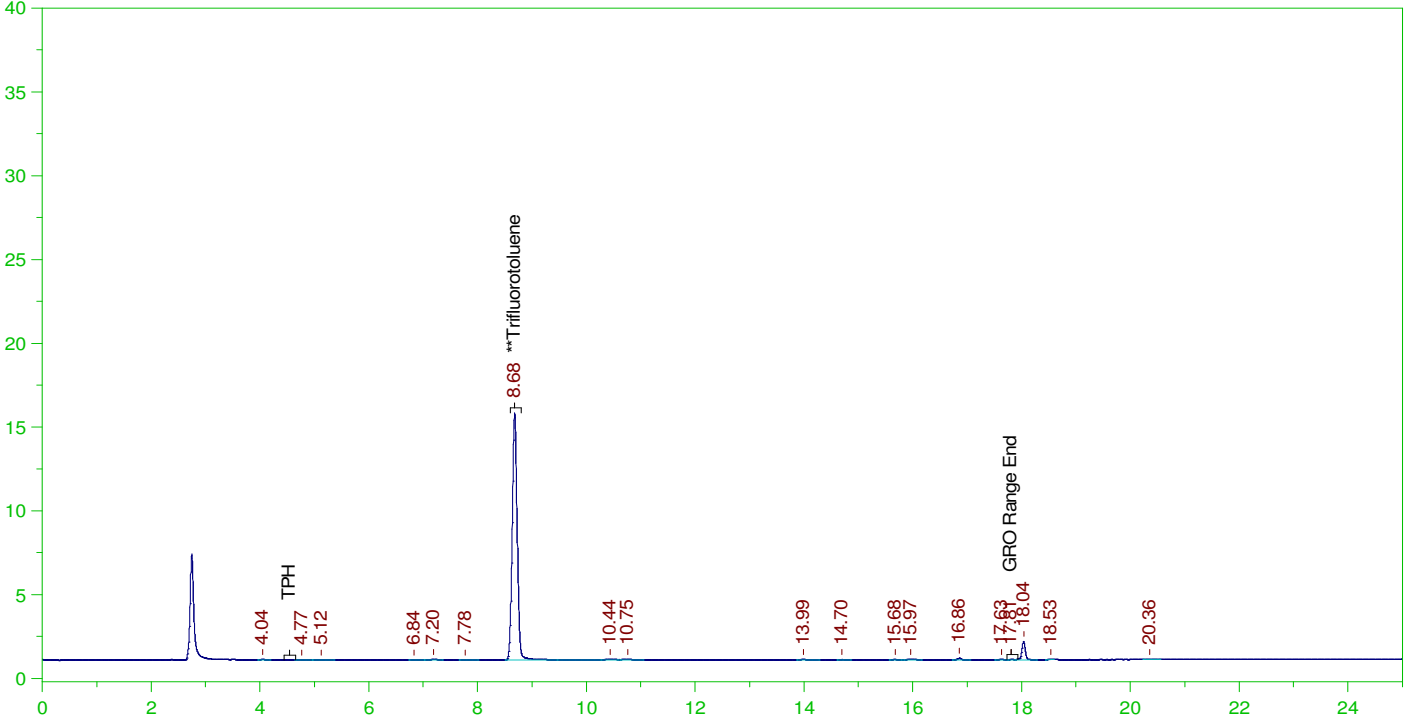
**SEMI-VOLATILE ORGANIC COMPOUNDS**

1,2,4-Trichlorobenzene	120-82-1
1,2-Dichlorobenzene	95-50-1
1,3-Dichlorobenzene	541-73-1
1,4-Dichlorobenzene	106-46-7
2,4,5-Trichlorophenol	95-95-4
2,4,6-Trichlorophenol	88-06-2
2,4-Dichlorophenol	120-83-2
2,4-Dimethylphenol	105-67-9
2,4-Dinitrophenol	51-28-5
2,4-Dinitrotoluene	121-14-2
2,6-Dinitrotoluene	606-20-2
2-Chloronaphthalene	91-58-7
2-Chlorophenol	95-57-8
2-Nitrophenol	88-75-5
3,3'-Dichlorobenzidine	91-94-1
4,6-Dinitro-2-methylphenol	534-52-1
4-Bromophenyl phenyl ether	101-55-3
4-Chloro-3-methylphenol	59-50-7
4-Chlorophenol	106-48-9
4-Chlorophenyl phenyl ether	7005-72-3
4-Nitrophenol	100-02-7
Azobenzene	103-33-3
bis(-2-chloroethoxy)Methane	111-91-1
bis(-2-chloroethyl)Ether	111-44-4
bis(2-chloroisopropyl)Ether	108-60-1
bis(2-ethylhexyl)Phthalate	117-81-7
Butylbenzylphthalate	85-68-7
Di-n-butyl phthalate	84-74-2
Di-n-octyl phthalate	117-84-0
Diethyl phthalate	84-66-2
Dimethyl phthalate	131-11-3
Hexachlorobenzene	118-74-1
Hexachlorobutadiene	87-68-3
Hexachlorocyclopentadiene	77-47-4
Hexachloroethane	67-72-1
Isophorone	78-59-1
m+p-Cresols	15831-10-4
n-Nitroso-di-n-propylamine	621-64-7
n-Nitrosodimethylamine	62-75-9
n-Nitrosodiphenylamine	86-30-6
Nitrobenzene	98-95-3
o-Cresol	95-48-7
Pentachlorophenol	87-86-5
Phenol	108-95-2
Pyridine	110-86-1

ERH2460 (RHMW12A)

G:\Org\PE1\DAT\PE1012022\_b\0120PE1B.0046.RAW

B22011131-001G ;0120PE1 , \$HC-8015-GRO-W,



**GASOLINE RANGE ORGANICS CHROMATOGRAM REPORT**

Sample Name: B22011131-001G ;0120PE1 , \$HC-8015-GRO-W,  
Raw File: G:\Org\PE1\DAT\PE1012022\_b\0120PE1B.0046.RAW  
Date & Time Acquired: 1/21/2022 10:16:57 AM  
Method File: G:\Org\PE1\Methods\211208G1131-1DoDB%.MET  
Calibration File: G:\Org\PE1\Cals\211208GRO8015CB.CAL  
Sample Weight: 5 Dilution: 1 S.A.: 1

Mean RF for C6 to C10: 945.9678  
Mean RF for TPH: 909.3915  
Rt range for Gasoline Range Organics: 4.45 to 17.93

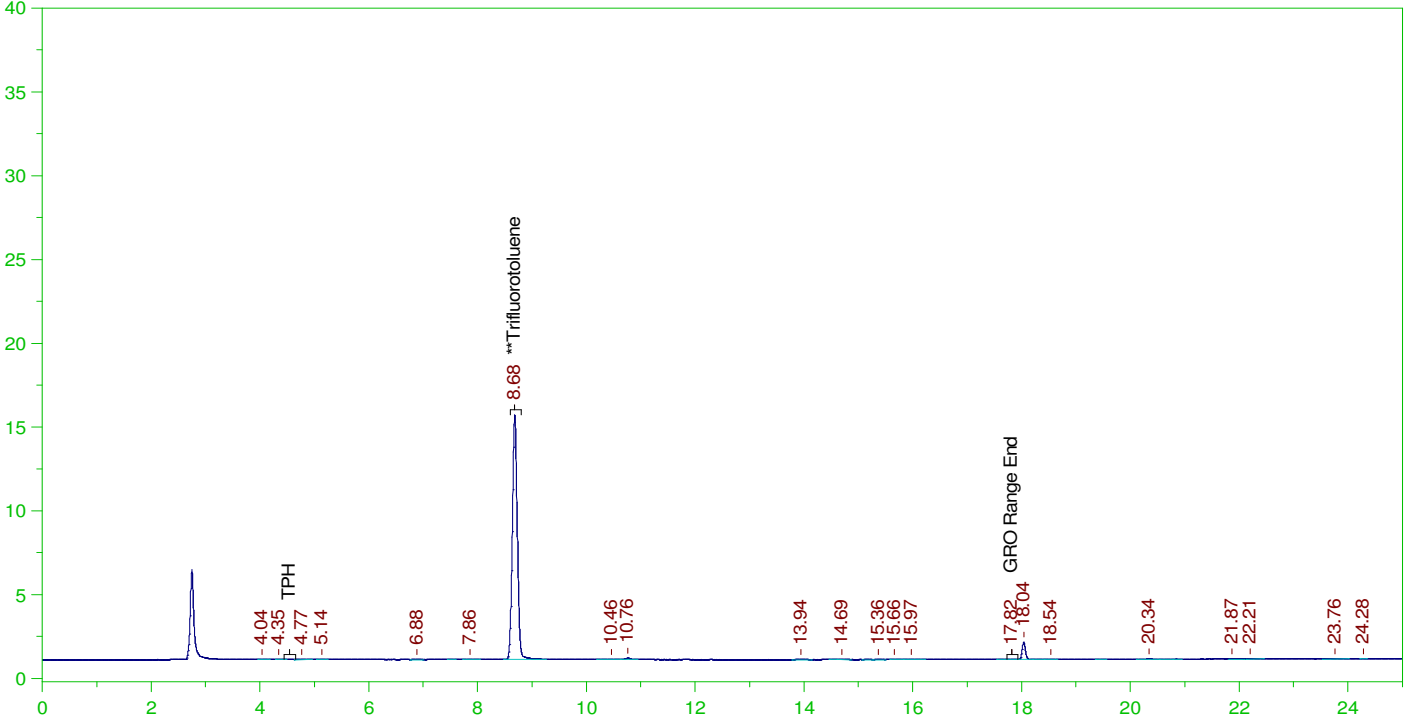
SURROGATE COMPOUND	RT	ACTUAL	MEASURED	%REC
**Trifluorotoluene	8.683	25.	20.122	80.49

C6 to C10 Area:3155.497 C6 to C10 Amount: 0.6671468  
TPH Area:8431.092 TPH Amount: 1.854227

ERH2459 (Trip Blank)-14733

G:\Org\PE1\DAT\PE1012022\_b\0120PE1B.0016.RAW

B22011131-003A ;0120PE1 , \$HC-8015-GRO-W,



**GASOLINE RANGE ORGANICS CHROMATOGRAM REPORT**

Sample Name: B22011131-003A ;0120PE1 , \$HC-8015-GRO-W,  
Raw File: G:\Org\PE1\DAT\PE1012022\_b\0120PE1B.0016.RAW  
Date & Time Acquired: 1/20/2022 5:07:12 PM  
Method File: G:\Org\PE1\Methods\211208G1131-3DoDB%.MET  
Calibration File: G:\Org\PE1\Cals\211208GRO8015CB.CAL  
Sample Weight: 5 Dilution: 1 S.A.: 1

Mean RF for C6 to C10: 945.9678  
Mean RF for TPH: 909.3915  
Rt range for Gasoline Range Organics: 4.45 to 17.93

SURROGATE COMPOUND	RT	ACTUAL	MEASURED	%REC
**Trifluorotoluene	8.685	25.	19.776	79.11

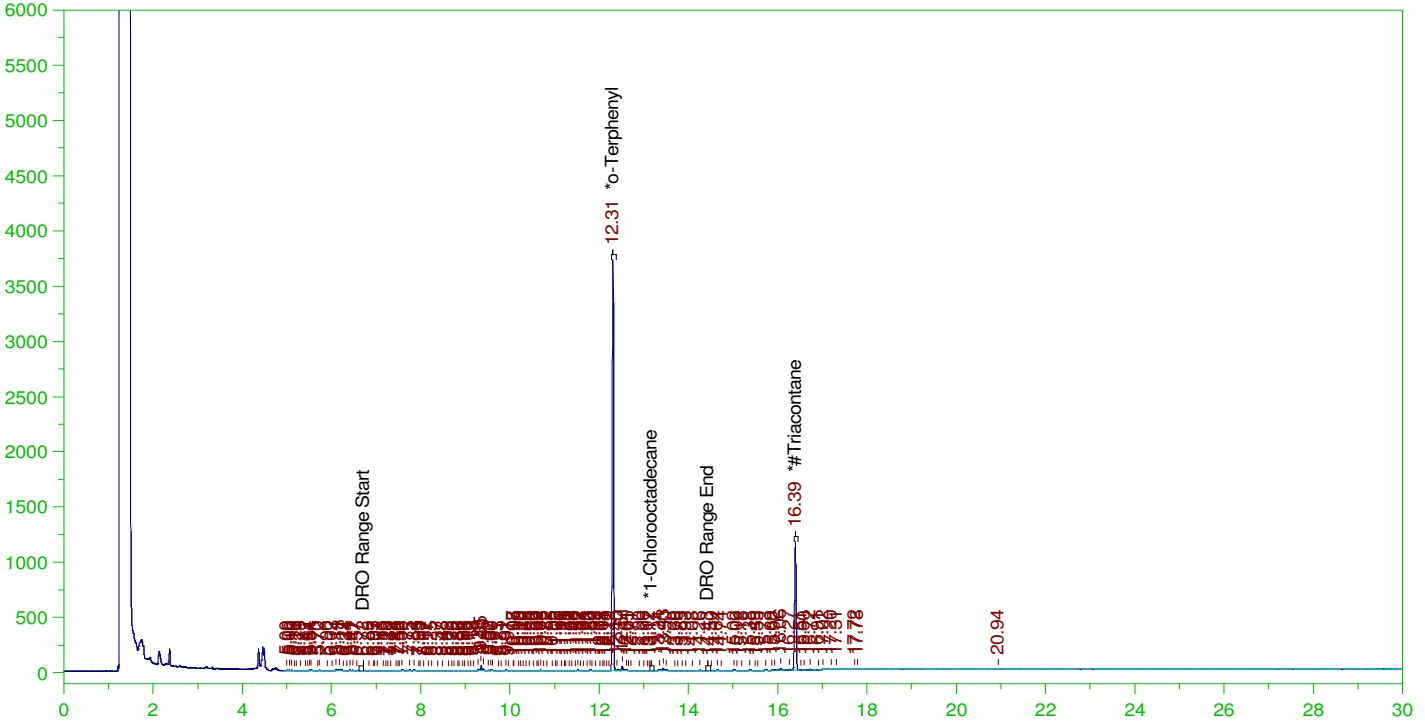
C6 to C10 Area:2526.943 C6 to C10 Amount: 0.5342556  
TPH Area:7899.134 TPH Amount: 1.737235

ERH2460 (RHMW12A)

Batch ID: 163074

G:\org\HP5\DAT\HP5012222\_b\0122HP5.0023.RAW

B22011131-001D ;0122HP5 , \$HC-8015-DRO-W,



**DIESEL RANGE ORGANICS CHROMATOGRAM REPORT**

Sample Name: B22011131-001D ;0122HP5 , \$HC-8015-DRO-W,  
Raw File: G:\org\HP5\DAT\HP5012222\_b\0122HP5.0023.RAW  
Date & Time Acquired: 1/23/2022 1:56:33 AM  
Method File: G:\Org\HP5\Methods\DR\_8015-C24T-JB-L%.met  
Calibration File: G:\Org\HP5\Cals\SW8015C\_DRO220111JB-C24-T.CAL  
Sample Weight: 1050 Dilution: 1 S.A.: 1

Mean RF for TEH: 32675.36

Rt range for Diesel Range Organics: 6.62 to 14.49

SURROGATE COMPOUND	RT	ACTUAL	MEASURED	%REC	
*o-Terphenyl	12.305	.19	.183	95.87	-
*1-Chlorooctadecane	13.172	.19	.	.03	-
*#Triacontane	16.392	.19	.096	50.6	-

DRO Area:838382.9 DRO Amount: 2.443615E-02  
TEH Area:1264552 TEH Amount: 3.685759E-02

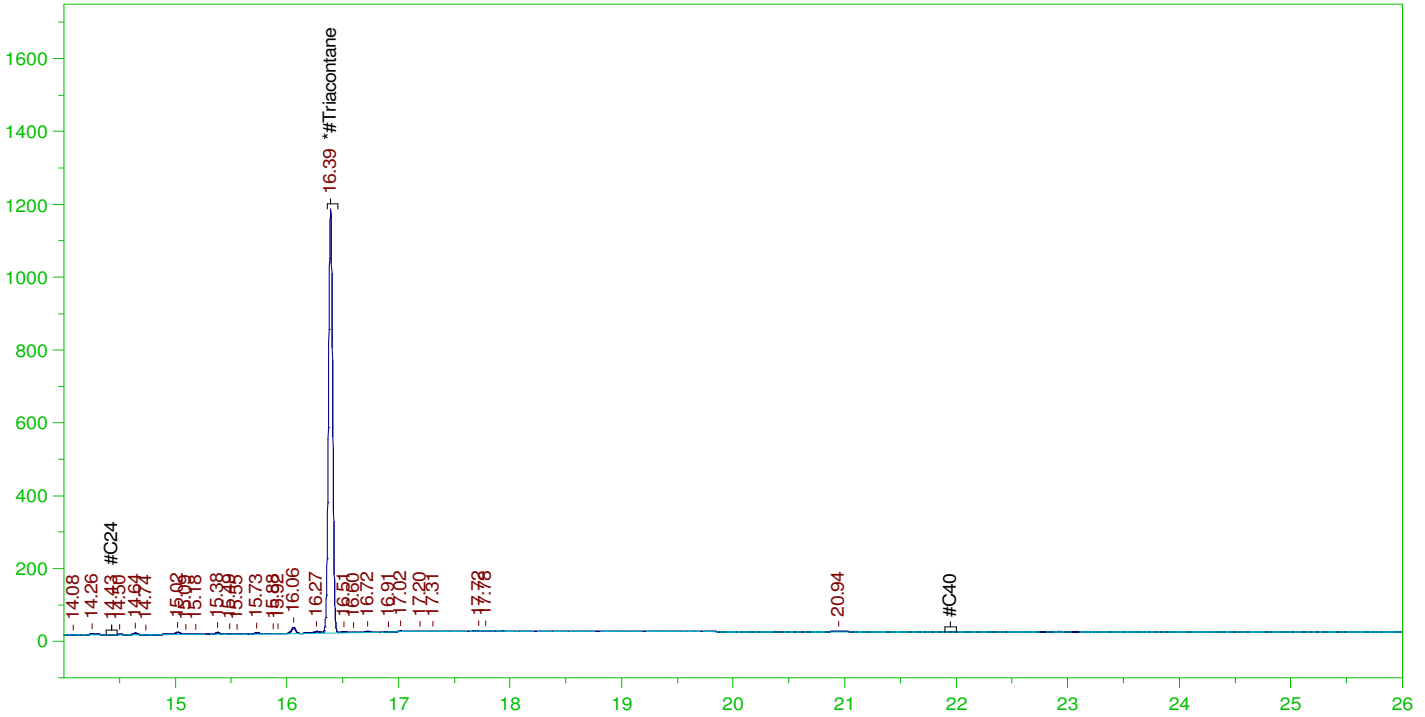


ERH2460 (RHMW12A)

Batch ID: 163074

G:\org\HP5\DAT\HP5012222\_b\0122HP5.0023.RAW

B22011131-001D ;0122HP5 , \$HC-8015-DRO-W,



**RESIDUAL RANGE ORGANICS CHROMATOGRAM**

Sample Name: B22011131-001D ;0122HP5 , \$HC-8015-DRO-W,  
Raw File: G:\org\HP5\DAT\HP5012222\_b\0122HP5.0023.RAW  
Date & Time Acquired: 1/23/2022 1:56:33 AM  
Method File: G:\Org\HP5\Methods\DR\_OROS-BBb-L%.MET  
Calibration File: G:\Org\HP5\Cals\SW8015C\_ORO220111BBb\_SAMP.CAL  
Sample Weight: 1050 Dilution: 1 S.A.: 1

Mean RF for for Residual Range Organics Calculations: 26424.55  
Rt range for Residual Range Organics: 14.38 to 22

SURROGATE COMPOUND	RT	ACTUAL	MEASURED	%REC
*#Triacontane_____	16.392	.476	.096	20.24

RRO Area:177800.6 RRO AMOUNT: 6.408204E-03

---

**From:** Ramos, Alethea <alethea.ramos@aecom.com>  
**Sent:** Monday, December 13, 2021 3:11 PM  
**To:** Tabitha Edwards  
**Cc:** Pascua, Margie; billingsPM@energylab.com  
**Subject:** RE: [EXTERNAL] FW: CV18F0126: Expedited NOI Groundwater Samples, Saturday 12/12 Submission

**Categories:** Must Attend

Hi Tabitha,

I believe Casper WY is DoD ELAP accredited in the TOC 9060 method. I spoke to Shari and she indicated there is a daily courier between Billings and Casper, and would be appx. a day delay. Under those stipulations, please subcontract these samples and inform on expedited TAT.

Thank you,

**Alethea Ramos, CIH**  
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M +1-808-389-5383  
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**From:** Tabitha Edwards <tedwards@energylab.com>  
**Sent:** Monday, December 13, 2021 7:05 AM  
**To:** Ramos, Alethea <alethea.ramos@aecom.com>  
**Cc:** Pascua, Margie <Margie.Pascua@aecom.com>; billingsPM@energylab.com  
**Subject:** [EXTERNAL] FW: CV18F0126: Expedited NOI Groundwater Samples, Saturday 12/12 Submission  
**Importance:** High

Alethea,

The TOC by 9060 must be subcontracted to our office in Casper, WY. I need authorization from you to subcontract these. Once that has been received we will discuss the TAT with them and let you know what is achievable.

Thank you,

**Energy Laboratories, Inc.**

Trust our People. Trust our Data.

**Tabitha Edwards** | Office Manager | Billings, MT

O: 406-869-6286 | [tedwards@energylab.com](mailto:tedwards@energylab.com) | [www.energylab.com](http://www.energylab.com)

This transmission may contain confidential information and is for the use of the intended recipient(s). If you received this in error, please contact the sender and delete this email and all copies.

***We want to help you ship successfully!** Please plan ahead and allow extra time to receive supplies from the lab and for the lab to receive your samples. All carriers are in full-swing holiday peak season operating with double the volume and limited capacity. We appreciate your business so please contact your local branch or Project Manager to discuss adjustments to your shipping schedule or to ask questions.*

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**From:** Ramos, Alethea [<mailto:alethea.ramos@aecom.com>]

**Sent:** Saturday, December 11, 2021 3:20 AM

**To:** Shari Endy; [billingsPM@energylab.com](mailto:billingsPM@energylab.com)

**Cc:** Jillian Miller; Pascua, Margie; KaaihiliChoy, Terri Ann

**Subject:** CV18F0126: Expedited NOI Groundwater Samples, Saturday 12/12 Submission

**Importance:** High

Hi Shari and Billings PM,

You will be receiving a Saturday shipment (12/12) of groundwater samples indicated in the attached COCs. We will need results by **Wednesday, December 15<sup>th</sup>**, and will pay any fees incurred for an expedited TAT. Please proceed with analysis without preservation traceability. Please see below tracking information links:

<https://www.fedex.com/fedextrack/?trknbr=287337969629&trkqual=2459558000~287337969629~FX>

<https://www.fedex.com/fedextrack/?trknbr=287343101019&trkqual=2459559000~287343101019~FX>

Thank you,

**Alethea Ramos, CIH**

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