

# ANALYTICAL REPORT

## PREPARED FOR

Attn: Mr. Erwin Kawata  
City & County of Honolulu  
630 South Beretania Street  
Public Service Bldg. Room 308  
Honolulu, Hawaii 96843

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## JOB DESCRIPTION

RED-HILL  
RUSH Weekly Red Hill

## JOB NUMBER

380-32096-1

# Eurofins Eaton Monrovia

## Job Notes

Laboratory certifies that the test results meet all TNI 2016 and ISO/IEC 17025:2017 requirements unless noted under the individual analysis.

Following the cover page are State Certification List, ISO 17025 Accredited Method List, Acknowledgement of Samples Received, Comments, Hits Report, Data Report, QC Summary, QC Report and Regulatory Forms, as applicable.

Test results relate only to the sample(s) tested.

Test results apply to the sample(s) as received, unless otherwise noted in the comments report (ISO/IEC 17025:2017).

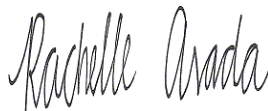
This report shall not be reproduced except in full, without the written approval of the laboratory.

This report includes ISO/IEC 17025 and non-ISO 17025 accredited methods.

## Compliance Statement

1. Laboratory is accredited in accordance with TNI 2016 Standards and ISO/IEC 17025:2017.
2. Laboratory certifies that the test results meet all TNI 2016 and ISO/IEC 17025:2017 requirements unless noted under the individual analysis
3. Test results relate only to the sample(s) tested.
4. This report shall not be reproduced except in full, without the written approval of the laboratory.
5. Unless otherwise noted, all analytes for this laboratory were covered under each accreditation/certification below. (DW, Water matrices)

## Authorization



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Authorized for release by  
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(626)386-1106



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# Definitions/Glossary

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Qualifiers

### GC/MS Semi VOA

| Qualifier | Qualifier Description  |
|-----------|--|
| ^3+       | Reporting Limit Check Standard is outside acceptance limits, high biased                                       |
| F1        | MS and/or MSD recovery exceeds control limits.   |
| J         | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value. |

### GC/MS Semi VOA TICs

| Qualifier | Qualifier Description   |
|-----------|---|
| J         | Indicates an Estimated Value for TICs                                     |
| T         | Result is a tentatively identified compound (TIC) and an estimated value. |

### Subcontract

| Qualifier | Qualifier Description          |
|-----------|--------------------------------|
| U         | This analyte was not detected. |

## Glossary

| Abbreviation   | These commonly used abbreviations may or may not be present in this report.                                 |
|----------------|---|
| ¤              | Listed under the "D" column to designate that the result is reported on a dry weight basis                  |
| %R             | Percent Recovery  |
| CFL            | Contains Free Liquid  |
| CFU            | Colony Forming Unit   |
| CNF            | Contains No Free Liquid   |
| DER            | Duplicate Error Ratio (normalized absolute difference)  |
| Dil Fac        | Dilution Factor   |
| DL             | Detection Limit (DoD/DOE)   |
| DL, RA, RE, IN | Indicates a Dilution, Re-analysis, Re-extraction, or additional Initial metals/anion analysis of the sample |
| DLC            | Decision Level Concentration (Radiochemistry)   |
| EDL            | Estimated Detection Limit (Dioxin)  |
| LOD            | Limit of Detection (DoD/DOE)  |
| LOQ            | Limit of Quantitation (DoD/DOE)   |
| MCL            | EPA recommended "Maximum Contaminant Level"   |
| MDA            | Minimum Detectable Activity (Radiochemistry)  |
| MDC            | Minimum Detectable Concentration (Radiochemistry)   |
| MDL            | Method Detection Limit  |
| ML             | Minimum Level (Dioxin)  |
| MPN            | Most Probable Number  |
| MQL            | Method Quantitation Limit   |
| NC             | Not Calculated  |
| ND             | Not Detected at the reporting limit (or MDL or EDL if shown)  |
| NEG            | Negative / Absent   |
| POS            | Positive / Present  |
| PQL            | Practical Quantitation Limit  |
| PRES           | Presumptive   |
| QC             | Quality Control   |
| RER            | Relative Error Ratio (Radiochemistry)   |
| RL             | Reporting Limit or Requested Limit (Radiochemistry)   |
| RPD            | Relative Percent Difference, a measure of the relative difference between two points                        |
| TEF            | Toxicity Equivalent Factor (Dioxin)   |
| TEQ            | Toxicity Equivalent Quotient (Dioxin)   |
| TNTC           | Too Numerous To Count   |

# Case Narrative

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

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## Job ID: 380-32096-1

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### Laboratory: Eurofins Eaton Monrovia

#### Narrative

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#### Job Narrative 380-32096-1

#### Comments

No additional comments.

#### Receipt

The samples were received on 12/21/2022 10:00 AM. Unless otherwise noted below, the samples arrived in good condition, and where required, properly preserved and on ice. The temperatures of the 2 coolers at receipt time were 1.6° C and 1.8° C.

#### Receipt Exceptions

The container label for the following samples did not match the information listed on the Chain-of-Custody (COC): Aiea Gulch Wells Pump 2 (380-32096-1), Aiea Wells Pump 2 (380-32096-2), Halawa Wells Pump 1 (380-32096-3), TB: Aiea Gulch Wells Pump 2 (380-32096-4), TB: Aiea Wells Pump 2 (380-32096-5) and TB: Halawa Wells Pump 1 (380-32096-6). Sample Aiea Wel Pump 2's respective containers' labels indicate a sample ID of AIEA WELLS PUMPS 1&2 (260), and sample Halawa Wells Pum 1's respective containers' labels indicate a sample ID of HALAWA WELLS UNITS 1&2.

#### GC/MS Semi VOA

No analytical or quality issues were noted, other than those described in the Definitions/Glossary page.

#### Subcontract non-Sister

See attached subcontract report.

#### Organic Prep

No analytical or quality issues were noted, other than those described in the Definitions/Glossary page.

#### Subcontract Work

Methods 8015 Gas (Purgeable) LL (EAL), 8015 LL DRO/MRO/JP5/JP8: These methods were subcontracted to EMAX Laboratories Inc. The subcontract laboratory certifications are different from that of the facility issuing the final report.

Method 625 PAH Physis LL (EAL) + TICs: This method was subcontracted to Physis Environmental Laboratories. The subcontract laboratory certification is different from that of the facility issuing the final report.

# Detection Summary

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

**Client Sample ID: Aiea Gulch Wells Pump 2**  
**PWSID Number: HI0000331**

**Lab Sample ID: 380-32096-1**

No Detections.

**Client Sample ID: Aiea Wells Pump 2**  
**PWSID Number: HI0000331**

**Lab Sample ID: 380-32096-2**

No Detections.

**Client Sample ID: Halawa Wells Pump 1**  
**PWSID Number: HI0000331**

**Lab Sample ID: 380-32096-3**

No Detections.

**Client Sample ID: TB: Aiea Gulch Wells Pump 2**

**Lab Sample ID: 380-32096-4**

No Detections.

**Client Sample ID: TB: Aiea Wells Pump 2**

**Lab Sample ID: 380-32096-5**

No Detections.

**Client Sample ID: TB: Halawa Wells Pump 1**

**Lab Sample ID: 380-32096-6**

No Detections.

This Detection Summary does not include radiochemical test results.

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# Client Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

**Client Sample ID: Aiea Gulch Wells Pump 2**

**Lab Sample ID: 380-32096-1**

Date Collected: 12/19/22 10:00

Matrix: Drinking Water

Date Received: 12/21/22 10:00

PWSID Number: HI0000331

**Method: EPA 525.2 - Semivolatile Organic Compounds (GC/MS)**

| Analyte                          | Result | Qualifier | RL    | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------------------------------|--------|-----------|-------|------|---|----------------|----------------|---------|
| 2,4'-DDD                         | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| 2,4'-DDE                         | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| 2,4'-DDT                         | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| 2,4-Dinitrotoluene               | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| 2,6-Dinitrotoluene               | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| 4,4'-DDD                         | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| 4,4'-DDE                         | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| 4,4'-DDT                         | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Acenaphthene                     | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Acenaphthylene                   | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Acetochlor                       | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Alachlor                         | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| alpha-BHC                        | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| alpha-Chlordane                  | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Anthracene                       | ND     |           | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Atrazine                         | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Benz(a)anthracene                | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Benzo[a]pyrene                   | ND     |           | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Benzo[b]fluoranthene             | ND     |           | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Benzo[g,h,i]perylene             | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Benzo[k]fluoranthene             | ND     |           | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| beta-BHC                         | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Bromacil                         | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Butachlor                        | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Butylbenzylphthalate             | ND     |           | 0.50  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Caffeine                         | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Chlorobenzilate                  | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Chloroneb                        | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Chlorothalonil (Draconil, Bravo) | ND     | ^3+       | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Chlorpyrifos                     | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Chrysene                         | ND     |           | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| delta-BHC                        | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Di(2-ethylhexyl)adipate          | ND     |           | 0.60  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Bis(2-ethylhexyl) phthalate      | ND     |           | 0.60  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Diazinon (Qualitative)           | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Dibenz(a,h)anthracene            | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Diclorvos (DDVP)                 | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Dieldrin                         | ND     |           | 0.20  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Diethylphthalate                 | ND     |           | 0.50  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Dimethoate                       | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Dimethylphthalate                | ND     |           | 0.50  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Di-n-butyl phthalate             | ND     |           | 1.0   | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Di-n-octyl phthalate             | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Endosulfan I (Alpha)             | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Endosulfan II (Beta)             | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Endosulfan sulfate               | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Endrin                           | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Endrin aldehyde                  | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| EPTC                             | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |

Eurofins Eaton Monrovia

# Client Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

**Client Sample ID: Aiea Gulch Wells Pump 2**

**Lab Sample ID: 380-32096-1**

Date Collected: 12/19/22 10:00

Matrix: Drinking Water

Date Received: 12/21/22 10:00

PWSID Number: HI0000331

**Method: EPA 525.2 - Semivolatile Organic Compounds (GC/MS) (Continued)**

| Analyte                          | Result | Qualifier | RL    | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------------------------------|--------|-----------|-------|------|---|----------------|----------------|---------|
| Fluoranthene                     | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Fluorene                         | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| gamma-Chlordane                  | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Heptachlor                       | ND     |           | 0.040 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Heptachlor epoxide (isomer B)    | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Hexachlorobenzene                | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Hexachlorocyclopentadiene        | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Indeno[1,2,3-cd]pyrene           | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Isophorone                       | ND     |           | 0.50  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Lindane                          | ND     |           | 0.040 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Malathion                        | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Methoxychlor                     | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Metolachlor                      | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Metribuzin                       | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Molinate                         | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Naphthalene                      | ND     |           | 0.30  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Parathion                        | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Pendimethalin (Penoxaline)       | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Total Permethrin (mixed isomers) | ND     |           | 0.20  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Phenanthrene                     | ND     |           | 0.040 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Propachlor                       | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Pyrene                           | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Simazine                         | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Terbacil                         | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Terbutylazine                    | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Thiobencarb                      | ND     |           | 0.20  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| trans-Nonachlor                  | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Trifluralin                      | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 14:45 | 1       |

| Tentatively Identified Compound | Est. Result | Qualifier | Unit | D | RT   | CAS No. | Prepared       | Analyzed       | Dil Fac |
|---------------------------------|-------------|-----------|------|---|------|---------|----------------|----------------|---------|
| Unknown                         | 0.73        | T J       | ug/L |   | 6.38 |         | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Unknown                         | 1.9         | T J       | ug/L |   | 7.32 |         | 12/22/22 06:58 | 12/27/22 14:45 | 1       |

| Surrogate          | %Recovery | Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|--------------------|-----------|-----------|----------|----------------|----------------|---------|
| 2-Nitro-m-xylene   | 99        |           | 70 - 130 | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Triphenylphosphate | 113       |           | 70 - 130 | 12/22/22 06:58 | 12/27/22 14:45 | 1       |
| Perylene-d12       | 103       |           | 70 - 130 | 12/22/22 06:58 | 12/27/22 14:45 | 1       |

**Method: 625 PAH Physis LL (EAL) + TICs - EPA 625 Base/Neutral and Acid Organics i**

| Analyte                    | Result | Qualifier | RL    | MDL   | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------------------------|--------|-----------|-------|-------|------|---|----------------|----------------|---------|
| 1-Methylnaphthalene        | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| 1-Methylphenanthrene       | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| 2,3,5-Trimethylnaphthalene | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| 2,6-Dimethylnaphthalene    | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| 2-Methylnaphthalene        | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Acenaphthene               | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Acenaphthylene             | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Anthracene                 | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Benz[a]anthracene          | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Benzo[a]pyrene             | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |

Eurofins Eaton Monrovia



# Client Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

**Client Sample ID: Aiea Gulch Wells Pump 2**

**Lab Sample ID: 380-32096-1**

Date Collected: 12/19/22 10:00

Matrix: Drinking Water

Date Received: 12/21/22 10:00

PWSID Number: HI0000331

**Method: 625 PAH Physis LL (EAL) + TICs - EPA 625 Base/Neutral and Acid Organics i (Continued)**

| Analyte                      | Result | Qualifier | RL    | MDL   | Unit | D | Prepared       | Analyzed       | Dil Fac |
|------------------------------|--------|-----------|-------|-------|------|---|----------------|----------------|---------|
| Benzo[b]fluoranthene         | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Benzo[e]pyrene               | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Benzo[g,h,i]perylene         | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Benzo[k]fluoranthene         | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Biphenyl                     | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Chrysene                     | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Dibenz[a,h]anthracene        | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Dibenzo[a,i]pyrene           | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Dibenzothiophene             | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Disalicylidenepropanediamine | ND     |           | 0.1   | 0.05  | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Fluoranthene                 | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Fluorene                     | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Indeno[1,2,3-cd]pyrene       | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Naphthalene                  | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Perylene                     | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Phenanthrene                 | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| Pyrene                       | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 07:38 | 1       |

| Surrogate          | %Recovery | Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|--------------------|-----------|-----------|----------|----------------|----------------|---------|
| (d10-Acenaphthene) | 82        |           | 27 - 133 | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| (d10-Phenanthrene) | 87        |           | 43 - 129 | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| (d12-Chrysene)     | 94        |           | 52 - 144 | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| (d12-Perylene)     | 85        |           | 36 - 161 | 12/26/22 00:00 | 01/08/23 07:38 | 1       |
| (d8-Naphthalene)   | 72        |           | 25 - 125 | 12/26/22 00:00 | 01/08/23 07:38 | 1       |

**Method: 8015 Gas (Purgeable) LL (EAL) - SW846 8015B Gasoline Range Organics**

| Analyte  | Result | Qualifier | RL    | MDL | Unit | D | Prepared | Analyzed       | Dil Fac |
|----------|--------|-----------|-------|-----|------|---|----------|----------------|---------|
| GASOLINE | ND     | U         | 0.020 |     | mg/L |   |          | 12/22/22 19:09 | 1       |

| Surrogate          | %Recovery | Qualifier | Limits   | Prepared | Analyzed       | Dil Fac |
|--------------------|-----------|-----------|----------|----------|----------------|---------|
| BROMOFLUOROBENZENE | 81        |           | 60 - 140 |          | 12/22/22 19:09 | 1       |

**Method: 8015 LL DRO/MRO/JP5/JP8 - 8015 - TPH DRO/ORO**

| Analyte   | Result | Qualifier | RL    | MDL | Unit | D | Prepared | Analyzed       | Dil Fac |
|-----------|--------|-----------|-------|-----|------|---|----------|----------------|---------|
| DIESEL    | ND     | U         | 0.026 |     | mg/L |   |          | 01/05/23 21:25 | 1       |
| JP5       | ND     | U         | 0.052 |     | mg/L |   |          | 01/05/23 21:25 | 1       |
| JP8       | ND     | U         | 0.052 |     | mg/L |   |          | 01/05/23 21:25 | 1       |
| MOTOR OIL | ND     | U         | 0.052 |     | mg/L |   |          | 01/05/23 21:25 | 1       |

| Surrogate    | %Recovery | Qualifier | Limits   | Prepared | Analyzed       | Dil Fac |
|--------------|-----------|-----------|----------|----------|----------------|---------|
| BROMOBENZENE | 93        |           | 60 - 130 |          | 01/05/23 21:25 | 1       |
| HEXACOSANE   | 115       |           | 60 - 130 |          | 01/05/23 21:25 | 1       |

**Client Sample ID: Aiea Wells Pump 2**

**Lab Sample ID: 380-32096-2**

Date Collected: 12/19/22 10:45

Matrix: Drinking Water

Date Received: 12/21/22 10:00

PWSID Number: HI0000331

**Method: EPA 525.2 - Semivolatile Organic Compounds (GC/MS)**

| Analyte  | Result | Qualifier | RL    | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------|--------|-----------|-------|------|---|----------------|----------------|---------|
| 2,4'-DDD | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| 2,4'-DDE | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |

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# Client Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

**Client Sample ID: Aiea Wells Pump 2**

**Lab Sample ID: 380-32096-2**

Date Collected: 12/19/22 10:45

Matrix: Drinking Water

Date Received: 12/21/22 10:00

PWSID Number: HI0000331

**Method: EPA 525.2 - Semivolatile Organic Compounds (GC/MS) (Continued)**

| Analyte                          | Result | Qualifier | RL    | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------------------------------|--------|-----------|-------|------|---|----------------|----------------|---------|
| 2,4'-DDT                         | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| 2,4-Dinitrotoluene               | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| 2,6-Dinitrotoluene               | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| 4,4'-DDD                         | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| 4,4'-DDE                         | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| 4,4'-DDT                         | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Acenaphthene                     | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Acenaphthylene                   | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Acetochlor                       | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Alachlor                         | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| alpha-BHC                        | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| alpha-Chlordane                  | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Anthracene                       | ND     |           | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Atrazine                         | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Benz(a)anthracene                | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Benzo[a]pyrene                   | ND     |           | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Benzo[b]fluoranthene             | ND     |           | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Benzo[g,h,i]perylene             | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Benzo[k]fluoranthene             | ND     |           | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| beta-BHC                         | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Bromacil                         | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Butachlor                        | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Butylbenzylphthalate             | ND     |           | 0.50  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Caffeine                         | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Chlorobenzilate                  | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Chloroneb                        | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Chlorothalonil (Draconil, Bravo) | ND     | ^3+       | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Chlorpyrifos                     | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Chrysene                         | ND     |           | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| delta-BHC                        | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Di(2-ethylhexyl)adipate          | ND     |           | 0.60  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Bis(2-ethylhexyl) phthalate      | ND     |           | 0.60  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Diazinon (Qualitative)           | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Dibenz(a,h)anthracene            | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Diclorvos (DDVP)                 | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Dieldrin                         | ND     |           | 0.20  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Diethylphthalate                 | ND     |           | 0.50  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Dimethoate                       | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Dimethylphthalate                | ND     |           | 0.50  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Di-n-butyl phthalate             | ND     |           | 0.99  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Di-n-octyl phthalate             | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Endosulfan I (Alpha)             | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Endosulfan II (Beta)             | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Endosulfan sulfate               | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Endrin                           | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Endrin aldehyde                  | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| EPTC                             | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Fluoranthene                     | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Fluorene                         | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |

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# Client Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

**Client Sample ID: Aiea Wells Pump 2**

**Lab Sample ID: 380-32096-2**

Date Collected: 12/19/22 10:45

Matrix: Drinking Water

Date Received: 12/21/22 10:00

PWSID Number: HI0000331

**Method: EPA 525.2 - Semivolatile Organic Compounds (GC/MS) (Continued)**

| Analyte                          | Result | Qualifier | RL    | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------------------------------|--------|-----------|-------|------|---|----------------|----------------|---------|
| gamma-Chlordane                  | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Heptachlor                       | ND     |           | 0.040 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Heptachlor epoxide (isomer B)    | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Hexachlorobenzene                | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Hexachlorocyclopentadiene        | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Indeno[1,2,3-cd]pyrene           | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Isophorone                       | ND     |           | 0.50  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Lindane                          | ND     |           | 0.040 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Malathion                        | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Methoxychlor                     | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Metolachlor                      | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Metribuzin                       | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Molinate                         | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Naphthalene                      | ND     |           | 0.30  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Parathion                        | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Pendimethalin (Penoxaline)       | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Total Permethrin (mixed isomers) | ND     |           | 0.20  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Phenanthrene                     | ND     |           | 0.040 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Propachlor                       | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Pyrene                           | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Simazine                         | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Terbacil                         | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Terbutylazine                    | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Thiobencarb                      | ND     |           | 0.20  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| trans-Nonachlor                  | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Trifluralin                      | ND     |           | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:05 | 1       |

| Tentatively Identified Compound | Est. Result | Qualifier | Unit | D | RT | CAS No. | Prepared       | Analyzed       | Dil Fac |
|---------------------------------|-------------|-----------|------|---|----|---------|----------------|----------------|---------|
| Tentatively Identified Compound | None        |           | ug/L |   |    |         | 12/22/22 06:58 | 12/27/22 15:05 | 1       |

| Surrogate          | %Recovery | Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|--------------------|-----------|-----------|----------|----------------|----------------|---------|
| 2-Nitro-m-xylene   | 99        |           | 70 - 130 | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Triphenylphosphate | 116       |           | 70 - 130 | 12/22/22 06:58 | 12/27/22 15:05 | 1       |
| Perylene-d12       | 102       |           | 70 - 130 | 12/22/22 06:58 | 12/27/22 15:05 | 1       |

**Method: 625 PAH Physis LL (EAL) + TICs - EPA 625 Base/Neutral and Acid Organics i**

| Analyte                    | Result | Qualifier | RL    | MDL   | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------------------------|--------|-----------|-------|-------|------|---|----------------|----------------|---------|
| 1-Methylnaphthalene        | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| 1-Methylphenanthrene       | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| 2,3,5-Trimethylnaphthalene | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| 2,6-Dimethylnaphthalene    | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| 2-Methylnaphthalene        | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Acenaphthene               | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Acenaphthylene             | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Anthracene                 | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Benz[a]anthracene          | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Benzo[a]pyrene             | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Benzo[b]fluoranthene       | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Benzo[e]pyrene             | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Benzo[g,h,i]perylene       | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |

Eurofins Eaton Monrovia

# Client Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

**Client Sample ID: Aiea Wells Pump 2**

**Lab Sample ID: 380-32096-2**

Date Collected: 12/19/22 10:45

Matrix: Drinking Water

Date Received: 12/21/22 10:00

PWSID Number: HI0000331

**Method: 625 PAH Physis LL (EAL) + TICs - EPA 625 Base/Neutral and Acid Organics i (Continued)**

| Analyte                      | Result | Qualifier | RL    | MDL   | Unit | D | Prepared       | Analyzed       | Dil Fac |
|------------------------------|--------|-----------|-------|-------|------|---|----------------|----------------|---------|
| Benzo[k]fluoranthene         | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Biphenyl                     | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Chrysene                     | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Dibenz[a,h]anthracene        | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Dibenzo[a,l]pyrene           | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Dibenzothiophene             | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Disalicylidenepropanediamine | ND     |           | 0.1   | 0.05  | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Fluoranthene                 | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Fluorene                     | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Indeno[1,2,3-cd]pyrene       | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Naphthalene                  | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Perylene                     | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Phenanthrene                 | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| Pyrene                       | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 09:22 | 1       |

| Surrogate          | %Recovery | Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|--------------------|-----------|-----------|----------|----------------|----------------|---------|
| (d10-Acenaphthene) | 86        |           | 27 - 133 | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| (d10-Phenanthrene) | 91        |           | 43 - 129 | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| (d12-Chrysene)     | 98        |           | 52 - 144 | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| (d12-Perylene)     | 91        |           | 36 - 161 | 12/26/22 00:00 | 01/08/23 09:22 | 1       |
| (d8-Naphthalene)   | 78        |           | 25 - 125 | 12/26/22 00:00 | 01/08/23 09:22 | 1       |

**Method: 8015 Gas (Purgeable) LL (EAL) - SW846 8015B Gasoline Range Organics**

| Analyte  | Result | Qualifier | RL    | MDL | Unit | D | Prepared | Analyzed       | Dil Fac |
|----------|--------|-----------|-------|-----|------|---|----------|----------------|---------|
| GASOLINE | ND     | U         | 0.020 |     | mg/L |   |          | 12/22/22 20:57 | 1       |

| Surrogate          | %Recovery | Qualifier | Limits   | Prepared | Analyzed       | Dil Fac |
|--------------------|-----------|-----------|----------|----------|----------------|---------|
| BROMOFLUOROBENZENE | 81        |           | 60 - 140 |          | 12/22/22 20:57 | 1       |

**Method: 8015 LL DRO/MRO/JP5/JP8 - 8015 - TPH DRO/ORO**

| Analyte   | Result | Qualifier | RL    | MDL | Unit | D | Prepared | Analyzed       | Dil Fac |
|-----------|--------|-----------|-------|-----|------|---|----------|----------------|---------|
| DIESEL    | ND     | U         | 0.027 |     | mg/L |   |          | 01/05/23 21:44 | 1       |
| JP5       | ND     | U         | 0.053 |     | mg/L |   |          | 01/05/23 21:44 | 1       |
| JP8       | ND     | U         | 0.053 |     | mg/L |   |          | 01/05/23 21:44 | 1       |
| MOTOR OIL | ND     | U         | 0.053 |     | mg/L |   |          | 01/05/23 21:44 | 1       |

| Surrogate    | %Recovery | Qualifier | Limits   | Prepared | Analyzed       | Dil Fac |
|--------------|-----------|-----------|----------|----------|----------------|---------|
| BROMOBENZENE | 84        |           | 60 - 130 |          | 01/05/23 21:44 | 1       |
| HEXACOSANE   | 111       |           | 60 - 130 |          | 01/05/23 21:44 | 1       |

**Client Sample ID: Halawa Wells Pump 1**

**Lab Sample ID: 380-32096-3**

Date Collected: 12/19/22 11:18

Matrix: Drinking Water

Date Received: 12/21/22 10:00

PWSID Number: HI0000331

**Method: EPA 525.2 - Semivolatile Organic Compounds (GC/MS)**

| Analyte            | Result | Qualifier | RL   | Unit | D | Prepared       | Analyzed       | Dil Fac |
|--------------------|--------|-----------|------|------|---|----------------|----------------|---------|
| 2,4'-DDD           | ND     |           | 0.10 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| 2,4'-DDE           | ND     |           | 0.10 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| 2,4'-DDT           | ND     |           | 0.10 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| 2,4-Dinitrotoluene | ND     |           | 0.10 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| 2,6-Dinitrotoluene | ND     |           | 0.10 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |

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# Client Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

**Client Sample ID: Halawa Wells Pump 1**

**Lab Sample ID: 380-32096-3**

Date Collected: 12/19/22 11:18

Matrix: Drinking Water

Date Received: 12/21/22 10:00

PWSID Number: HI0000331

**Method: EPA 525.2 - Semivolatile Organic Compounds (GC/MS) (Continued)**

| Analyte                          | Result | Qualifier | RL    | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------------------------------|--------|-----------|-------|------|---|----------------|----------------|---------|
| 4,4'-DDD                         | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| 4,4'-DDE                         | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| 4,4'-DDT                         | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Acenaphthene                     | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Acenaphthylene                   | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Acetochlor                       | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Alachlor                         | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| alpha-BHC                        | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| alpha-Chlordane                  | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Anthracene                       | ND     |           | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Atrazine                         | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Benz(a)anthracene                | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Benzo[a]pyrene                   | ND     |           | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Benzo[b]fluoranthene             | ND     |           | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Benzo[g,h,i]perylene             | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Benzo[k]fluoranthene             | ND     |           | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| beta-BHC                         | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Bromacil                         | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Butachlor                        | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Butylbenzylphthalate             | ND     |           | 0.50  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Caffeine                         | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Chlorobenzilate                  | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Chloroneb                        | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Chlorothalonil (Draconil, Bravo) | ND     | ^3+       | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Chlorpyrifos                     | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Chrysene                         | ND     |           | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| delta-BHC                        | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Di(2-ethylhexyl)adipate          | ND     |           | 0.60  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Bis(2-ethylhexyl) phthalate      | ND     |           | 0.60  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Diazinon (Qualitative)           | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Dibenz(a,h)anthracene            | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Diclorvos (DDVP)                 | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Dieldrin                         | ND     |           | 0.20  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Diethylphthalate                 | ND     |           | 0.50  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Dimethoate                       | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Dimethylphthalate                | ND     |           | 0.50  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Di-n-butyl phthalate             | ND     |           | 1.0   | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Di-n-octyl phthalate             | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Endosulfan I (Alpha)             | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Endosulfan II (Beta)             | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Endosulfan sulfate               | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Endrin                           | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Endrin aldehyde                  | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| EPTC                             | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Fluoranthene                     | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Fluorene                         | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| gamma-Chlordane                  | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Heptachlor                       | ND     |           | 0.040 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Heptachlor epoxide (isomer B)    | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |

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# Client Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

**Client Sample ID: Halawa Wells Pump 1**

**Lab Sample ID: 380-32096-3**

Date Collected: 12/19/22 11:18

Matrix: Drinking Water

Date Received: 12/21/22 10:00

PWSID Number: HI0000331

**Method: EPA 525.2 - Semivolatile Organic Compounds (GC/MS) (Continued)**

| Analyte                          | Result | Qualifier | RL    | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------------------------------|--------|-----------|-------|------|---|----------------|----------------|---------|
| Hexachlorobenzene                | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Hexachlorocyclopentadiene        | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Indeno[1,2,3-cd]pyrene           | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Isophorone                       | ND     |           | 0.50  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Lindane                          | ND     |           | 0.040 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Malathion                        | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Methoxychlor                     | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Metolachlor                      | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Metribuzin                       | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Molinate                         | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Naphthalene                      | ND     |           | 0.30  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Parathion                        | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Pendimethalin (Penoxaline)       | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Total Permethrin (mixed isomers) | ND     |           | 0.20  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Phenanthrene                     | ND     |           | 0.040 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Propachlor                       | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Pyrene                           | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Simazine                         | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Terbacil                         | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Terbutylazine                    | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Thiobencarb                      | ND     |           | 0.20  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| trans-Nonachlor                  | ND     |           | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Trifluralin                      | ND     |           | 0.10  | ug/L |   | 12/22/22 06:58 | 12/27/22 15:26 | 1       |

| Tentatively Identified Compound | Est. Result | Qualifier | Unit | D | RT | CAS No. | Prepared       | Analyzed       | Dil Fac |
|---------------------------------|-------------|-----------|------|---|----|---------|----------------|----------------|---------|
| Tentatively Identified Compound | None        |           | ug/L |   |    |         | 12/22/22 06:58 | 12/27/22 15:26 | 1       |

| Surrogate          | %Recovery | Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|--------------------|-----------|-----------|----------|----------------|----------------|---------|
| 2-Nitro-m-xylene   | 98        |           | 70 - 130 | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Triphenylphosphate | 118       |           | 70 - 130 | 12/22/22 06:58 | 12/27/22 15:26 | 1       |
| Perylene-d12       | 104       |           | 70 - 130 | 12/22/22 06:58 | 12/27/22 15:26 | 1       |

**Method: 625 PAH Physis LL (EAL) + TICs - EPA 625 Base/Neutral and Acid Organics i**

| Analyte                    | Result | Qualifier | RL    | MDL   | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------------------------|--------|-----------|-------|-------|------|---|----------------|----------------|---------|
| 1-Methylnaphthalene        | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| 1-Methylphenanthrene       | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| 2,3,5-Trimethylnaphthalene | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| 2,6-Dimethylnaphthalene    | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| 2-Methylnaphthalene        | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Acenaphthene               | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Acenaphthylene             | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Anthracene                 | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Benz[a]anthracene          | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Benzo[a]pyrene             | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Benzo[b]fluoranthene       | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Benzo[e]pyrene             | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Benzo[g,h,i]perylene       | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Benzo[k]fluoranthene       | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Biphenyl                   | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Chrysene                   | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |

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# Client Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

**Client Sample ID: Halawa Wells Pump 1**

**Lab Sample ID: 380-32096-3**

Date Collected: 12/19/22 11:18

Matrix: Drinking Water

Date Received: 12/21/22 10:00

PWSID Number: HI0000331

**Method: 625 PAH Physis LL (EAL) + TICs - EPA 625 Base/Neutral and Acid Organics i (Continued)**

| Analyte                    | Result | Qualifier | RL    | MDL   | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------------------------|--------|-----------|-------|-------|------|---|----------------|----------------|---------|
| Dibenz[a,h]anthracene      | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Dibenzo[a,l]pyrene         | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Dibenzothiophene           | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Disalicylidenepranediamine | ND     |           | 0.1   | 0.05  | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Fluoranthene               | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Fluorene                   | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Indeno[1,2,3-cd]pyrene     | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Naphthalene                | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Perylene                   | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Phenanthrene               | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| Pyrene                     | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 11:07 | 1       |

| Surrogate          | %Recovery | Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|--------------------|-----------|-----------|----------|----------------|----------------|---------|
| (d10-Acenaphthene) | 90        |           | 27 - 133 | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| (d10-Phenanthrene) | 94        |           | 43 - 129 | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| (d12-Chrysene)     | 99        |           | 52 - 144 | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| (d12-Perylene)     | 86        |           | 36 - 161 | 12/26/22 00:00 | 01/08/23 11:07 | 1       |
| (d8-Naphthalene)   | 80        |           | 25 - 125 | 12/26/22 00:00 | 01/08/23 11:07 | 1       |

**Method: 8015 Gas (Purgeable) LL (EAL) - SW846 8015B Gasoline Range Organics**

| Analyte  | Result | Qualifier | RL    | MDL | Unit | D | Prepared | Analyzed       | Dil Fac |
|----------|--------|-----------|-------|-----|------|---|----------|----------------|---------|
| GASOLINE | ND     | U         | 0.020 |     | mg/L |   |          | 12/22/22 21:33 | 1       |

| Surrogate          | %Recovery | Qualifier | Limits   | Prepared | Analyzed       | Dil Fac |
|--------------------|-----------|-----------|----------|----------|----------------|---------|
| BROMOFLUOROBENZENE | 81        |           | 60 - 140 |          | 12/22/22 21:33 | 1       |

**Method: 8015 LL DRO/MRO/JP5/JP8 - 8015 - TPH DRO/ORO**

| Analyte   | Result | Qualifier | RL    | MDL | Unit | D | Prepared | Analyzed       | Dil Fac |
|-----------|--------|-----------|-------|-----|------|---|----------|----------------|---------|
| DIESEL    | ND     | U         | 0.027 |     | mg/L |   |          | 01/05/23 22:02 | 1       |
| JP5       | ND     | U         | 0.053 |     | mg/L |   |          | 01/05/23 22:02 | 1       |
| JP8       | ND     | U         | 0.053 |     | mg/L |   |          | 01/05/23 22:02 | 1       |
| MOTOR OIL | ND     | U         | 0.053 |     | mg/L |   |          | 01/05/23 22:02 | 1       |

| Surrogate    | %Recovery | Qualifier | Limits   | Prepared | Analyzed       | Dil Fac |
|--------------|-----------|-----------|----------|----------|----------------|---------|
| BROMOBENZENE | 83        |           | 60 - 130 |          | 01/05/23 22:02 | 1       |
| HEXACOSANE   | 97        |           | 60 - 130 |          | 01/05/23 22:02 | 1       |

**Client Sample ID: TB: Aiea Gulch Wells Pump 2**

**Lab Sample ID: 380-32096-4**

Date Collected: 12/19/22 10:00

Matrix: Drinking Water

Date Received: 12/21/22 10:00

**Method: 8015 Gas (Purgeable) LL (EAL) - SW846 8015B Gasoline Range Organics**

| Analyte  | Result | Qualifier | RL    | MDL | Unit | D | Prepared | Analyzed       | Dil Fac |
|----------|--------|-----------|-------|-----|------|---|----------|----------------|---------|
| GASOLINE | ND     | U         | 0.020 |     | mg/L |   |          | 12/22/22 22:08 | 1       |

| Surrogate          | %Recovery | Qualifier | Limits   | Prepared | Analyzed       | Dil Fac |
|--------------------|-----------|-----------|----------|----------|----------------|---------|
| BROMOFLUOROBENZENE | 80        |           | 60 - 140 |          | 12/22/22 22:08 | 1       |

# Client Sample Results

Client: City & County of Honolulu  
 Project/Site: RED-HILL

Job ID: 380-32096-1

## Client Sample ID: TB: Aiea Wells Pump 2

Lab Sample ID: 380-32096-5

Date Collected: 12/19/22 10:45

Matrix: Drinking Water

Date Received: 12/21/22 10:00

### Method: 8015 Gas (Purgeable) LL (EAL) - SW846 8015B Gasoline Range Organics

| Analyte            | Result    | Qualifier | RL       | MDL | Unit | D | Prepared | Analyzed       | Dil Fac |
|--------------------|-----------|-----------|----------|-----|------|---|----------|----------------|---------|
| GASOLINE           | ND        | U         | 0.020    |     | mg/L |   |          | 12/22/22 22:44 | 1       |
| Surrogate          | %Recovery | Qualifier | Limits   |     |      |   | Prepared | Analyzed       | Dil Fac |
| BROMOFLUOROBENZENE | 81        |           | 60 - 140 |     |      |   |          | 12/22/22 22:44 | 1       |

## Client Sample ID: TB: Halawa Wells Pump 1

Lab Sample ID: 380-32096-6

Date Collected: 12/19/22 11:18

Matrix: Drinking Water

Date Received: 12/21/22 10:00

### Method: 8015 Gas (Purgeable) LL (EAL) - SW846 8015B Gasoline Range Organics

| Analyte            | Result    | Qualifier | RL       | MDL | Unit | D | Prepared | Analyzed       | Dil Fac |
|--------------------|-----------|-----------|----------|-----|------|---|----------|----------------|---------|
| GASOLINE           | ND        | U         | 0.020    |     | mg/L |   |          | 12/22/22 23:20 | 1       |
| Surrogate          | %Recovery | Qualifier | Limits   |     |      |   | Prepared | Analyzed       | Dil Fac |
| BROMOFLUOROBENZENE | 80        |           | 60 - 140 |     |      |   |          | 12/22/22 23:20 | 1       |



# Action Limit Summary

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

**Client Sample ID: Aiea Gulch Wells Pump 2**

**Lab Sample ID: 380-32096-1**

**PWSID Number: HI0000331**

## Compliance Check

The results obtained from the analytical testing of this data set were checked against compliance limits received from the client. Any results at or above the compliance limits have been highlighted for your convenience.

| Analyte                       | Result | Qualifier | Unit | EPAMCL | RL    | Method | Prep Type |
|-------------------------------|--------|-----------|------|--------|-------|--------|-----------|
|                               |        |           |      | Limit  |       |        |           |
| Alachlor                      | ND     |           | ug/L | 2      | 0.050 | 525.2  | Total/NA  |
| Atrazine                      | ND     |           | ug/L | 3      | 0.050 | 525.2  | Total/NA  |
| Benzo[a]pyrene                | ND     |           | ug/L | 0.2    | 0.020 | 525.2  | Total/NA  |
| Di(2-ethylhexyl)adipate       | ND     |           | ug/L | 400    | 0.60  | 525.2  | Total/NA  |
| Bis(2-ethylhexyl) phthalate   | ND     |           | ug/L | 6      | 0.60  | 525.2  | Total/NA  |
| Endrin                        | ND     |           | ug/L | 2      | 0.10  | 525.2  | Total/NA  |
| Heptachlor                    | ND     |           | ug/L | 0.4    | 0.040 | 525.2  | Total/NA  |
| Heptachlor epoxide (isomer B) | ND     |           | ug/L | 0.2    | 0.050 | 525.2  | Total/NA  |
| Hexachlorobenzene             | ND     |           | ug/L | 1      | 0.050 | 525.2  | Total/NA  |
| Hexachlorocyclopentadiene     | ND     |           | ug/L | 50     | 0.050 | 525.2  | Total/NA  |
| Lindane                       | ND     |           | ug/L | 0.2    | 0.040 | 525.2  | Total/NA  |
| Methoxychlor                  | ND     |           | ug/L | 40     | 0.10  | 525.2  | Total/NA  |
| Simazine                      | ND     |           | ug/L | 4      | 0.050 | 525.2  | Total/NA  |

**Client Sample ID: Aiea Wells Pump 2**

**Lab Sample ID: 380-32096-2**

**PWSID Number: HI0000331**

## Compliance Check

The results obtained from the analytical testing of this data set were checked against compliance limits received from the client. Any results at or above the compliance limits have been highlighted for your convenience.

| Analyte                       | Result | Qualifier | Unit | EPAMCL | RL    | Method | Prep Type |
|-------------------------------|--------|-----------|------|--------|-------|--------|-----------|
|                               |        |           |      | Limit  |       |        |           |
| Alachlor                      | ND     |           | ug/L | 2      | 0.050 | 525.2  | Total/NA  |
| Atrazine                      | ND     |           | ug/L | 3      | 0.050 | 525.2  | Total/NA  |
| Benzo[a]pyrene                | ND     |           | ug/L | 0.2    | 0.020 | 525.2  | Total/NA  |
| Di(2-ethylhexyl)adipate       | ND     |           | ug/L | 400    | 0.60  | 525.2  | Total/NA  |
| Bis(2-ethylhexyl) phthalate   | ND     |           | ug/L | 6      | 0.60  | 525.2  | Total/NA  |
| Endrin                        | ND     |           | ug/L | 2      | 0.099 | 525.2  | Total/NA  |
| Heptachlor                    | ND     |           | ug/L | 0.4    | 0.040 | 525.2  | Total/NA  |
| Heptachlor epoxide (isomer B) | ND     |           | ug/L | 0.2    | 0.050 | 525.2  | Total/NA  |
| Hexachlorobenzene             | ND     |           | ug/L | 1      | 0.050 | 525.2  | Total/NA  |
| Hexachlorocyclopentadiene     | ND     |           | ug/L | 50     | 0.050 | 525.2  | Total/NA  |
| Lindane                       | ND     |           | ug/L | 0.2    | 0.040 | 525.2  | Total/NA  |
| Methoxychlor                  | ND     |           | ug/L | 40     | 0.099 | 525.2  | Total/NA  |
| Simazine                      | ND     |           | ug/L | 4      | 0.050 | 525.2  | Total/NA  |

**Client Sample ID: Halawa Wells Pump 1**

**Lab Sample ID: 380-32096-3**

**PWSID Number: HI0000331**

## Compliance Check

The results obtained from the analytical testing of this data set were checked against compliance limits received from the client. Any results at or above the compliance limits have been highlighted for your convenience.

| Analyte  | Result | Qualifier | Unit | EPAMCL | RL    | Method | Prep Type |
|----------|--------|-----------|------|--------|-------|--------|-----------|
|          |        |           |      | Limit  |       |        |           |
| Alachlor | ND     |           | ug/L | 2      | 0.050 | 525.2  | Total/NA  |
| Atrazine | ND     |           | ug/L | 3      | 0.050 | 525.2  | Total/NA  |

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# Action Limit Summary

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

**Client Sample ID: Halawa Wells Pump 1 (Continued)**

**Lab Sample ID: 380-32096-3**

**PWSID Number: HI0000331**

## Compliance Check

The results obtained from the analytical testing of this data set were checked against compliance limits received from the client. Any results at or above the compliance limits have been highlighted for your convenience.

| Analyte                       | Result | Qualifier | Unit | EPAMCL | RL    | Method | Prep Type |
|-------------------------------|--------|-----------|------|--------|-------|--------|-----------|
|                               |        |           |      | Limit  |       |        |           |
| Benzo[a]pyrene                | ND     |           | ug/L | 0.2    | 0.020 | 525.2  | Total/NA  |
| Di(2-ethylhexyl)adipate       | ND     |           | ug/L | 400    | 0.60  | 525.2  | Total/NA  |
| Bis(2-ethylhexyl) phthalate   | ND     |           | ug/L | 6      | 0.60  | 525.2  | Total/NA  |
| Endrin                        | ND     |           | ug/L | 2      | 0.10  | 525.2  | Total/NA  |
| Heptachlor                    | ND     |           | ug/L | 0.4    | 0.040 | 525.2  | Total/NA  |
| Heptachlor epoxide (isomer B) | ND     |           | ug/L | 0.2    | 0.050 | 525.2  | Total/NA  |
| Hexachlorobenzene             | ND     |           | ug/L | 1      | 0.050 | 525.2  | Total/NA  |
| Hexachlorocyclopentadiene     | ND     |           | ug/L | 50     | 0.050 | 525.2  | Total/NA  |
| Lindane                       | ND     |           | ug/L | 0.2    | 0.040 | 525.2  | Total/NA  |
| Methoxychlor                  | ND     |           | ug/L | 40     | 0.10  | 525.2  | Total/NA  |
| Simazine                      | ND     |           | ug/L | 4      | 0.050 | 525.2  | Total/NA  |

# Surrogate Summary

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 525.2 - Semivolatile Organic Compounds (GC/MS)

Matrix: Drinking Water

Prep Type: Total/NA

| Lab Sample ID | Client Sample ID        | Percent Surrogate Recovery (Acceptance Limits) |                 |                 |
|---------------|-------------------------|--|-----------------|-----------------|
|               |                         | 2NMX<br>(70-130)                               | TPP<br>(70-130) | PRY<br>(70-130) |
| 380-32096-1   | Aiea Gulch Wells Pump 2 | 99   | 113             | 103             |
| 380-32096-2   | Aiea Wells Pump 2       | 99   | 116             | 102             |
| 380-32096-3   | Halawa Wells Pump 1     | 98   | 118             | 104             |

**Surrogate Legend**

2NMX = 2-Nitro-m-xylene  
TPP = Triphenylphosphate  
PRY = Perylene-d12

## Method: 525.2 - Semivolatile Organic Compounds (GC/MS)

Matrix: Water

Prep Type: Total/NA

| Lab Sample ID       | Client Sample ID       | Percent Surrogate Recovery (Acceptance Limits) |                 |                 |
|---------------------|------------------------|--|-----------------|-----------------|
|                     |                        | 2NMX<br>(70-130)                               | TPP<br>(70-130) | PRY<br>(70-130) |
| 380-32064-C-1-A MS  | Matrix Spike           | 101  | 106             | 101             |
| 380-32064-I-1-A MSD | Matrix Spike Duplicate | 98   | 108             | 102             |
| LCS 380-27725/3-A   | Lab Control Sample     | 98   | 113             | 100             |
| LCS 380-27725/4-A   | Lab Control Sample Dup | 101  | 108             | 99              |
| MB 380-27725/1-A    | Method Blank           | 100  | 106             | 94              |
| MRL 380-27725/2-A   | Lab Control Sample     | 100  | 116             | 98              |

**Surrogate Legend**

2NMX = 2-Nitro-m-xylene  
TPP = Triphenylphosphate  
PRY = Perylene-d12

## Method: 625 PAH Physis LL (EAL) + TICs - EPA 625 Base/Neutral and Acid Organics i

Matrix: BlankMatrix

Prep Type: Total/NA

| Lab Sample ID | Client Sample ID       | Percent Surrogate Recovery (Acceptance Limits) |                      |                 |                 |                 |
|---------------|------------------------|--|----------------------|-----------------|-----------------|-----------------|
|               |                        | Acenaphtl<br>(27-133)                          | Phenanth<br>(43-129) | CRY<br>(52-144) | NPT<br>(25-125) | PRY<br>(36-161) |
| 102964-B1     | Method Blank           | 92   | 94                   | 99              | 87              | 103             |
| 102964-BS1    | Lab Control Sample     | 93   | 96                   | 97              | 84              | 99              |
| 102964-BS2    | Lab Control Sample Dup | 92   | 95                   | 95              | 84              | 94              |

**Surrogate Legend**

(d10-Acenaphthene) = (d10-Acenaphthene)  
(d10-Phenanthrene) = (d10-Phenanthrene)  
CRY = (d12-Chrysene)  
NPT = (d8-Naphthalene)  
PRY = (d12-Perylene)

## Method: 625 PAH Physis LL (EAL) + TICs - EPA 625 Base/Neutral and Acid Organics i

Matrix: Drinking Water

Prep Type: Total/NA

| Lab Sample ID | Client Sample ID        | Percent Surrogate Recovery (Acceptance Limits) |                      |                 |                 |                 |
|---------------|-------------------------|--|----------------------|-----------------|-----------------|-----------------|
|               |                         | Acenaphtl<br>(27-133)                          | Phenanth<br>(43-129) | CRY<br>(52-144) | NPT<br>(25-125) | PRY<br>(36-161) |
| 380-32096-1   | Aiea Gulch Wells Pump 2 | 82   | 87                   | 94              | 72              | 85              |
| 380-32096-2   | Aiea Wells Pump 2       | 86   | 91                   | 98              | 78              | 91              |

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# Surrogate Summary

Client: City & County of Honolulu  
 Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 625 PAH Physis LL (EAL) + TICs - EPA 625 Base/Neutral and Acid Organics i (Continued)

Matrix: Drinking Water

Prep Type: Total/NA

| Lab Sample ID | Client Sample ID    | Percent Surrogate Recovery (Acceptance Limits) |                      |                 |                 |                 |
|---------------|---------------------|--|----------------------|-----------------|-----------------|-----------------|
|               |                     | Acenaphtl<br>(27-133)                          | Phenanth<br>(43-129) | CRY<br>(52-144) | NPT<br>(25-125) | PRY<br>(36-161) |
| 380-32096-3   | Halawa Wells Pump 1 | 90   | 94                   | 99              | 80              | 86              |

**Surrogate Legend**

(d10-Acenaphthene) = (d10-Acenaphthene)

(d10-Phenanthrene) = (d10-Phenanthrene)

CRY = (d12-Chrysene)

NPT = (d8-Naphthalene)

PRY = (d12-Perylene)

## Method: 8015 Gas (Purgeable) LL (EAL) - SW846 8015B Gasoline Range Organics

Matrix: Drinking Water

Prep Type: Total/NA

| Lab Sample ID | Client Sample ID            | Percent Surrogate Recovery (Acceptance Limits) |
|---------------|-----------------------------|--|
|               |                             | BFB<br>(60-140)                                |
| 380-32096-1   | Aiea Gulch Wells Pump 2     | 81   |
| 380-32096-2   | Aiea Wells Pump 2           | 81   |
| 380-32096-3   | Halawa Wells Pump 1         | 81   |
| 380-32096-4   | TB: Aiea Gulch Wells Pump 2 | 80   |
| 380-32096-5   | TB: Aiea Wells Pump 2       | 81   |
| 380-32096-6   | TB: Halawa Wells Pump 1     | 80   |

**Surrogate Legend**

BFB = BROMOFLUOROBENZENE

## Method: 8015 Gas (Purgeable) LL (EAL) - SW846 8015B Gasoline Range Organics

Matrix: WATER

Prep Type: Total/NA

| Lab Sample ID | Client Sample ID   | Percent Surrogate Recovery (Acceptance Limits) |
|---------------|--------------------|--|
|               |                    | BFB<br>(70-130)                                |
| 22VG39L08C    | LCD                | 109  |
| 22VG39L08L    | Lab Control Sample | 110  |

**Surrogate Legend**

BFB = BROMOFLUOROBENZENE

## Method: 8015 Gas (Purgeable) LL (EAL) - SW846 8015B Gasoline Range Organics

Matrix: WATER

Prep Type: Total/NA

| Lab Sample ID | Client Sample ID       | Percent Surrogate Recovery (Acceptance Limits) |
|---------------|------------------------|--|
|               |                        | BFB<br>(60-140)                                |
| 22L294-01M    | Matrix Spike           | 108  |
| 22L294-01S    | Matrix Spike Duplicate | 110  |

**Surrogate Legend**

BFB = BROMOFLUOROBENZENE

# Surrogate Summary

Client: City & County of Honolulu  
 Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 8015 Gas (Purgeable) LL (EAL) - SW846 8015B Gasoline Range Organics

Matrix: WATER

Prep Type: Total/NA

### Percent Surrogate Recovery (Acceptance Limits)

BFB

| Lab Sample ID | Client Sample ID |
|---------------|------------------|
| 22VG39L08B    | Method Blank     |

**Surrogate Legend**

BFB = BROMOFLUOROBENZENE

## Method: 8015 LL DRO/MRO/JP5/JP8 - 8015 - TPH DRO/ORO

Matrix: Drinking Water

Prep Type: Total/NA

### Percent Surrogate Recovery (Acceptance Limits)

| Lab Sample ID | Client Sample ID        | BB<br>(60-130) | XACOSAI<br>(60-130) |
|---------------|-------------------------|----------------|---------------------|
| 380-32096-1   | Aiea Gulch Wells Pump 2 | 93             | 115                 |
| 380-32096-2   | Aiea Wells Pump 2       | 84             | 111                 |
| 380-32096-3   | Halawa Wells Pump 1     | 83             | 97                  |

**Surrogate Legend**

BB = BROMOBENZENE

HEXACOSANE = HEXACOSANE

## Method: 8015 LL DRO/MRO/JP5/JP8 - 8015 - TPH DRO/ORO

Matrix: WATER

Prep Type: Total/NA

### Percent Surrogate Recovery (Acceptance Limits)

BB XACOSAI

| Lab Sample ID | Client Sample ID |
|---------------|------------------|
| 22DSL040WB    | Method Blank     |

**Surrogate Legend**

BB = BROMOBENZENE

HEXACOSANE = HEXACOSANE

## Method: 8015 LL DRO/MRO/JP5/JP8 - 8015 - TPH DRO/ORO

Matrix: WATER

Prep Type: Total/NA

### Percent Surrogate Recovery (Acceptance Limits)

| Lab Sample ID | Client Sample ID   | BB<br>(60-130) | XACOSAI<br>(60-130) |
|---------------|--------------------|----------------|---------------------|
| 22DSL040WL    | Lab Control Sample | 102            | 106                 |
| 22J5L040WL    | Lab Control Sample | 105            | 97                  |
| 22J8L040WL    | Lab Control Sample | 98             | 102                 |

**Surrogate Legend**

BB = BROMOBENZENE

HEXACOSANE = HEXACOSANE

# QC Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 525.2 - Semivolatile Organic Compounds (GC/MS)

**Lab Sample ID: MB 380-27725/1-A**  
**Matrix: Water**  
**Analysis Batch: 28001**

**Client Sample ID: Method Blank**  
**Prep Type: Total/NA**  
**Prep Batch: 27725**

| Analyte                          | MB<br>Result | MB<br>Qualifier | RL    | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------------------------------|--------------|-----------------|-------|------|---|----------------|----------------|---------|
| 2,4'-DDD                         | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| 2,4'-DDE                         | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| 2,4'-DDT                         | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| 2,4-Dinitrotoluene               | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| 2,6-Dinitrotoluene               | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| 4,4'-DDD                         | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| 4,4'-DDE                         | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| 4,4'-DDT                         | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Acenaphthene                     | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Acenaphthylene                   | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Acetochlor                       | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Alachlor                         | ND           |                 | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| alpha-BHC                        | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| alpha-Chlordane                  | ND           |                 | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Anthracene                       | ND           |                 | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Atrazine                         | ND           |                 | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Benz(a)anthracene                | ND           |                 | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Benzo[a]pyrene                   | ND           |                 | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Benzo[b]fluoranthene             | ND           |                 | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Benzo[g,h,i]perylene             | ND           |                 | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Benzo[k]fluoranthene             | ND           |                 | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| beta-BHC                         | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Bromacil                         | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Butachlor                        | ND           |                 | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Butylbenzylphthalate             | ND           |                 | 0.50  | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Caffeine                         | ND           |                 | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Chlorobenzilate                  | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Chloroneb                        | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Chlorothalonil (Draconil, Bravo) | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Chlorpyrifos                     | ND           |                 | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Chrysene                         | ND           |                 | 0.020 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| delta-BHC                        | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Di(2-ethylhexyl)adipate          | ND           |                 | 0.59  | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Bis(2-ethylhexyl) phthalate      | ND           |                 | 0.59  | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Diazinon (Qualitative)           | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Dibenz(a,h)anthracene            | ND           |                 | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Diclorvos (DDVP)                 | ND           |                 | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Dieldrin                         | ND           |                 | 0.20  | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Diethylphthalate                 | ND           |                 | 0.50  | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Dimethoate                       | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Dimethylphthalate                | ND           |                 | 0.50  | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Di-n-butyl phthalate             | ND           |                 | 0.99  | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Di-n-octyl phthalate             | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Endosulfan I (Alpha)             | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Endosulfan II (Beta)             | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Endosulfan sulfate               | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Endrin                           | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Endrin aldehyde                  | ND           |                 | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |

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# QC Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 525.2 - Semivolatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: MB 380-27725/1-A**  
**Matrix: Water**  
**Analysis Batch: 28001**

**Client Sample ID: Method Blank**  
**Prep Type: Total/NA**  
**Prep Batch: 27725**

| Analyte                          | MB Result | MB Qualifier | RL    | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------------------------------|-----------|--------------|-------|------|---|----------------|----------------|---------|
| EPTC                             | ND        |              | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Fluoranthene                     | ND        |              | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Fluorene                         | ND        |              | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| gamma-Chlordane                  | ND        |              | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Heptachlor                       | ND        |              | 0.040 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Heptachlor epoxide (isomer B)    | ND        |              | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Hexachlorobenzene                | ND        |              | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Hexachlorocyclopentadiene        | ND        |              | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Indeno[1,2,3-cd]pyrene           | ND        |              | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Isophorone                       | ND        |              | 0.50  | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Lindane                          | ND        |              | 0.040 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Malathion                        | ND        |              | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Methoxychlor                     | ND        |              | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Metolachlor                      | ND        |              | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Metribuzin                       | ND        |              | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Molinate                         | ND        |              | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Naphthalene                      | ND        |              | 0.30  | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Parathion                        | ND        |              | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Pendimethalin (Penoxaline)       | ND        |              | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Total Permethrin (mixed isomers) | ND        |              | 0.20  | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Phenanthrene                     | ND        |              | 0.040 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Propachlor                       | ND        |              | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Pyrene                           | ND        |              | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Simazine                         | ND        |              | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Terbacil                         | ND        |              | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Terbutylazine                    | ND        |              | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Thiobencarb                      | ND        |              | 0.20  | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| trans-Nonachlor                  | ND        |              | 0.050 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| Trifluralin                      | ND        |              | 0.099 | ug/L |   | 12/22/22 06:58 | 12/27/22 11:24 | 1       |

| <i>Tentatively Identified Compound</i> | MB Est. Result | MB Qualifier | Unit | D | RT | CAS No. | Prepared       | Analyzed       | Dil Fac |
|--|----------------|--------------|------|---|----|---------|----------------|----------------|---------|
| <i>Tentatively Identified Compound</i> | None           |              | ug/L |   |    |         | 12/22/22 06:58 | 12/27/22 11:24 | 1       |

| <i>Surrogate</i>          | MB %Recovery | MB Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|---------------------------|--------------|--------------|----------|----------------|----------------|---------|
| <i>2-Nitro-m-xylene</i>   | 100          |              | 70 - 130 | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| <i>Triphenylphosphate</i> | 106          |              | 70 - 130 | 12/22/22 06:58 | 12/27/22 11:24 | 1       |
| <i>Perylene-d12</i>       | 94           |              | 70 - 130 | 12/22/22 06:58 | 12/27/22 11:24 | 1       |

**Lab Sample ID: LCS 380-27725/3-A**  
**Matrix: Water**  
**Analysis Batch: 28001**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 27725**

| Analyte            | Spike Added | LCS Result | LCS Qualifier | Unit | D | %Rec | %Rec Limits |
|--------------------|-------------|------------|---------------|------|---|------|-------------|
| 2,4'-DDD           | 1.98        | 1.96       |               | ug/L |   | 99   | 70 - 130    |
| 2,4'-DDE           | 1.98        | 1.91       |               | ug/L |   | 96   | 70 - 130    |
| 2,4'-DDT           | 1.98        | 2.09       |               | ug/L |   | 105  | 70 - 130    |
| 2,4-Dinitrotoluene | 1.98        | 1.80       |               | ug/L |   | 91   | 70 - 130    |
| 2,6-Dinitrotoluene | 1.98        | 1.93       |               | ug/L |   | 98   | 70 - 130    |

Eurofins Eaton Monrovia

# QC Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 525.2 - Semivolatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: LCS 380-27725/3-A**  
**Matrix: Water**  
**Analysis Batch: 28001**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 27725**

| Analyte                          | Spike Added | LCS Result | LCS Qualifier | Unit | D | %Rec | %Rec Limits |
|----------------------------------|-------------|------------|---------------|------|---|------|-------------|
| 4,4'-DDD                         | 1.98        | 2.18       |               | ug/L |   | 110  | 70 - 130    |
| 4,4'-DDE                         | 1.98        | 2.06       |               | ug/L |   | 104  | 70 - 130    |
| 4,4'-DDT                         | 1.98        | 2.17       |               | ug/L |   | 110  | 70 - 130    |
| Acenaphthene                     | 1.98        | 1.93       |               | ug/L |   | 98   | 70 - 130    |
| Acenaphthylene                   | 1.98        | 1.90       |               | ug/L |   | 96   | 70 - 130    |
| Acetochlor                       | 1.98        | 2.09       |               | ug/L |   | 105  | 70 - 130    |
| Alachlor                         | 1.98        | 2.00       |               | ug/L |   | 101  | 70 - 130    |
| alpha-BHC                        | 1.98        | 1.96       |               | ug/L |   | 99   | 70 - 130    |
| alpha-Chlordane                  | 1.98        | 1.87       |               | ug/L |   | 94   | 70 - 130    |
| Anthracene                       | 1.98        | 1.88       |               | ug/L |   | 95   | 70 - 130    |
| Atrazine                         | 1.98        | 1.98       |               | ug/L |   | 100  | 70 - 130    |
| Benz(a)anthracene                | 1.98        | 2.19       |               | ug/L |   | 111  | 70 - 130    |
| Benzo[a]pyrene                   | 1.98        | 2.13       |               | ug/L |   | 107  | 70 - 130    |
| Benzo[b]fluoranthene             | 1.98        | 2.16       |               | ug/L |   | 109  | 70 - 130    |
| Benzo[g,h,i]perylene             | 1.98        | 2.05       |               | ug/L |   | 103  | 70 - 130    |
| Benzo[k]fluoranthene             | 1.98        | 2.10       |               | ug/L |   | 106  | 70 - 130    |
| beta-BHC                         | 1.98        | 2.00       |               | ug/L |   | 101  | 70 - 130    |
| Bromacil                         | 1.98        | 2.26       |               | ug/L |   | 114  | 70 - 130    |
| Butachlor                        | 1.98        | 2.15       |               | ug/L |   | 109  | 70 - 130    |
| Butylbenzylphthalate             | 1.98        | 2.26       |               | ug/L |   | 114  | 70 - 130    |
| Caffeine                         | 1.98        | 1.62       |               | ug/L |   | 82   | 45 - 137    |
| Chlorobenzilate                  | 1.98        | 2.21       |               | ug/L |   | 111  | 70 - 130    |
| Chloroneb                        | 1.98        | 1.89       |               | ug/L |   | 95   | 70 - 130    |
| Chlorothalonil (Draconil, Bravo) | 1.98        | 1.92       |               | ug/L |   | 97   | 70 - 130    |
| Chlorpyrifos                     | 1.98        | 2.02       |               | ug/L |   | 102  | 70 - 130    |
| Chrysene                         | 1.98        | 1.91       |               | ug/L |   | 96   | 70 - 130    |
| delta-BHC                        | 1.98        | 2.05       |               | ug/L |   | 103  | 70 - 130    |
| Di(2-ethylhexyl)adipate          | 1.98        | 2.24       |               | ug/L |   | 113  | 70 - 130    |
| Bis(2-ethylhexyl) phthalate      | 1.98        | 1.88       |               | ug/L |   | 95   | 70 - 130    |
| Diazinon (Qualitative)           | 1.98        | 1.79       |               | ug/L |   | 90   | 15 - 132    |
| Dibenz(a,h)anthracene            | 1.98        | 2.21       |               | ug/L |   | 112  | 70 - 130    |
| Diclorvos (DDVP)                 | 1.98        | 2.19       |               | ug/L |   | 111  | 70 - 130    |
| Dieldrin                         | 1.98        | 2.06       |               | ug/L |   | 104  | 70 - 130    |
| Diethylphthalate                 | 1.98        | 1.93       |               | ug/L |   | 97   | 70 - 130    |
| Dimethoate                       | 1.98        | 1.30       |               | ug/L |   | 65   | 35 - 100    |
| Dimethylphthalate                | 1.98        | 2.02       |               | ug/L |   | 102  | 70 - 130    |
| Di-n-butyl phthalate             | 3.96        | 3.90       |               | ug/L |   | 98   | 70 - 130    |
| Di-n-octyl phthalate             | 1.98        | 1.93       |               | ug/L |   | 98   | 70 - 130    |
| Endosulfan I (Alpha)             | 1.98        | 1.82       |               | ug/L |   | 92   | 70 - 130    |
| Endosulfan II (Beta)             | 1.98        | 2.17       |               | ug/L |   | 110  | 70 - 130    |
| Endosulfan sulfate               | 1.98        | 2.23       |               | ug/L |   | 113  | 70 - 130    |
| Endrin                           | 1.98        | 2.21       |               | ug/L |   | 112  | 70 - 130    |
| Endrin aldehyde                  | 1.98        | 1.92       |               | ug/L |   | 97   | 70 - 130    |
| EPTC                             | 1.98        | 1.96       |               | ug/L |   | 99   | 70 - 130    |
| Fluoranthene                     | 1.98        | 2.02       |               | ug/L |   | 102  | 70 - 130    |
| Fluorene                         | 1.98        | 1.97       |               | ug/L |   | 99   | 70 - 130    |
| gamma-Chlordane                  | 1.98        | 1.94       |               | ug/L |   | 98   | 70 - 130    |
| Heptachlor                       | 1.98        | 1.89       |               | ug/L |   | 95   | 70 - 130    |
| Heptachlor epoxide (isomer B)    | 1.98        | 1.97       |               | ug/L |   | 99   | 70 - 130    |

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# QC Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 525.2 - Semivolatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: LCS 380-27725/3-A**  
**Matrix: Water**  
**Analysis Batch: 28001**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 27725**

| Analyte                    | Spike<br>Added | LCS<br>Result | LCS<br>Qualifier | Unit | D | %Rec | %Rec<br>Limits |
|----------------------------|----------------|---------------|------------------|------|---|------|----------------|
| Hexachlorobenzene          | 1.98           | 1.85          |                  | ug/L |   | 94   | 70 - 130       |
| Hexachlorocyclopentadiene  | 1.98           | 1.99          |                  | ug/L |   | 100  | 70 - 130       |
| Indeno[1,2,3-cd]pyrene     | 1.98           | 2.20          |                  | ug/L |   | 111  | 70 - 130       |
| Isophorone                 | 1.98           | 2.05          |                  | ug/L |   | 104  | 70 - 130       |
| Lindane                    | 1.98           | 1.89          |                  | ug/L |   | 95   | 70 - 130       |
| Malathion                  | 1.98           | 2.17          |                  | ug/L |   | 110  | 70 - 130       |
| Methoxychlor               | 1.98           | 2.20          |                  | ug/L |   | 111  | 70 - 130       |
| Metolachlor                | 1.98           | 2.04          |                  | ug/L |   | 103  | 70 - 130       |
| Metribuzin                 | 1.98           | 2.13          |                  | ug/L |   | 107  | 70 - 130       |
| Molinate                   | 1.98           | 1.99          |                  | ug/L |   | 100  | 70 - 130       |
| Naphthalene                | 1.98           | 1.88          |                  | ug/L |   | 95   | 70 - 130       |
| Parathion                  | 1.98           | 2.20          |                  | ug/L |   | 111  | 70 - 130       |
| Pendimethalin (Penoxaline) | 1.98           | 2.00          |                  | ug/L |   | 101  | 70 - 130       |
| Phenanthrene               | 1.98           | 1.91          |                  | ug/L |   | 96   | 70 - 130       |
| Propachlor                 | 1.98           | 1.98          |                  | ug/L |   | 100  | 70 - 130       |
| Pyrene                     | 1.98           | 2.05          |                  | ug/L |   | 103  | 70 - 130       |
| Simazine                   | 1.98           | 2.06          |                  | ug/L |   | 104  | 70 - 130       |
| Terbacil                   | 1.98           | 2.17          |                  | ug/L |   | 109  | 70 - 130       |
| Terbutylazine              | 1.98           | 1.99          |                  | ug/L |   | 101  | 70 - 130       |
| Thiobencarb                | 1.98           | 2.08          |                  | ug/L |   | 105  | 70 - 130       |
| trans-Nonachlor            | 1.98           | 1.93          |                  | ug/L |   | 97   | 70 - 130       |
| Trifluralin                | 1.98           | 1.79          |                  | ug/L |   | 90   | 70 - 130       |

| Surrogate          | LCS<br>%Recovery | LCS<br>Qualifier | Limits   |
|--------------------|------------------|------------------|----------|
| 2-Nitro-m-xylene   | 98               |                  | 70 - 130 |
| Triphenylphosphate | 113              |                  | 70 - 130 |
| Perylene-d12       | 100              |                  | 70 - 130 |

**Lab Sample ID: LCSD 380-27725/4-A**  
**Matrix: Water**  
**Analysis Batch: 28001**

**Client Sample ID: Lab Control Sample Dup**  
**Prep Type: Total/NA**  
**Prep Batch: 27725**

| Analyte            | Spike<br>Added | LCSD<br>Result | LCSD<br>Qualifier | Unit | D | %Rec | %Rec<br>Limits | RPD | RPD<br>Limit |
|--------------------|----------------|----------------|-------------------|------|---|------|----------------|-----|--------------|
| 2,4'-DDD           | 1.99           | 1.92           |                   | ug/L |   | 97   | 70 - 130       | 2   | 20           |
| 2,4'-DDE           | 1.99           | 1.92           |                   | ug/L |   | 97   | 70 - 130       | 0   | 20           |
| 2,4'-DDT           | 1.99           | 2.06           |                   | ug/L |   | 104  | 70 - 130       | 1   | 20           |
| 2,4-Dinitrotoluene | 1.99           | 1.95           |                   | ug/L |   | 98   | 70 - 130       | 8   | 20           |
| 2,6-Dinitrotoluene | 1.99           | 2.01           |                   | ug/L |   | 101  | 70 - 130       | 4   | 20           |
| 4,4'-DDD           | 1.99           | 2.14           |                   | ug/L |   | 108  | 70 - 130       | 2   | 20           |
| 4,4'-DDE           | 1.99           | 2.00           |                   | ug/L |   | 101  | 70 - 130       | 3   | 20           |
| 4,4'-DDT           | 1.99           | 2.06           |                   | ug/L |   | 103  | 70 - 130       | 6   | 20           |
| Acenaphthene       | 1.99           | 2.03           |                   | ug/L |   | 102  | 70 - 130       | 5   | 20           |
| Acenaphthylene     | 1.99           | 1.99           |                   | ug/L |   | 100  | 70 - 130       | 5   | 20           |
| Acetochlor         | 1.99           | 2.14           |                   | ug/L |   | 108  | 70 - 130       | 2   | 20           |
| Alachlor           | 1.99           | 2.09           |                   | ug/L |   | 105  | 70 - 130       | 4   | 20           |
| alpha-BHC          | 1.99           | 2.01           |                   | ug/L |   | 101  | 70 - 130       | 3   | 20           |
| alpha-Chlordane    | 1.99           | 1.92           |                   | ug/L |   | 97   | 70 - 130       | 3   | 20           |
| Anthracene         | 1.99           | 1.95           |                   | ug/L |   | 98   | 70 - 130       | 4   | 20           |

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# QC Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 525.2 - Semivolatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: LCSD 380-27725/4-A**  
**Matrix: Water**  
**Analysis Batch: 28001**

**Client Sample ID: Lab Control Sample Dup**  
**Prep Type: Total/NA**  
**Prep Batch: 27725**

| Analyte                          | Spike Added | LCSD Result | LCSD Qualifier | Unit | D | %Rec | %Rec     |     | RPD | RPD Limit |
|----------------------------------|-------------|-------------|----------------|------|---|------|----------|-----|-----|-----------|
|                                  |             |             |                |      |   |      | Limits   | RPD |     |           |
| Atrazine                         | 1.99        | 1.98        |                | ug/L |   | 100  | 70 - 130 | 0   | 20  |           |
| Benz(a)anthracene                | 1.99        | 2.15        |                | ug/L |   | 108  | 70 - 130 | 2   | 20  |           |
| Benzo[a]pyrene                   | 1.99        | 2.26        |                | ug/L |   | 114  | 70 - 130 | 6   | 20  |           |
| Benzo[b]fluoranthene             | 1.99        | 2.26        |                | ug/L |   | 114  | 70 - 130 | 4   | 20  |           |
| Benzo[g,h,i]perylene             | 1.99        | 2.19        |                | ug/L |   | 110  | 70 - 130 | 7   | 20  |           |
| Benzo[k]fluoranthene             | 1.99        | 2.26        |                | ug/L |   | 114  | 70 - 130 | 7   | 20  |           |
| beta-BHC                         | 1.99        | 2.01        |                | ug/L |   | 101  | 70 - 130 | 0   | 20  |           |
| Bromacil                         | 1.99        | 2.25        |                | ug/L |   | 113  | 70 - 130 | 0   | 20  |           |
| Butachlor                        | 1.99        | 2.11        |                | ug/L |   | 106  | 70 - 130 | 2   | 20  |           |
| Butylbenzylphthalate             | 1.99        | 2.27        |                | ug/L |   | 114  | 70 - 130 | 0   | 20  |           |
| Caffeine                         | 1.99        | 1.63        |                | ug/L |   | 82   | 45 - 137 | 1   | 20  |           |
| Chlorobenzilate                  | 1.99        | 2.21        |                | ug/L |   | 112  | 70 - 130 | 0   | 20  |           |
| Chloroneb                        | 1.99        | 1.96        |                | ug/L |   | 98   | 70 - 130 | 4   | 20  |           |
| Chlorothalonil (Draconil, Bravo) | 1.99        | 1.94        |                | ug/L |   | 98   | 70 - 130 | 1   | 20  |           |
| Chlorpyrifos                     | 1.99        | 2.06        |                | ug/L |   | 104  | 70 - 130 | 2   | 20  |           |
| Chrysene                         | 1.99        | 2.05        |                | ug/L |   | 103  | 70 - 130 | 7   | 20  |           |
| delta-BHC                        | 1.99        | 2.10        |                | ug/L |   | 106  | 70 - 130 | 3   | 20  |           |
| Di(2-ethylhexyl)adipate          | 1.99        | 2.10        |                | ug/L |   | 106  | 70 - 130 | 7   | 20  |           |
| Bis(2-ethylhexyl) phthalate      | 1.99        | 1.84        |                | ug/L |   | 93   | 70 - 130 | 2   | 20  |           |
| Diazinon (Qualitative)           | 1.99        | 1.84        |                | ug/L |   | 92   | 15 - 132 | 3   | 20  |           |
| Dibenz(a,h)anthracene            | 1.99        | 2.23        |                | ug/L |   | 112  | 70 - 130 | 1   | 20  |           |
| Diclorvos (DDVP)                 | 1.99        | 2.37        |                | ug/L |   | 119  | 70 - 130 | 8   | 20  |           |
| Dieldrin                         | 1.99        | 2.07        |                | ug/L |   | 104  | 70 - 130 | 0   | 20  |           |
| Diethylphthalate                 | 1.99        | 2.05        |                | ug/L |   | 103  | 70 - 130 | 6   | 20  |           |
| Dimethoate                       | 1.99        | 1.39        |                | ug/L |   | 70   | 35 - 100 | 7   | 20  |           |
| Dimethylphthalate                | 1.99        | 2.08        |                | ug/L |   | 105  | 70 - 130 | 3   | 20  |           |
| Di-n-butyl phthalate             | 3.97        | 4.12        |                | ug/L |   | 104  | 70 - 130 | 5   | 20  |           |
| Di-n-octyl phthalate             | 1.99        | 1.82        |                | ug/L |   | 91   | 70 - 130 | 6   | 20  |           |
| Endosulfan I (Alpha)             | 1.99        | 1.88        |                | ug/L |   | 94   | 70 - 130 | 3   | 20  |           |
| Endosulfan II (Beta)             | 1.99        | 2.21        |                | ug/L |   | 111  | 70 - 130 | 2   | 20  |           |
| Endosulfan sulfate               | 1.99        | 2.23        |                | ug/L |   | 112  | 70 - 130 | 0   | 20  |           |
| Endrin                           | 1.99        | 2.18        |                | ug/L |   | 110  | 70 - 130 | 2   | 20  |           |
| Endrin aldehyde                  | 1.99        | 1.91        |                | ug/L |   | 96   | 70 - 130 | 0   | 20  |           |
| EPTC                             | 1.99        | 2.00        |                | ug/L |   | 101  | 70 - 130 | 2   | 20  |           |
| Fluoranthene                     | 1.99        | 2.02        |                | ug/L |   | 102  | 70 - 130 | 0   | 20  |           |
| Fluorene                         | 1.99        | 2.01        |                | ug/L |   | 101  | 70 - 130 | 2   | 20  |           |
| gamma-Chlordane                  | 1.99        | 1.95        |                | ug/L |   | 98   | 70 - 130 | 0   | 20  |           |
| Heptachlor                       | 1.99        | 2.00        |                | ug/L |   | 101  | 70 - 130 | 6   | 20  |           |
| Heptachlor epoxide (isomer B)    | 1.99        | 2.03        |                | ug/L |   | 102  | 70 - 130 | 3   | 20  |           |
| Hexachlorobenzene                | 1.99        | 1.88        |                | ug/L |   | 94   | 70 - 130 | 1   | 20  |           |
| Hexachlorocyclopentadiene        | 1.99        | 2.06        |                | ug/L |   | 104  | 70 - 130 | 3   | 20  |           |
| Indeno[1,2,3-cd]pyrene           | 1.99        | 2.28        |                | ug/L |   | 115  | 70 - 130 | 4   | 20  |           |
| Isophorone                       | 1.99        | 2.21        |                | ug/L |   | 111  | 70 - 130 | 7   | 20  |           |
| Lindane                          | 1.99        | 1.97        |                | ug/L |   | 99   | 70 - 130 | 4   | 20  |           |
| Malathion                        | 1.99        | 2.23        |                | ug/L |   | 112  | 70 - 130 | 3   | 20  |           |
| Methoxychlor                     | 1.99        | 2.32        |                | ug/L |   | 117  | 70 - 130 | 6   | 20  |           |
| Metolachlor                      | 1.99        | 2.09        |                | ug/L |   | 105  | 70 - 130 | 3   | 20  |           |
| Metribuzin                       | 1.99        | 2.15        |                | ug/L |   | 108  | 70 - 130 | 1   | 20  |           |
| Molinate                         | 1.99        | 1.99        |                | ug/L |   | 100  | 70 - 130 | 0   | 20  |           |

Eurofins Eaton Monrovia

# QC Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 525.2 - Semivolatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: LCSD 380-27725/4-A**  
**Matrix: Water**  
**Analysis Batch: 28001**

**Client Sample ID: Lab Control Sample Dup**  
**Prep Type: Total/NA**  
**Prep Batch: 27725**

| Analyte                    | Spike Added | LCSD Result | LCSD Qualifier | Unit | D | %Rec | %Rec Limits | RPD | RPD Limit |
|----------------------------|-------------|-------------|----------------|------|---|------|-------------|-----|-----------|
| Naphthalene                | 1.99        | 2.01        |                | ug/L |   | 101  | 70 - 130    | 7   | 20        |
| Parathion                  | 1.99        | 2.23        |                | ug/L |   | 112  | 70 - 130    | 1   | 20        |
| Pendimethalin (Penoxaline) | 1.99        | 2.08        |                | ug/L |   | 105  | 70 - 130    | 4   | 20        |
| Phenanthrene               | 1.99        | 1.97        |                | ug/L |   | 99   | 70 - 130    | 3   | 20        |
| Propachlor                 | 1.99        | 2.10        |                | ug/L |   | 106  | 70 - 130    | 6   | 20        |
| Pyrene                     | 1.99        | 2.03        |                | ug/L |   | 102  | 70 - 130    | 1   | 20        |
| Simazine                   | 1.99        | 2.09        |                | ug/L |   | 105  | 70 - 130    | 1   | 20        |
| Terbacil                   | 1.99        | 2.16        |                | ug/L |   | 109  | 70 - 130    | 0   | 20        |
| Terbutylazine              | 1.99        | 2.06        |                | ug/L |   | 104  | 70 - 130    | 3   | 20        |
| Thiobencarb                | 1.99        | 2.08        |                | ug/L |   | 105  | 70 - 130    | 0   | 20        |
| trans-Nonachlor            | 1.99        | 1.87        |                | ug/L |   | 94   | 70 - 130    | 3   | 20        |
| Trifluralin                | 1.99        | 1.88        |                | ug/L |   | 95   | 70 - 130    | 5   | 20        |

| Surrogate          | LCSD %Recovery | LCSD Qualifier | LCSD Limits |
|--------------------|----------------|----------------|-------------|
| 2-Nitro-m-xylene   | 101            |                | 70 - 130    |
| Triphenylphosphate | 108            |                | 70 - 130    |
| Perylene-d12       | 99             |                | 70 - 130    |

**Lab Sample ID: MRL 380-27725/2-A**  
**Matrix: Water**  
**Analysis Batch: 28001**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 27725**

| Analyte              | Spike Added | MRL Result | MRL Qualifier | Unit | D | %Rec | %Rec Limits |
|----------------------|-------------|------------|---------------|------|---|------|-------------|
| 2,4'-DDD             | 0.0993      | 0.147      |               | ug/L |   | 148  | 50 - 150    |
| 2,4'-DDE             | 0.0993      | 0.109      |               | ug/L |   | 109  | 50 - 150    |
| 2,4'-DDT             | 0.0993      | 0.112      |               | ug/L |   | 113  | 50 - 150    |
| 2,4-Dinitrotoluene   | 0.0993      | 0.0813     | J             | ug/L |   | 82   | 50 - 150    |
| 2,6-Dinitrotoluene   | 0.0993      | 0.0885     | J             | ug/L |   | 89   | 50 - 150    |
| 4,4'-DDD             | 0.0993      | 0.119      |               | ug/L |   | 120  | 50 - 150    |
| 4,4'-DDE             | 0.0993      | 0.103      |               | ug/L |   | 103  | 50 - 150    |
| 4,4'-DDT             | 0.0993      | 0.115      |               | ug/L |   | 116  | 50 - 150    |
| Acenaphthene         | 0.0993      | 0.0998     |               | ug/L |   | 100  | 50 - 150    |
| Acenaphthylene       | 0.0993      | 0.0900     | J             | ug/L |   | 91   | 50 - 150    |
| Acetochlor           | 0.0497      | 0.0531     | J             | ug/L |   | 107  | 50 - 150    |
| Alachlor             | 0.0497      | 0.0590     |               | ug/L |   | 119  | 50 - 150    |
| alpha-BHC            | 0.0993      | 0.0988     | J             | ug/L |   | 99   | 50 - 150    |
| alpha-Chlordane      | 0.0248      | 0.0297     | J             | ug/L |   | 120  | 50 - 150    |
| Anthracene           | 0.0199      | 0.0212     |               | ug/L |   | 107  | 50 - 150    |
| Atrazine             | 0.0497      | 0.0505     |               | ug/L |   | 102  | 50 - 150    |
| Benz(a)anthracene    | 0.0497      | 0.0526     |               | ug/L |   | 106  | 50 - 150    |
| Benzo[a]pyrene       | 0.0199      | 0.0182     | J             | ug/L |   | 91   | 50 - 150    |
| Benzo[b]fluoranthene | 0.0199      | 0.0240     |               | ug/L |   | 121  | 50 - 150    |
| Benzo[g,h,i]perylene | 0.0497      | 0.0496     | J             | ug/L |   | 100  | 50 - 150    |
| Benzo[k]fluoranthene | 0.0199      | 0.0218     |               | ug/L |   | 110  | 50 - 150    |
| beta-BHC             | 0.0993      | 0.104      |               | ug/L |   | 105  | 50 - 150    |
| Bromacil             | 0.0993      | 0.114      |               | ug/L |   | 115  | 50 - 150    |
| Butachlor            | 0.0497      | 0.0602     |               | ug/L |   | 121  | 50 - 150    |
| Butylbenzylphthalate | 0.149       | 0.214      | J             | ug/L |   | 143  | 50 - 150    |

Eurofins Eaton Monrovia

# QC Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 525.2 - Semivolatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: MRL 380-27725/2-A**  
**Matrix: Water**  
**Analysis Batch: 28001**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 27725**

| Analyte                          | Spike Added | MRL    | MRL       | Unit | D | %Rec | %Rec Limits |
|----------------------------------|-------------|--------|-----------|------|---|------|-------------|
|                                  |             | Result | Qualifier |      |   |      |             |
| Caffeine                         | 0.0497      | 0.0446 | J         | ug/L |   | 90   | 50 - 150    |
| Chlorobenzilate                  | 0.0993      | 0.110  |           | ug/L |   | 111  | 50 - 150    |
| Chloroneb                        | 0.0993      | 0.103  |           | ug/L |   | 104  | 50 - 150    |
| Chlorothalonil (Draconil, Bravo) | 0.0993      | 0.201  | ^3+       | ug/L |   | 202  | 50 - 150    |
| Chlorpyrifos                     | 0.0497      | 0.0550 |           | ug/L |   | 111  | 50 - 150    |
| Chrysene                         | 0.0199      | 0.0204 |           | ug/L |   | 103  | 50 - 150    |
| delta-BHC                        | 0.0993      | 0.125  |           | ug/L |   | 126  | 50 - 150    |
| Di(2-ethylhexyl)adipate          | 0.298       | 0.365  | J         | ug/L |   | 123  | 50 - 150    |
| Bis(2-ethylhexyl) phthalate      | 0.596       | 0.663  |           | ug/L |   | 111  | 50 - 150    |
| Diazinon (Qualitative)           | 0.0993      | 0.0892 | J         | ug/L |   | 90   | 15 - 132    |
| Dibenz(a,h)anthracene            | 0.0497      | 0.0535 |           | ug/L |   | 108  | 50 - 150    |
| Diclorvos (DDVP)                 | 0.0497      | 0.0600 |           | ug/L |   | 121  | 50 - 150    |
| Dieldrin                         | 0.0993      | 0.146  | J         | ug/L |   | 147  | 50 - 150    |
| Diethylphthalate                 | 0.149       | 0.138  | J         | ug/L |   | 93   | 50 - 150    |
| Dimethoate                       | 0.0993      | 0.0697 | J         | ug/L |   | 70   | 35 - 100    |
| Dimethylphthalate                | 0.298       | 0.299  | J         | ug/L |   | 100  | 50 - 150    |
| Di-n-butyl phthalate             | 0.298       | 0.412  | J         | ug/L |   | 138  | 49 - 243    |
| Di-n-octyl phthalate             | 0.0993      | 0.0945 | J         | ug/L |   | 95   | 50 - 150    |
| Endosulfan I (Alpha)             | 0.0993      | 0.108  |           | ug/L |   | 109  | 50 - 150    |
| Endosulfan II (Beta)             | 0.0993      | 0.142  |           | ug/L |   | 143  | 50 - 150    |
| Endosulfan sulfate               | 0.0993      | 0.123  |           | ug/L |   | 124  | 50 - 150    |
| Endrin                           | 0.0993      | 0.125  |           | ug/L |   | 126  | 50 - 150    |
| Endrin aldehyde                  | 0.0993      | 0.0998 |           | ug/L |   | 101  | 50 - 150    |
| EPTC                             | 0.0993      | 0.0969 | J         | ug/L |   | 98   | 50 - 150    |
| Fluoranthene                     | 0.0497      | 0.0550 | J         | ug/L |   | 111  | 50 - 150    |
| Fluorene                         | 0.0497      | ND     |           | ug/L |   | 99   | 50 - 150    |
| gamma-Chlordane                  | 0.0248      | 0.0325 | J         | ug/L |   | 131  | 50 - 150    |
| Heptachlor                       | 0.0397      | 0.0493 |           | ug/L |   | 124  | 50 - 150    |
| Heptachlor epoxide (isomer B)    | 0.0497      | 0.0632 |           | ug/L |   | 127  | 50 - 150    |
| Hexachlorobenzene                | 0.0497      | 0.0431 | J         | ug/L |   | 87   | 50 - 150    |
| Hexachlorocyclopentadiene        | 0.0497      | 0.0474 | J         | ug/L |   | 96   | 50 - 150    |
| Indeno[1,2,3-cd]pyrene           | 0.0497      | 0.0564 |           | ug/L |   | 114  | 50 - 150    |
| Isophorone                       | 0.0993      | 0.109  | J         | ug/L |   | 109  | 50 - 150    |
| Lindane                          | 0.0397      | 0.0445 |           | ug/L |   | 112  | 50 - 150    |
| Malathion                        | 0.0993      | 0.104  |           | ug/L |   | 104  | 50 - 150    |
| Methoxychlor                     | 0.0993      | 0.0969 | J         | ug/L |   | 98   | 50 - 150    |
| Metolachlor                      | 0.0497      | 0.0573 |           | ug/L |   | 115  | 50 - 150    |
| Metribuzin                       | 0.0497      | 0.0554 |           | ug/L |   | 112  | 50 - 150    |
| Molinate                         | 0.0993      | 0.102  |           | ug/L |   | 102  | 50 - 150    |
| Naphthalene                      | 0.0993      | 0.102  | J         | ug/L |   | 102  | 50 - 150    |
| Parathion                        | 0.0993      | 0.0939 | J         | ug/L |   | 95   | 50 - 150    |
| Pendimethalin (Penoxaline)       | 0.0993      | 0.0815 | J         | ug/L |   | 82   | 50 - 150    |
| Phenanthrene                     | 0.0199      | 0.0226 | J         | ug/L |   | 114  | 50 - 150    |
| Propachlor                       | 0.0497      | 0.0421 | J         | ug/L |   | 85   | 50 - 150    |
| Pyrene                           | 0.0497      | 0.0550 |           | ug/L |   | 111  | 50 - 150    |
| Simazine                         | 0.0497      | 0.0570 |           | ug/L |   | 115  | 50 - 150    |
| Terbacil                         | 0.0993      | 0.126  |           | ug/L |   | 126  | 50 - 150    |
| Terbutylazine                    | 0.0993      | 0.0952 | J         | ug/L |   | 96   | 50 - 150    |
| Thiobencarb                      | 0.0993      | 0.110  | J         | ug/L |   | 110  | 50 - 150    |

Eurofins Eaton Monrovia

# QC Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 525.2 - Semivolatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: MRL 380-27725/2-A**  
**Matrix: Water**  
**Analysis Batch: 28001**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 27725**

| Analyte         | Spike Added | MRL Result | MRL Qualifier | Unit | D | %Rec | %Rec Limits |
|-----------------|-------------|------------|---------------|------|---|------|-------------|
| trans-Nonachlor | 0.0248      | 0.0309     | J             | ug/L |   | 124  | 50 - 150    |
| Trifluralin     | 0.0993      | 0.0692     | J             | ug/L |   | 70   | 50 - 150    |

| Surrogate          | MRL %Recovery | MRL Qualifier | Limits   |
|--------------------|---------------|---------------|----------|
| 2-Nitro-m-xylene   | 100           |               | 70 - 130 |
| Triphenylphosphate | 116           |               | 70 - 130 |
| Perylene-d12       | 98            |               | 70 - 130 |

**Lab Sample ID: 380-32064-C-1-A MS**  
**Matrix: Water**  
**Analysis Batch: 28001**

**Client Sample ID: Matrix Spike**  
**Prep Type: Total/NA**  
**Prep Batch: 27725**

| Analyte                          | Sample Result | Sample Qualifier | Spike Added | MS Result | MS Qualifier | Unit | D | %Rec | %Rec Limits |
|----------------------------------|---------------|------------------|-------------|-----------|--------------|------|---|------|-------------|
| 2,4'-DDD                         | ND            |                  | 2.03        | 1.94      |              | ug/L |   | 96   | 70 - 130    |
| 2,4'-DDE                         | ND            |                  | 2.03        | 1.94      |              | ug/L |   | 96   | 70 - 130    |
| 2,4'-DDT                         | ND            |                  | 2.03        | 2.10      |              | ug/L |   | 104  | 70 - 130    |
| 2,4-Dinitrotoluene               | ND            |                  | 2.03        | 2.10      |              | ug/L |   | 104  | 70 - 130    |
| 2,6-Dinitrotoluene               | ND            |                  | 2.03        | 2.10      |              | ug/L |   | 104  | 70 - 130    |
| 4,4'-DDD                         | ND            |                  | 2.03        | 2.16      |              | ug/L |   | 107  | 70 - 130    |
| 4,4'-DDE                         | ND            |                  | 2.03        | 2.02      |              | ug/L |   | 100  | 70 - 130    |
| 4,4'-DDT                         | ND            |                  | 2.03        | 2.08      |              | ug/L |   | 103  | 70 - 130    |
| Acenaphthene                     | ND            |                  | 2.03        | 2.07      |              | ug/L |   | 102  | 70 - 130    |
| Acenaphthylene                   | ND            |                  | 2.03        | 2.10      |              | ug/L |   | 103  | 70 - 130    |
| Acetochlor                       | ND            |                  | 2.03        | 2.14      |              | ug/L |   | 106  | 70 - 130    |
| Alachlor                         | ND            |                  | 2.03        | 2.08      |              | ug/L |   | 103  | 70 - 130    |
| alpha-BHC                        | ND            |                  | 2.03        | 2.12      |              | ug/L |   | 105  | 70 - 130    |
| alpha-Chlordane                  | ND            |                  | 2.03        | 1.88      |              | ug/L |   | 93   | 70 - 130    |
| Anthracene                       | ND            |                  | 2.03        | 1.54      |              | ug/L |   | 76   | 70 - 130    |
| Atrazine                         | ND            |                  | 2.03        | 2.09      |              | ug/L |   | 103  | 70 - 130    |
| Benz(a)anthracene                | ND            |                  | 2.03        | 2.05      |              | ug/L |   | 101  | 70 - 130    |
| Benzo[a]pyrene                   | ND            |                  | 2.03        | 1.99      |              | ug/L |   | 98   | 70 - 130    |
| Benzo[b]fluoranthene             | ND            |                  | 2.03        | 2.27      |              | ug/L |   | 112  | 70 - 130    |
| Benzo[g,h,i]perylene             | ND            |                  | 2.03        | 2.19      |              | ug/L |   | 108  | 70 - 130    |
| Benzo[k]fluoranthene             | ND            |                  | 2.03        | 2.20      |              | ug/L |   | 109  | 70 - 130    |
| beta-BHC                         | ND            |                  | 2.03        | 2.14      |              | ug/L |   | 106  | 70 - 130    |
| Bromacil                         | ND            |                  | 2.03        | 2.32      |              | ug/L |   | 115  | 70 - 130    |
| Butachlor                        | ND            |                  | 2.03        | 2.26      |              | ug/L |   | 112  | 70 - 130    |
| Butylbenzylphthalate             | ND            |                  | 2.03        | 2.26      |              | ug/L |   | 111  | 70 - 130    |
| Caffeine                         | ND            |                  | 2.03        | 1.64      |              | ug/L |   | 81   | 46 - 144    |
| Chlorobenzilate                  | ND            |                  | 2.03        | 2.25      |              | ug/L |   | 111  | 70 - 130    |
| Chloroneb                        | ND            |                  | 2.03        | 2.00      |              | ug/L |   | 99   | 70 - 130    |
| Chlorothalonil (Draconil, Bravo) | ND            | ^3+              | 2.03        | 1.92      |              | ug/L |   | 95   | 70 - 130    |
| Chlorpyrifos                     | ND            |                  | 2.03        | 2.10      |              | ug/L |   | 104  | 70 - 130    |
| Chrysene                         | ND            |                  | 2.03        | 2.05      |              | ug/L |   | 101  | 70 - 130    |
| delta-BHC                        | ND            |                  | 2.03        | 2.06      |              | ug/L |   | 102  | 70 - 130    |
| Di(2-ethylhexyl)adipate          | ND            |                  | 2.03        | 2.21      |              | ug/L |   | 109  | 70 - 130    |
| Bis(2-ethylhexyl) phthalate      | ND            |                  | 2.03        | 2.17      |              | ug/L |   | 107  | 70 - 130    |
| Diazinon (Qualitative)           | ND            |                  | 2.03        | 1.95      |              | ug/L |   | 96   | 15 - 132    |

# QC Sample Results

Client: City & County of Honolulu  
 Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 525.2 - Semivolatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: 380-32064-C-1-A MS**  
**Matrix: Water**  
**Analysis Batch: 28001**

**Client Sample ID: Matrix Spike**  
**Prep Type: Total/NA**  
**Prep Batch: 27725**

| Analyte                       | Sample | Sample           | Spike            | MS            | MS        | Unit | D | %Rec | %Rec<br>Limits |
|-------------------------------|--------|------------------|------------------|---------------|-----------|------|---|------|----------------|
|                               | Result | Qualifier        | Added            | Result        | Qualifier |      |   |      |                |
| Dibenz(a,h)anthracene         | ND     |                  | 2.03             | 2.40          |           | ug/L |   | 118  | 70 - 130       |
| Diclorvos (DDVP)              | ND     |                  | 2.03             | 2.45          |           | ug/L |   | 121  | 70 - 130       |
| Dieldrin                      | ND     |                  | 2.03             | 2.22          |           | ug/L |   | 110  | 70 - 130       |
| Diethylphthalate              | ND     |                  | 2.03             | 2.14          |           | ug/L |   | 105  | 70 - 130       |
| Dimethoate                    | ND     |                  | 2.03             | 1.45          |           | ug/L |   | 72   | 34 - 111       |
| Dimethylphthalate             | ND     |                  | 2.03             | 2.14          |           | ug/L |   | 106  | 70 - 130       |
| Di-n-butyl phthalate          | ND     |                  | 4.05             | 4.13          |           | ug/L |   | 102  | 70 - 130       |
| Di-n-octyl phthalate          | ND     |                  | 2.03             | 2.21          |           | ug/L |   | 109  | 70 - 130       |
| Endosulfan I (Alpha)          | ND     |                  | 2.03             | 1.85          |           | ug/L |   | 91   | 70 - 130       |
| Endosulfan II (Beta)          | ND     |                  | 2.03             | 2.18          |           | ug/L |   | 108  | 70 - 130       |
| Endosulfan sulfate            | ND     |                  | 2.03             | 2.28          |           | ug/L |   | 113  | 70 - 130       |
| Endrin                        | ND     |                  | 2.03             | 2.21          |           | ug/L |   | 109  | 70 - 130       |
| Endrin aldehyde               | ND     | F1               | 2.03             | 1.31          | F1        | ug/L |   | 65   | 70 - 130       |
| EPTC                          | ND     |                  | 2.03             | 2.29          |           | ug/L |   | 113  | 70 - 130       |
| Fluoranthene                  | ND     |                  | 2.03             | 2.08          |           | ug/L |   | 103  | 70 - 130       |
| Fluorene                      | ND     |                  | 2.03             | 2.14          |           | ug/L |   | 106  | 70 - 130       |
| gamma-Chlordane               | ND     |                  | 2.03             | 1.90          |           | ug/L |   | 94   | 70 - 130       |
| Heptachlor                    | ND     |                  | 2.03             | 1.99          |           | ug/L |   | 98   | 70 - 130       |
| Heptachlor epoxide (isomer B) | ND     |                  | 2.03             | 2.03          |           | ug/L |   | 100  | 70 - 130       |
| Hexachlorobenzene             | ND     |                  | 2.03             | 1.96          |           | ug/L |   | 97   | 70 - 130       |
| Hexachlorocyclopentadiene     | ND     |                  | 2.03             | 2.10          |           | ug/L |   | 104  | 70 - 130       |
| Indeno[1,2,3-cd]pyrene        | ND     |                  | 2.03             | 2.34          |           | ug/L |   | 116  | 70 - 130       |
| Isophorone                    | ND     |                  | 2.03             | 2.25          |           | ug/L |   | 111  | 70 - 130       |
| Lindane                       | ND     |                  | 2.03             | 2.02          |           | ug/L |   | 100  | 70 - 130       |
| Malathion                     | ND     |                  | 2.03             | 2.33          |           | ug/L |   | 115  | 70 - 130       |
| Methoxychlor                  | ND     |                  | 2.03             | 2.43          |           | ug/L |   | 120  | 70 - 130       |
| Metolachlor                   | ND     |                  | 2.03             | 2.16          |           | ug/L |   | 106  | 70 - 130       |
| Metribuzin                    | ND     |                  | 2.03             | 2.35          |           | ug/L |   | 116  | 70 - 130       |
| Molinate                      | ND     |                  | 2.03             | 2.28          |           | ug/L |   | 112  | 70 - 130       |
| Naphthalene                   | ND     |                  | 2.03             | 2.05          |           | ug/L |   | 101  | 70 - 130       |
| Parathion                     | ND     |                  | 2.03             | 2.37          |           | ug/L |   | 117  | 70 - 130       |
| Pendimethalin (Penoxaline)    | ND     |                  | 2.03             | 2.15          |           | ug/L |   | 106  | 70 - 130       |
| Phenanthrene                  | ND     |                  | 2.03             | 2.04          |           | ug/L |   | 101  | 70 - 130       |
| Propachlor                    | ND     |                  | 2.03             | 2.15          |           | ug/L |   | 106  | 70 - 130       |
| Pyrene                        | ND     |                  | 2.03             | 2.09          |           | ug/L |   | 103  | 70 - 130       |
| Simazine                      | ND     |                  | 2.03             | 2.20          |           | ug/L |   | 109  | 70 - 130       |
| Terbacil                      | ND     |                  | 2.03             | 2.22          |           | ug/L |   | 109  | 70 - 130       |
| Terbutylazine                 | ND     |                  | 2.03             | 2.15          |           | ug/L |   | 106  | 70 - 130       |
| Thiobencarb                   | ND     |                  | 2.03             | 2.24          |           | ug/L |   | 110  | 70 - 130       |
| trans-Nonachlor               | ND     |                  | 2.03             | 1.90          |           | ug/L |   | 94   | 70 - 130       |
| Trifluralin                   | ND     |                  | 2.03             | 1.96          |           | ug/L |   | 97   | 70 - 130       |
|                               |        | <b>MS MS</b>     |                  |               |           |      |   |      |                |
| <b>Surrogate</b>              |        | <b>%Recovery</b> | <b>Qualifier</b> | <b>Limits</b> |           |      |   |      |                |
| 2-Nitro-m-xylene              |        | 101              |                  | 70 - 130      |           |      |   |      |                |
| Triphenylphosphate            |        | 106              |                  | 70 - 130      |           |      |   |      |                |
| Perylene-d12                  |        | 101              |                  | 70 - 130      |           |      |   |      |                |

# QC Sample Results

Client: City & County of Honolulu  
 Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 525.2 - Semivolatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: 380-32064-I-1-A MSD**

**Matrix: Water**

**Analysis Batch: 28001**

**Client Sample ID: Matrix Spike Duplicate**

**Prep Type: Total/NA**

**Prep Batch: 27725**

| Analyte                          | Sample | Sample    | Spike | MSD    | MSD       | Unit | D | %Rec | %Rec     | RPD | RPD   |
|----------------------------------|--------|-----------|-------|--------|-----------|------|---|------|----------|-----|-------|
|                                  | Result | Qualifier | Added | Result | Qualifier |      |   |      | Limits   |     | Limit |
| 2,4'-DDD                         | ND     |           | 2.03  | 1.98   |           | ug/L |   | 98   | 70 - 130 | 2   | 20    |
| 2,4'-DDE                         | ND     |           | 2.03  | 1.99   |           | ug/L |   | 98   | 70 - 130 | 3   | 20    |
| 2,4'-DDT                         | ND     |           | 2.03  | 2.12   |           | ug/L |   | 104  | 70 - 130 | 1   | 20    |
| 2,4-Dinitrotoluene               | ND     |           | 2.03  | 2.08   |           | ug/L |   | 102  | 70 - 130 | 1   | 20    |
| 2,6-Dinitrotoluene               | ND     |           | 2.03  | 2.14   |           | ug/L |   | 105  | 70 - 130 | 2   | 20    |
| 4,4'-DDD                         | ND     |           | 2.03  | 2.22   |           | ug/L |   | 110  | 70 - 130 | 3   | 20    |
| 4,4'-DDE                         | ND     |           | 2.03  | 2.11   |           | ug/L |   | 104  | 70 - 130 | 5   | 20    |
| 4,4'-DDT                         | ND     |           | 2.03  | 2.18   |           | ug/L |   | 108  | 70 - 130 | 5   | 20    |
| Acenaphthene                     | ND     |           | 2.03  | 2.06   |           | ug/L |   | 102  | 70 - 130 | 0   | 20    |
| Acenaphthylene                   | ND     |           | 2.03  | 2.10   |           | ug/L |   | 104  | 70 - 130 | 0   | 20    |
| Acetochlor                       | ND     |           | 2.03  | 2.10   |           | ug/L |   | 104  | 70 - 130 | 2   | 20    |
| Alachlor                         | ND     |           | 2.03  | 2.11   |           | ug/L |   | 104  | 70 - 130 | 1   | 20    |
| alpha-BHC                        | ND     |           | 2.03  | 2.11   |           | ug/L |   | 104  | 70 - 130 | 1   | 20    |
| alpha-Chlordane                  | ND     |           | 2.03  | 1.94   |           | ug/L |   | 96   | 70 - 130 | 3   | 20    |
| Anthracene                       | ND     |           | 2.03  | 1.55   |           | ug/L |   | 76   | 70 - 130 | 1   | 20    |
| Atrazine                         | ND     |           | 2.03  | 2.03   |           | ug/L |   | 100  | 70 - 130 | 3   | 20    |
| Benz(a)anthracene                | ND     |           | 2.03  | 2.13   |           | ug/L |   | 105  | 70 - 130 | 4   | 20    |
| Benzo[a]pyrene                   | ND     |           | 2.03  | 2.07   |           | ug/L |   | 102  | 70 - 130 | 4   | 20    |
| Benzo[b]fluoranthene             | ND     |           | 2.03  | 2.28   |           | ug/L |   | 113  | 70 - 130 | 0   | 20    |
| Benzo[g,h,i]perylene             | ND     |           | 2.03  | 2.25   |           | ug/L |   | 111  | 70 - 130 | 3   | 20    |
| Benzo[k]fluoranthene             | ND     |           | 2.03  | 2.29   |           | ug/L |   | 113  | 70 - 130 | 4   | 20    |
| beta-BHC                         | ND     |           | 2.03  | 2.03   |           | ug/L |   | 100  | 70 - 130 | 5   | 20    |
| Bromacil                         | ND     |           | 2.03  | 2.31   |           | ug/L |   | 114  | 70 - 130 | 0   | 20    |
| Butachlor                        | ND     |           | 2.03  | 2.21   |           | ug/L |   | 109  | 70 - 130 | 2   | 20    |
| Butylbenzylphthalate             | ND     |           | 2.03  | 2.33   |           | ug/L |   | 115  | 70 - 130 | 3   | 20    |
| Caffeine                         | ND     |           | 2.03  | 1.67   |           | ug/L |   | 83   | 46 - 144 | 2   | 20    |
| Chlorobenzilate                  | ND     |           | 2.03  | 2.27   |           | ug/L |   | 112  | 70 - 130 | 1   | 20    |
| Chloroneb                        | ND     |           | 2.03  | 2.04   |           | ug/L |   | 100  | 70 - 130 | 2   | 20    |
| Chlorothalonil (Draconil, Bravo) | ND     | ^3+       | 2.03  | 1.97   |           | ug/L |   | 97   | 70 - 130 | 2   | 20    |
| Chlorpyrifos                     | ND     |           | 2.03  | 2.06   |           | ug/L |   | 102  | 70 - 130 | 2   | 20    |
| Chrysene                         | ND     |           | 2.03  | 2.05   |           | ug/L |   | 101  | 70 - 130 | 0   | 20    |
| delta-BHC                        | ND     |           | 2.03  | 2.12   |           | ug/L |   | 105  | 70 - 130 | 3   | 20    |
| Di(2-ethylhexyl)adipate          | ND     |           | 2.03  | 2.33   |           | ug/L |   | 115  | 70 - 130 | 5   | 20    |
| Bis(2-ethylhexyl) phthalate      | ND     |           | 2.03  | 2.10   |           | ug/L |   | 104  | 70 - 130 | 3   | 20    |
| Diazinon (Qualitative)           | ND     |           | 2.03  | 2.06   |           | ug/L |   | 101  | 15 - 132 | 5   | 20    |
| Dibenz(a,h)anthracene            | ND     |           | 2.03  | 2.44   |           | ug/L |   | 120  | 70 - 130 | 2   | 20    |
| Diclorvos (DDVP)                 | ND     |           | 2.03  | 2.40   |           | ug/L |   | 118  | 70 - 130 | 2   | 20    |
| Dieldrin                         | ND     |           | 2.03  | 2.23   |           | ug/L |   | 110  | 70 - 130 | 0   | 20    |
| Diethylphthalate                 | ND     |           | 2.03  | 2.10   |           | ug/L |   | 103  | 70 - 130 | 2   | 20    |
| Dimethoate                       | ND     |           | 2.03  | 1.53   |           | ug/L |   | 76   | 34 - 111 | 5   | 20    |
| Dimethylphthalate                | ND     |           | 2.03  | 2.15   |           | ug/L |   | 106  | 70 - 130 | 0   | 20    |
| Di-n-butyl phthalate             | ND     |           | 4.05  | 4.15   |           | ug/L |   | 102  | 70 - 130 | 1   | 20    |
| Di-n-octyl phthalate             | ND     |           | 2.03  | 2.23   |           | ug/L |   | 110  | 70 - 130 | 1   | 20    |
| Endosulfan I (Alpha)             | ND     |           | 2.03  | 1.92   |           | ug/L |   | 95   | 70 - 130 | 4   | 20    |
| Endosulfan II (Beta)             | ND     |           | 2.03  | 2.22   |           | ug/L |   | 110  | 70 - 130 | 2   | 20    |
| Endosulfan sulfate               | ND     |           | 2.03  | 2.36   |           | ug/L |   | 117  | 70 - 130 | 4   | 20    |
| Endrin                           | ND     |           | 2.03  | 2.25   |           | ug/L |   | 111  | 70 - 130 | 2   | 20    |
| Endrin aldehyde                  | ND     | F1        | 2.03  | 1.17   | F1        | ug/L |   | 58   | 70 - 130 | 11  | 20    |

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# QC Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 525.2 - Semivolatile Organic Compounds (GC/MS) (Continued)

**Lab Sample ID: 380-32064-I-1-A MSD**  
**Matrix: Water**  
**Analysis Batch: 28001**

**Client Sample ID: Matrix Spike Duplicate**  
**Prep Type: Total/NA**  
**Prep Batch: 27725**

| Analyte                       | Sample | Sample           | Spike            | MSD    | MSD           | Unit | D | %Rec | %Rec     | RPD | Limit |
|-------------------------------|--------|------------------|------------------|--------|---------------|------|---|------|----------|-----|-------|
|                               | Result | Qualifier        | Added            | Result | Qualifier     |      |   |      | Limits   |     |       |
| EPTC                          | ND     |                  | 2.03             | 2.29   |               | ug/L |   | 113  | 70 - 130 | 0   | 20    |
| Fluoranthene                  | ND     |                  | 2.03             | 2.08   |               | ug/L |   | 102  | 70 - 130 | 0   | 20    |
| Fluorene                      | ND     |                  | 2.03             | 2.10   |               | ug/L |   | 104  | 70 - 130 | 2   | 20    |
| gamma-Chlordane               | ND     |                  | 2.03             | 1.98   |               | ug/L |   | 98   | 70 - 130 | 4   | 20    |
| Heptachlor                    | ND     |                  | 2.03             | 2.01   |               | ug/L |   | 99   | 70 - 130 | 1   | 20    |
| Heptachlor epoxide (isomer B) | ND     |                  | 2.03             | 2.06   |               | ug/L |   | 102  | 70 - 130 | 1   | 20    |
| Hexachlorobenzene             | ND     |                  | 2.03             | 2.00   |               | ug/L |   | 98   | 70 - 130 | 2   | 20    |
| Hexachlorocyclopentadiene     | ND     |                  | 2.03             | 2.08   |               | ug/L |   | 103  | 70 - 130 | 1   | 20    |
| Indeno[1,2,3-cd]pyrene        | ND     |                  | 2.03             | 2.42   |               | ug/L |   | 119  | 70 - 130 | 3   | 20    |
| Isophorone                    | ND     |                  | 2.03             | 2.21   |               | ug/L |   | 109  | 70 - 130 | 2   | 20    |
| Lindane                       | ND     |                  | 2.03             | 2.01   |               | ug/L |   | 99   | 70 - 130 | 0   | 20    |
| Malathion                     | ND     |                  | 2.03             | 2.24   |               | ug/L |   | 111  | 70 - 130 | 4   | 20    |
| Methoxychlor                  | ND     |                  | 2.03             | 2.34   |               | ug/L |   | 115  | 70 - 130 | 4   | 20    |
| Metolachlor                   | ND     |                  | 2.03             | 2.12   |               | ug/L |   | 105  | 70 - 130 | 2   | 20    |
| Metribuzin                    | ND     |                  | 2.03             | 2.42   |               | ug/L |   | 119  | 70 - 130 | 3   | 20    |
| Molinate                      | ND     |                  | 2.03             | 2.34   |               | ug/L |   | 116  | 70 - 130 | 3   | 20    |
| Naphthalene                   | ND     |                  | 2.03             | 1.98   |               | ug/L |   | 98   | 70 - 130 | 3   | 20    |
| Parathion                     | ND     |                  | 2.03             | 2.34   |               | ug/L |   | 116  | 70 - 130 | 1   | 20    |
| Pendimethalin (Penoxaline)    | ND     |                  | 2.03             | 2.10   |               | ug/L |   | 104  | 70 - 130 | 2   | 20    |
| Phenanthrene                  | ND     |                  | 2.03             | 2.01   |               | ug/L |   | 99   | 70 - 130 | 2   | 20    |
| Propachlor                    | ND     |                  | 2.03             | 2.19   |               | ug/L |   | 108  | 70 - 130 | 2   | 20    |
| Pyrene                        | ND     |                  | 2.03             | 2.07   |               | ug/L |   | 102  | 70 - 130 | 1   | 20    |
| Simazine                      | ND     |                  | 2.03             | 2.21   |               | ug/L |   | 109  | 70 - 130 | 0   | 20    |
| Terbacil                      | ND     |                  | 2.03             | 2.19   |               | ug/L |   | 108  | 70 - 130 | 1   | 20    |
| Terbutylazine                 | ND     |                  | 2.03             | 2.11   |               | ug/L |   | 104  | 70 - 130 | 2   | 20    |
| Thiobencarb                   | ND     |                  | 2.03             | 2.22   |               | ug/L |   | 109  | 70 - 130 | 1   | 20    |
| trans-Nonachlor               | ND     |                  | 2.03             | 1.93   |               | ug/L |   | 95   | 70 - 130 | 1   | 20    |
| Trifluralin                   | ND     |                  | 2.03             | 2.01   |               | ug/L |   | 99   | 70 - 130 | 3   | 20    |
|                               |        | <b>MSD</b>       | <b>MSD</b>       |        |               |      |   |      |          |     |       |
| <b>Surrogate</b>              |        | <b>%Recovery</b> | <b>Qualifier</b> |        | <b>Limits</b> |      |   |      |          |     |       |
| 2-Nitro-m-xylene              |        | 98               |                  |        | 70 - 130      |      |   |      |          |     |       |
| Triphenylphosphate            |        | 108              |                  |        | 70 - 130      |      |   |      |          |     |       |
| Perylene-d12                  |        | 102              |                  |        | 70 - 130      |      |   |      |          |     |       |

## Method: 625 PAH Physis LL (EAL) + TICs - EPA 625 Base/Neutral and Acid Organics i

**Lab Sample ID: 102964-B1**  
**Matrix: BlankMatrix**  
**Analysis Batch: O-40074**

**Client Sample ID: Method Blank**  
**Prep Type: Total/NA**  
**Prep Batch: O-40074\_P**

| Analyte                    | Blank  | Blank     | RL    | MDL   | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------------------------|--------|-----------|-------|-------|------|---|----------------|----------------|---------|
|                            | Result | Qualifier |       |       |      |   |                |                |         |
| 1-Methylnaphthalene        | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| 1-Methylphenanthrene       | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| 2,3,5-Trimethylnaphthalene | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| 2,6-Dimethylnaphthalene    | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| 2-Methylnaphthalene        | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Acenaphthene               | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Acenaphthylene             | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Anthracene                 | ND     |           | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |

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# QC Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 625 PAH Physis LL (EAL) + TICs - EPA 625 Base/Neutral and Acid Organics i (Continued)

**Lab Sample ID: 102964-B1**  
**Matrix: BlankMatrix**  
**Analysis Batch: O-40074**

**Client Sample ID: Method Blank**  
**Prep Type: Total/NA**  
**Prep Batch: O-40074\_P**

| Analyte                      | Blank Result | Blank Qualifier | RL    | MDL   | Unit | D | Prepared       | Analyzed       | Dil Fac |
|------------------------------|--------------|-----------------|-------|-------|------|---|----------------|----------------|---------|
| Benz[a]anthracene            | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Benzo[a]pyrene               | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Benzo[b]fluoranthene         | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Benzo[e]pyrene               | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Benzo[g,h,i]perylene         | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Benzo[k]fluoranthene         | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Biphenyl                     | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Chrysene                     | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Dibenz[a,h]anthracene        | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Dibenzo[a,l]pyrene           | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Dibenzothiophene             | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Disalicylidenepropanediamine | ND           |                 | 0.1   | 0.05  | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Fluoranthene                 | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Fluorene                     | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Indeno[1,2,3-cd]pyrene       | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Naphthalene                  | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Perylene                     | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Phenanthrene                 | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| Pyrene                       | ND           |                 | 0.005 | 0.001 | µg/L |   | 12/26/22 00:00 | 01/08/23 02:25 | 1       |

| Surrogate          | Blank %Recovery | Blank Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|--------------------|-----------------|-----------------|----------|----------------|----------------|---------|
| (d10-Acenaphthene) | 92              |                 | 27 - 133 | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| (d10-Phenanthrene) | 94              |                 | 43 - 129 | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| (d12-Chrysene)     | 99              |                 | 52 - 144 | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| (d12-Perylene)     | 103             |                 | 36 - 161 | 12/26/22 00:00 | 01/08/23 02:25 | 1       |
| (d8-Naphthalene)   | 87              |                 | 25 - 125 | 12/26/22 00:00 | 01/08/23 02:25 | 1       |

**Lab Sample ID: 102964-BS1**  
**Matrix: BlankMatrix**  
**Analysis Batch: O-40074**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: O-40074\_P**

| Analyte                    | Spike Added | LCS Result | LCS Qualifier | Unit | D | %Rec | %Rec Limits |
|----------------------------|-------------|------------|---------------|------|---|------|-------------|
| 1-Methylnaphthalene        | 0.5         | 0.477      |               | µg/L |   | 95   | 31 - 128    |
| 1-Methylphenanthrene       | 0.5         | 0.448      |               | µg/L |   | 90   | 66 - 127    |
| 2,3,5-Trimethylnaphthalene | 0.5         | 0.498      |               | µg/L |   | 100  | 55 - 122    |
| 2,6-Dimethylnaphthalene    | 0.5         | 0.493      |               | µg/L |   | 99   | 48 - 120    |
| 2-Methylnaphthalene        | 1.5         | 1.44       |               | µg/L |   | 96   | 47 - 130    |
| Acenaphthene               | 1.5         | 1.49       |               | µg/L |   | 99   | 53 - 131    |
| Acenaphthylene             | 1.5         | 1.52       |               | µg/L |   | 101  | 43 - 140    |
| Anthracene                 | 1.5         | 1.52       |               | µg/L |   | 101  | 58 - 135    |
| Benz[a]anthracene          | 1.5         | 1.56       |               | µg/L |   | 104  | 55 - 145    |
| Benzo[a]pyrene             | 1.5         | 1.54       |               | µg/L |   | 103  | 51 - 143    |
| Benzo[b]fluoranthene       | 1.5         | 1.63       |               | µg/L |   | 109  | 46 - 165    |
| Benzo[e]pyrene             | 0.5         | 0.502      |               | µg/L |   | 100  | 42 - 152    |
| Benzo[g,h,i]perylene       | 1.5         | 1.59       |               | µg/L |   | 106  | 63 - 133    |
| Benzo[k]fluoranthene       | 1.5         | 1.64       |               | µg/L |   | 109  | 56 - 145    |
| Biphenyl                   | 0.5         | 0.523      |               | µg/L |   | 105  | 56 - 119    |
| Chrysene                   | 1.5         | 1.55       |               | µg/L |   | 103  | 56 - 141    |

Eurofins Eaton Monrovia

# QC Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 625 PAH Physis LL (EAL) + TICs - EPA 625 Base/Neutral and Acid Organics i (Continued)

**Lab Sample ID: 102964-BS1**  
**Matrix: BlankMatrix**  
**Analysis Batch: O-40074**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: O-40074\_P**

| Analyte                      | Spike Added | LCS Result | LCS Qualifier | Unit | D | %Rec | %Rec Limits |
|------------------------------|-------------|------------|---------------|------|---|------|-------------|
| Dibenz[a,h]anthracene        | 1.5         | 1.55       |               | µg/L |   | 103  | 55 - 150    |
| Dibenzo[a,l]pyrene           | 0.5         | 0.539      |               | µg/L |   | 108  | 50 - 150    |
| Dibenzothiophene             | 0.5         | 0.459      |               | µg/L |   | 92   | 46 - 126    |
| Disalicylideneprapanediamine | 50          | 42.9       |               | µg/L |   | 86   | 50 - 150    |
| Fluoranthene                 | 1.5         | 1.68       |               | µg/L |   | 112  | 60 - 146    |
| Fluorene                     | 1.5         | 1.58       |               | µg/L |   | 105  | 58 - 131    |
| Indeno[1,2,3-cd]pyrene       | 1.5         | 1.61       |               | µg/L |   | 107  | 50 - 151    |
| Naphthalene                  | 1.5         | 1.38       |               | µg/L |   | 92   | 41 - 126    |
| Perylene                     | 0.5         | 0.478      |               | µg/L |   | 96   | 48 - 141    |
| Phenanthrene                 | 1.5         | 1.57       |               | µg/L |   | 105  | 67 - 127    |
| Pyrene                       | 1.5         | 1.62       |               | µg/L |   | 108  | 54 - 156    |

| Surrogate          | LCS %Recovery | LCS Qualifier | Limits   |
|--------------------|---------------|---------------|----------|
| (d10-Acenaphthene) | 93            |               | 27 - 133 |
| (d10-Phenanthrene) | 96            |               | 43 - 129 |
| (d12-Chrysene)     | 97            |               | 52 - 144 |
| (d12-Perylene)     | 99            |               | 36 - 161 |
| (d8-Naphthalene)   | 84            |               | 25 - 125 |

**Lab Sample ID: 102964-BS2**  
**Matrix: BlankMatrix**  
**Analysis Batch: O-40074**

**Client Sample ID: Lab Control Sample Dup**  
**Prep Type: Total/NA**  
**Prep Batch: O-40074\_P**

| Analyte                      | Spike Added | LCS DUP Result | LCS DUP Qualifier | Unit | D | %Rec | %Rec Limits | RPD | RPD Limit |
|------------------------------|-------------|----------------|-------------------|------|---|------|-------------|-----|-----------|
| 1-Methylnaphthalene          | 0.5         | 0.49           |                   | µg/L |   | 98   | 31 - 128    | 3   | 30        |
| 1-Methylphenanthrene         | 0.5         | 0.452          |                   | µg/L |   | 90   | 66 - 127    | 0   | 30        |
| 2,3,5-Trimethylnaphthalene   | 0.5         | 0.516          |                   | µg/L |   | 103  | 55 - 122    | 3   | 30        |
| 2,6-Dimethylnaphthalene      | 0.5         | 0.505          |                   | µg/L |   | 101  | 48 - 120    | 2   | 30        |
| 2-Methylnaphthalene          | 1.5         | 1.48           |                   | µg/L |   | 99   | 47 - 130    | 3   | 30        |
| Acenaphthene                 | 1.5         | 1.54           |                   | µg/L |   | 103  | 53 - 131    | 4   | 30        |
| Acenaphthylene               | 1.5         | 1.57           |                   | µg/L |   | 105  | 43 - 140    | 4   | 30        |
| Anthracene                   | 1.5         | 1.56           |                   | µg/L |   | 104  | 58 - 135    | 3   | 30        |
| Benz[a]anthracene            | 1.5         | 1.58           |                   | µg/L |   | 105  | 55 - 145    | 1   | 30        |
| Benzo[a]pyrene               | 1.5         | 1.66           |                   | µg/L |   | 111  | 51 - 143    | 7   | 30        |
| Benzo[b]fluoranthene         | 1.5         | 1.66           |                   | µg/L |   | 111  | 46 - 165    | 2   | 30        |
| Benzo[e]pyrene               | 0.5         | 0.504          |                   | µg/L |   | 101  | 42 - 152    | 1   | 30        |
| Benzo[g,h,i]perylene         | 1.5         | 1.57           |                   | µg/L |   | 105  | 63 - 133    | 1   | 30        |
| Benzo[k]fluoranthene         | 1.5         | 1.67           |                   | µg/L |   | 111  | 56 - 145    | 2   | 30        |
| Biphenyl                     | 0.5         | 0.54           |                   | µg/L |   | 108  | 56 - 119    | 3   | 30        |
| Chrysene                     | 1.5         | 1.57           |                   | µg/L |   | 105  | 56 - 141    | 2   | 30        |
| Dibenz[a,h]anthracene        | 1.5         | 1.5            |                   | µg/L |   | 100  | 55 - 150    | 3   | 30        |
| Dibenzo[a,l]pyrene           | 0.5         | 0.52           |                   | µg/L |   | 104  | 50 - 150    | 4   | 30        |
| Dibenzothiophene             | 0.5         | 0.469          |                   | µg/L |   | 94   | 46 - 126    | 2   | 30        |
| Disalicylideneprapanediamine | 50          | 45.8           |                   | µg/L |   | 92   | 50 - 150    | 6   | 30        |
| Fluoranthene                 | 1.5         | 1.72           |                   | µg/L |   | 115  | 60 - 146    | 3   | 30        |
| Fluorene                     | 1.5         | 1.64           |                   | µg/L |   | 109  | 58 - 131    | 4   | 30        |
| Indeno[1,2,3-cd]pyrene       | 1.5         | 1.67           |                   | µg/L |   | 111  | 50 - 151    | 4   | 30        |
| Naphthalene                  | 1.5         | 1.42           |                   | µg/L |   | 95   | 41 - 126    | 3   | 30        |

Eurofins Eaton Monrovia

# QC Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 625 PAH Physis LL (EAL) + TICs - EPA 625 Base/Neutral and Acid Organics i (Continued)

**Lab Sample ID: 102964-BS2**  
**Matrix: BlankMatrix**  
**Analysis Batch: O-40074**

**Client Sample ID: Lab Control Sample Dup**  
**Prep Type: Total/NA**  
**Prep Batch: O-40074\_P**

| Analyte      | Spike Added | LCS DUP Result | LCS DUP Qualifier | Unit | D | %Rec | %Rec Limits | RPD | RPD Limit |
|--------------|-------------|----------------|-------------------|------|---|------|-------------|-----|-----------|
| Perylene     | 0.5         | 0.506          |                   | µg/L |   | 101  | 48 - 141    | 5   | 30        |
| Phenanthrene | 1.5         | 1.6            |                   | µg/L |   | 107  | 67 - 127    | 2   | 30        |
| Pyrene       | 1.5         | 1.65           |                   | µg/L |   | 110  | 54 - 156    | 2   | 30        |

| Surrogate          | LCS DUP %Recovery | LCS DUP Qualifier | Limits   |
|--------------------|-------------------|-------------------|----------|
| (d10-Acenaphthene) | 92                |                   | 27 - 133 |
| (d10-Phenanthrene) | 95                |                   | 43 - 129 |
| (d12-Chrysene)     | 95                |                   | 52 - 144 |
| (d12-Perylene)     | 94                |                   | 36 - 161 |
| (d8-Naphthalene)   | 84                |                   | 25 - 125 |

## Method: 8015 Gas (Purgeable) LL (EAL) - SW846 8015B Gasoline Range Organics

**Lab Sample ID: 22VG39L08B**  
**Matrix: WATER**  
**Analysis Batch: 22VG39L08**

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

| Analyte  | MB Result | MB Qualifier | RL    | MDL | Unit | D | Prepared | Analyzed       | Dil Fac |
|----------|-----------|--------------|-------|-----|------|---|----------|----------------|---------|
| GASOLINE | ND        | U            | 0.020 |     | mg/L |   |          | 12/22/22 11:58 | 1       |

| Surrogate          | MB %Recovery | MB Qualifier | Limits | Prepared | Analyzed       | Dil Fac |
|--------------------|--------------|--------------|--------|----------|----------------|---------|
| BROMOFLUOROBENZENE |              |              |        |          | 12/22/22 11:58 | 1       |

**Lab Sample ID: 22VG39L08L**  
**Matrix: WATER**  
**Analysis Batch: 22VG39L08**

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

| Analyte  | Spike Added | LCS Result | LCS Qualifier | Unit | D | %Rec | %Rec Limits |
|----------|-------------|------------|---------------|------|---|------|-------------|
| GASOLINE | 0.500       | 0.471      |               | mg/L |   | 94   | 60 - 130    |

| Surrogate          | LCS %Recovery | LCS Qualifier | Limits   |
|--------------------|---------------|---------------|----------|
| BROMOFLUOROBENZENE | 110           |               | 70 - 130 |

**Lab Sample ID: 22L294-01M**  
**Matrix: WATER**  
**Analysis Batch: 22VG39L08**

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

| Analyte  | Sample Result | Sample Qualifier | Spike Added | MS Result | MS Qualifier | Unit | D | %Rec | %Rec Limits |
|----------|---------------|------------------|-------------|-----------|--------------|------|---|------|-------------|
| GASOLINE | ND            |                  | 0.500       | 0.494     |              | mg/L |   | 99   | 50 - 130    |

| Surrogate          | MS %Recovery | MS Qualifier | Limits   |
|--------------------|--------------|--------------|----------|
| BROMOFLUOROBENZENE | 108          |              | 60 - 140 |

# QC Sample Results

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 8015 Gas (Purgeable) LL (EAL) - SW846 8015B Gasoline Range Organics (Continued)

**Lab Sample ID: 22L294-01S**  
**Matrix: WATER**  
**Analysis Batch: 22VG39L08**

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

| Analyte            | Sample Result    | Sample Qualifier     | Spike Added       | MSD Result | MSD Qualifier | Unit | D | %Rec | %Rec Limits | RPD | RPD Limit |
|--------------------|------------------|----------------------|-------------------|------------|---------------|------|---|------|-------------|-----|-----------|
| GASOLINE           | ND               |                      | 0.500             | 0.484      |               | mg/L |   | 97   | 50 - 130    | 2   | 30        |
| <b>Surrogate</b>   | <b>%Recovery</b> | <b>MSD Qualifier</b> | <b>MSD Limits</b> |            |               |      |   |      |             |     |           |
| BROMOFLUOROBENZENE | 110              |                      | 60 - 140          |            |               |      |   |      |             |     |           |

## Method: 8015 LL DRO/MRO/JP5/JP8 - 8015 - TPH DRO/ORO

**Lab Sample ID: 22DSL040WB**  
**Matrix: WATER**  
**Analysis Batch: 22DSL040W**

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

| Analyte          | MB Result        | MB Qualifier        | RL               | MDL | Unit | D               | Prepared        | Analyzed       | Dil Fac |
|------------------|------------------|---------------------|------------------|-----|------|-----------------|-----------------|----------------|---------|
| DIESEL           | ND               | U                   | 0.025            |     | mg/L |                 |                 | 01/05/23 20:12 | 1       |
| JP5              | ND               | U                   | 0.050            |     | mg/L |                 |                 | 01/05/23 20:12 | 1       |
| JP8              | ND               | U                   | 0.050            |     | mg/L |                 |                 | 01/05/23 20:12 | 1       |
| MOTOR OIL        | ND               | U                   | 0.050            |     | mg/L |                 |                 | 01/05/23 20:12 | 1       |
| <b>Surrogate</b> | <b>%Recovery</b> | <b>MB Qualifier</b> | <b>MB Limits</b> |     |      | <b>Prepared</b> | <b>Analyzed</b> | <b>Dil Fac</b> |         |
| BROMOBENZENE     |                  |                     |                  |     |      |                 | 01/05/23 20:12  | 1              |         |
| HEXACOSANE       |                  |                     |                  |     |      |                 | 01/05/23 20:12  | 1              |         |

**Lab Sample ID: 22DSL040WL**  
**Matrix: WATER**  
**Analysis Batch: 22DSL040W**

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

| Analyte          | Spike Added      | LCS Result           | LCS Qualifier     | Unit | D | %Rec | %Rec Limits |
|------------------|------------------|----------------------|-------------------|------|---|------|-------------|
| DIESEL           | 2.50             | 2.83                 |                   | mg/L |   | 113  | 50 - 130    |
| <b>Surrogate</b> | <b>%Recovery</b> | <b>LCS Qualifier</b> | <b>LCS Limits</b> |      |   |      |             |
| BROMOBENZENE     | 102              |                      | 60 - 130          |      |   |      |             |
| HEXACOSANE       | 106              |                      | 60 - 130          |      |   |      |             |

**Lab Sample ID: 22J5L040WL**  
**Matrix: WATER**  
**Analysis Batch: 22DSL040W**

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

| Analyte          | Spike Added      | LCS Result           | LCS Qualifier     | Unit | D | %Rec | %Rec Limits |
|------------------|------------------|----------------------|-------------------|------|---|------|-------------|
| JP5              | 2.50             | 2.43                 |                   | mg/L |   | 97   | 30 - 160    |
| <b>Surrogate</b> | <b>%Recovery</b> | <b>LCS Qualifier</b> | <b>LCS Limits</b> |      |   |      |             |
| BROMOBENZENE     | 105              |                      | 60 - 130          |      |   |      |             |
| HEXACOSANE       | 97               |                      | 60 - 130          |      |   |      |             |

# QC Sample Results

Client: City & County of Honolulu  
 Project/Site: RED-HILL

Job ID: 380-32096-1

## Method: 8015 LL DRO/MRO/JP5/JP8 - 8015 - TPH DRO/ORO (Continued)

**Lab Sample ID: 22J8L040WL**  
**Matrix: WATER**  
**Analysis Batch: 22DSL040W**

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

| Analyte          | Spike<br>Added   | LCS<br>Result    | LCS<br>Qualifier | Unit | D | %Rec | %Rec<br>Limits |
|------------------|------------------|------------------|------------------|------|---|------|----------------|
| JP8              | 2.50             | 2.90             |                  | mg/L |   | 116  | 30 - 160       |
| <b>Surrogate</b> |                  |                  |                  |      |   |      |                |
|                  | <b>LCS</b>       | <b>LCS</b>       |                  |      |   |      |                |
|                  | <b>%Recovery</b> | <b>Qualifier</b> |                  |      |   |      | <b>Limits</b>  |
| BROMOBENZENE     | 98               |                  |                  |      |   |      | 60 - 130       |
| HEXACOSANE       | 102              |                  |                  |      |   |      | 60 - 130       |



# QC Association Summary

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## GC/MS Semi VOA

### Prep Batch: 27725

| Lab Sample ID       | Client Sample ID        | Prep Type | Matrix         | Method | Prep Batch |
|---------------------|-------------------------|-----------|----------------|--------|------------|
| 380-32096-1         | Aiea Gulch Wells Pump 2 | Total/NA  | Drinking Water | 525.2  |            |
| 380-32096-2         | Aiea Wells Pump 2       | Total/NA  | Drinking Water | 525.2  |            |
| 380-32096-3         | Halawa Wells Pump 1     | Total/NA  | Drinking Water | 525.2  |            |
| MB 380-27725/1-A    | Method Blank            | Total/NA  | Water          | 525.2  |            |
| LCS 380-27725/3-A   | Lab Control Sample      | Total/NA  | Water          | 525.2  |            |
| LCSD 380-27725/4-A  | Lab Control Sample Dup  | Total/NA  | Water          | 525.2  |            |
| MRL 380-27725/2-A   | Lab Control Sample      | Total/NA  | Water          | 525.2  |            |
| 380-32064-C-1-A MS  | Matrix Spike            | Total/NA  | Water          | 525.2  |            |
| 380-32064-I-1-A MSD | Matrix Spike Duplicate  | Total/NA  | Water          | 525.2  |            |

### Analysis Batch: 28001

| Lab Sample ID       | Client Sample ID        | Prep Type | Matrix         | Method | Prep Batch |
|---------------------|-------------------------|-----------|----------------|--------|------------|
| 380-32096-1         | Aiea Gulch Wells Pump 2 | Total/NA  | Drinking Water | 525.2  | 27725      |
| 380-32096-2         | Aiea Wells Pump 2       | Total/NA  | Drinking Water | 525.2  | 27725      |
| 380-32096-3         | Halawa Wells Pump 1     | Total/NA  | Drinking Water | 525.2  | 27725      |
| MB 380-27725/1-A    | Method Blank            | Total/NA  | Water          | 525.2  | 27725      |
| LCS 380-27725/3-A   | Lab Control Sample      | Total/NA  | Water          | 525.2  | 27725      |
| LCSD 380-27725/4-A  | Lab Control Sample Dup  | Total/NA  | Water          | 525.2  | 27725      |
| MRL 380-27725/2-A   | Lab Control Sample      | Total/NA  | Water          | 525.2  | 27725      |
| 380-32064-C-1-A MS  | Matrix Spike            | Total/NA  | Water          | 525.2  | 27725      |
| 380-32064-I-1-A MSD | Matrix Spike Duplicate  | Total/NA  | Water          | 525.2  | 27725      |

## Subcontract

### Analysis Batch: O-40074

| Lab Sample ID | Client Sample ID        | Prep Type | Matrix         | Method                            | Prep Batch |
|---------------|-------------------------|-----------|----------------|-----------------------------------|------------|
| 380-32096-1   | Aiea Gulch Wells Pump 2 | Total/NA  | Drinking Water | 625 PAH Physis<br>LL (EAL) + TICs | O-40074_P  |
| 380-32096-2   | Aiea Wells Pump 2       | Total/NA  | Drinking Water | 625 PAH Physis<br>LL (EAL) + TICs | O-40074_P  |
| 380-32096-3   | Halawa Wells Pump 1     | Total/NA  | Drinking Water | 625 PAH Physis<br>LL (EAL) + TICs | O-40074_P  |
| 102964-B1     | Method Blank            | Total/NA  | BlankMatrix    | 625 PAH Physis<br>LL (EAL) + TICs | O-40074_P  |
| 102964-BS1    | Lab Control Sample      | Total/NA  | BlankMatrix    | 625 PAH Physis<br>LL (EAL) + TICs | O-40074_P  |
| 102964-BS2    | Lab Control Sample Dup  | Total/NA  | BlankMatrix    | 625 PAH Physis<br>LL (EAL) + TICs | O-40074_P  |

### Analysis Batch: 22DSL040W

| Lab Sample ID | Client Sample ID        | Prep Type | Matrix         | Method                         | Prep Batch |
|---------------|-------------------------|-----------|----------------|--------------------------------|------------|
| 380-32096-1   | Aiea Gulch Wells Pump 2 | Total/NA  | Drinking Water | 8015 LL<br>DRO/MRO/JP5/J<br>P8 |            |
| 380-32096-2   | Aiea Wells Pump 2       | Total/NA  | Drinking Water | 8015 LL<br>DRO/MRO/JP5/J<br>P8 |            |
| 380-32096-3   | Halawa Wells Pump 1     | Total/NA  | Drinking Water | 8015 LL<br>DRO/MRO/JP5/J<br>P8 |            |
| 22DSL040WB    | Method Blank            | Total/NA  | WATER          | 8015 LL<br>DRO/MRO/JP5/J<br>P8 |            |

Eurofins Eaton Monrovia

# QC Association Summary

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Subcontract (Continued)

### Analysis Batch: 22DSL040W (Continued)

| Lab Sample ID | Client Sample ID   | Prep Type | Matrix | Method                         | Prep Batch |
|---------------|--------------------|-----------|--------|--------------------------------|------------|
| 22DSL040WL    | Lab Control Sample | Total/NA  | WATER  | 8015 LL<br>DRO/MRO/JP5/J<br>P8 |            |
| 22J5L040WL    | Lab Control Sample | Total/NA  | WATER  | 8015 LL<br>DRO/MRO/JP5/J<br>P8 |            |
| 22J8L040WL    | Lab Control Sample | Total/NA  | WATER  | 8015 LL<br>DRO/MRO/JP5/J<br>P8 |            |

### Analysis Batch: 22VG39L08

| Lab Sample ID | Client Sample ID            | Prep Type | Matrix         | Method                              | Prep Batch |
|---------------|-----------------------------|-----------|----------------|-------------------------------------|------------|
| 380-32096-1   | Aiea Gulch Wells Pump 2     | Total/NA  | Drinking Water | 8015 Gas<br>(Purgeable) LL<br>(EAL) |            |
| 380-32096-2   | Aiea Wells Pump 2           | Total/NA  | Drinking Water | 8015 Gas<br>(Purgeable) LL<br>(EAL) |            |
| 380-32096-3   | Halawa Wells Pump 1         | Total/NA  | Drinking Water | 8015 Gas<br>(Purgeable) LL<br>(EAL) |            |
| 380-32096-4   | TB: Aiea Gulch Wells Pump 2 | Total/NA  | Drinking Water | 8015 Gas<br>(Purgeable) LL<br>(EAL) |            |
| 380-32096-5   | TB: Aiea Wells Pump 2       | Total/NA  | Drinking Water | 8015 Gas<br>(Purgeable) LL<br>(EAL) |            |
| 380-32096-6   | TB: Halawa Wells Pump 1     | Total/NA  | Drinking Water | 8015 Gas<br>(Purgeable) LL<br>(EAL) |            |
| 22VG39L08B    | Method Blank                | Total/NA  | WATER          | 8015 Gas<br>(Purgeable) LL<br>(EAL) |            |
| 22VG39L08L    | Lab Control Sample          | Total/NA  | WATER          | 8015 Gas<br>(Purgeable) LL<br>(EAL) |            |
| 22L294-01M    | Matrix Spike                | Total/NA  | WATER          | 8015 Gas<br>(Purgeable) LL<br>(EAL) |            |
| 22L294-01S    | Matrix Spike Duplicate      | Total/NA  | WATER          | 8015 Gas<br>(Purgeable) LL<br>(EAL) |            |

### Prep Batch: O-40074\_P

| Lab Sample ID | Client Sample ID        | Prep Type | Matrix         | Method  | Prep Batch |
|---------------|-------------------------|-----------|----------------|---------|------------|
| 380-32096-1   | Aiea Gulch Wells Pump 2 | Total/NA  | Drinking Water | EPA_625 |            |
| 380-32096-2   | Aiea Wells Pump 2       | Total/NA  | Drinking Water | EPA_625 |            |
| 380-32096-3   | Halawa Wells Pump 1     | Total/NA  | Drinking Water | EPA_625 |            |
| 102964-B1     | Method Blank            | Total/NA  | BlankMatrix    | EPA_625 |            |
| 102964-BS1    | Lab Control Sample      | Total/NA  | BlankMatrix    | EPA_625 |            |
| 102964-BS2    | Lab Control Sample Dup  | Total/NA  | BlankMatrix    | EPA_625 |            |

# Lab Chronicle

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Client Sample ID: Aiea Gulch Wells Pump 2

Date Collected: 12/19/22 10:00

Date Received: 12/21/22 10:00

## Lab Sample ID: 380-32096-1

Matrix: Drinking Water

| Prep Type | Batch Type | Batch Method                   | Run | Dilution Factor | Batch Number | Analyst | Lab    | Prepared or Analyzed |
|-----------|------------|--------------------------------|-----|-----------------|--------------|---------|--------|----------------------|
| Total/NA  | Prep       | 525.2                          |     |                 | 27725        | OTM3    | EA MON | 12/22/22 06:58       |
| Total/NA  | Analysis   | 525.2                          |     | 1               | 28001        | Q8LA    | EA MON | 12/27/22 14:45       |
| Total/NA  | Prep       | EPA_625                        |     | 1               | O-40074_P    |         |        | 12/26/22 00:00       |
| Total/NA  | Analysis   | 625 PAH Physis LL (EAL) + TICs |     | 1               | O-40074      | YC      |        | 01/08/23 07:38       |
| Total/NA  | Analysis   | 8015 Gas (Purgeable) LL (EAL)  |     | 1               | 22VG39L08    | SCerva  |        | 12/22/22 19:09       |
| Total/NA  | Analysis   | 8015 LL DRO/MRO/JP5/JP8        |     | 1               | 22DSL040W    | SDees   |        | 01/05/23 21:25       |

## Client Sample ID: Aiea Wells Pump 2

Date Collected: 12/19/22 10:45

Date Received: 12/21/22 10:00

## Lab Sample ID: 380-32096-2

Matrix: Drinking Water

| Prep Type | Batch Type | Batch Method                   | Run | Dilution Factor | Batch Number | Analyst | Lab    | Prepared or Analyzed |
|-----------|------------|--------------------------------|-----|-----------------|--------------|---------|--------|----------------------|
| Total/NA  | Prep       | 525.2                          |     |                 | 27725        | OTM3    | EA MON | 12/22/22 06:58       |
| Total/NA  | Analysis   | 525.2                          |     | 1               | 28001        | Q8LA    | EA MON | 12/27/22 15:05       |
| Total/NA  | Prep       | EPA_625                        |     | 1               | O-40074_P    |         |        | 12/26/22 00:00       |
| Total/NA  | Analysis   | 625 PAH Physis LL (EAL) + TICs |     | 1               | O-40074      | YC      |        | 01/08/23 09:22       |
| Total/NA  | Analysis   | 8015 Gas (Purgeable) LL (EAL)  |     | 1               | 22VG39L08    | SCerva  |        | 12/22/22 20:57       |
| Total/NA  | Analysis   | 8015 LL DRO/MRO/JP5/JP8        |     | 1               | 22DSL040W    | SDees   |        | 01/05/23 21:44       |

## Client Sample ID: Halawa Wells Pump 1

Date Collected: 12/19/22 11:18

Date Received: 12/21/22 10:00

## Lab Sample ID: 380-32096-3

Matrix: Drinking Water

| Prep Type | Batch Type | Batch Method                   | Run | Dilution Factor | Batch Number | Analyst | Lab    | Prepared or Analyzed |
|-----------|------------|--------------------------------|-----|-----------------|--------------|---------|--------|----------------------|
| Total/NA  | Prep       | 525.2                          |     |                 | 27725        | OTM3    | EA MON | 12/22/22 06:58       |
| Total/NA  | Analysis   | 525.2                          |     | 1               | 28001        | Q8LA    | EA MON | 12/27/22 15:26       |
| Total/NA  | Prep       | EPA_625                        |     | 1               | O-40074_P    |         |        | 12/26/22 00:00       |
| Total/NA  | Analysis   | 625 PAH Physis LL (EAL) + TICs |     | 1               | O-40074      | YC      |        | 01/08/23 11:07       |
| Total/NA  | Analysis   | 8015 Gas (Purgeable) LL (EAL)  |     | 1               | 22VG39L08    | SCerva  |        | 12/22/22 21:33       |
| Total/NA  | Analysis   | 8015 LL DRO/MRO/JP5/JP8        |     | 1               | 22DSL040W    | SDees   |        | 01/05/23 22:02       |

## Client Sample ID: TB: Aiea Gulch Wells Pump 2

Date Collected: 12/19/22 10:00

Date Received: 12/21/22 10:00

## Lab Sample ID: 380-32096-4

Matrix: Drinking Water

| Prep Type | Batch Type | Batch Method                  | Run | Dilution Factor | Batch Number | Analyst | Lab | Prepared or Analyzed |
|-----------|------------|-------------------------------|-----|-----------------|--------------|---------|-----|----------------------|
| Total/NA  | Analysis   | 8015 Gas (Purgeable) LL (EAL) |     | 1               | 22VG39L08    | SCerva  |     | 12/22/22 22:08       |



# Lab Chronicle

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

## Client Sample ID: TB: Aiea Wells Pump 2

Date Collected: 12/19/22 10:45

Date Received: 12/21/22 10:00

Lab Sample ID: 380-32096-5

Matrix: Drinking Water

| Prep Type | Batch Type | Batch Method                     | Run | Dilution Factor | Batch Number | Analyst | Lab | Prepared or Analyzed |
|-----------|------------|----------------------------------|-----|-----------------|--------------|---------|-----|----------------------|
| Total/NA  | Analysis   | 8015 Gas<br>(Purgeable) LL (EAL) |     | 1               | 22VG39L08    | SCerva  |     | 12/22/22 22:44       |

## Client Sample ID: TB: Halawa Wells Pump 1

Date Collected: 12/19/22 11:18

Date Received: 12/21/22 10:00

Lab Sample ID: 380-32096-6

Matrix: Drinking Water

| Prep Type | Batch Type | Batch Method                     | Run | Dilution Factor | Batch Number | Analyst | Lab | Prepared or Analyzed |
|-----------|------------|----------------------------------|-----|-----------------|--------------|---------|-----|----------------------|
| Total/NA  | Analysis   | 8015 Gas<br>(Purgeable) LL (EAL) |     | 1               | 22VG39L08    | SCerva  |     | 12/22/22 23:20       |

### Laboratory References:

= Physis Environmental Laboratories, 1904 Wright Circle, Anaheim, CA 92806

EA MON = Eurofins Eaton Monrovia, 750 Royal Oaks Drive, Suite 100, Monrovia, CA 91016, TEL (626)386-1100



# Accreditation/Certification Summary

Client: City & County of Honolulu  
 Project/Site: RED-HILL

Job ID: 380-32096-1

## Laboratory: Eurofins Eaton Monrovia

Unless otherwise noted, all analytes for this laboratory were covered under each accreditation/certification below.

| Authority | Program | Identification Number | Expiration Date |
|-----------|---------|-----------------------|-----------------|
| Hawaii    | State   | CA00006               | 01-08-23        |

The following analytes are included in this report, but the laboratory is not certified by the governing authority. This list may include analytes for which the agency does not offer certification.

| Analysis Method | Prep Method | Matrix         | Analyte                          |
|-----------------|-------------|----------------|----------------------------------|
| 525.2           | 525.2       | Drinking Water | 2,4'-DDD                         |
| 525.2           | 525.2       | Drinking Water | 2,4'-DDE                         |
| 525.2           | 525.2       | Drinking Water | 2,4'-DDT                         |
| 525.2           | 525.2       | Drinking Water | 2,4-Dinitrotoluene               |
| 525.2           | 525.2       | Drinking Water | 2,6-Dinitrotoluene               |
| 525.2           | 525.2       | Drinking Water | 4,4'-DDD                         |
| 525.2           | 525.2       | Drinking Water | 4,4'-DDE                         |
| 525.2           | 525.2       | Drinking Water | 4,4'-DDT                         |
| 525.2           | 525.2       | Drinking Water | Acenaphthene                     |
| 525.2           | 525.2       | Drinking Water | Acenaphthylene                   |
| 525.2           | 525.2       | Drinking Water | Acetochlor                       |
| 525.2           | 525.2       | Drinking Water | alpha-BHC                        |
| 525.2           | 525.2       | Drinking Water | alpha-Chlordane                  |
| 525.2           | 525.2       | Drinking Water | Anthracene                       |
| 525.2           | 525.2       | Drinking Water | Benz(a)anthracene                |
| 525.2           | 525.2       | Drinking Water | Benzo[b]fluoranthene             |
| 525.2           | 525.2       | Drinking Water | Benzo[g,h,i]perylene             |
| 525.2           | 525.2       | Drinking Water | Benzo[k]fluoranthene             |
| 525.2           | 525.2       | Drinking Water | beta-BHC                         |
| 525.2           | 525.2       | Drinking Water | Bromacil                         |
| 525.2           | 525.2       | Drinking Water | Butylbenzylphthalate             |
| 525.2           | 525.2       | Drinking Water | Caffeine                         |
| 525.2           | 525.2       | Drinking Water | Chlorobenzilate                  |
| 525.2           | 525.2       | Drinking Water | Chloroneb                        |
| 525.2           | 525.2       | Drinking Water | Chlorothalonil (Draconil, Bravo) |
| 525.2           | 525.2       | Drinking Water | Chlorpyrifos                     |
| 525.2           | 525.2       | Drinking Water | Chrysene                         |
| 525.2           | 525.2       | Drinking Water | delta-BHC                        |
| 525.2           | 525.2       | Drinking Water | Diazinon (Qualitative)           |
| 525.2           | 525.2       | Drinking Water | Dibenz(a,h)anthracene            |
| 525.2           | 525.2       | Drinking Water | Diclorvos (DDVP)                 |
| 525.2           | 525.2       | Drinking Water | Diethylphthalate                 |
| 525.2           | 525.2       | Drinking Water | Dimethoate                       |
| 525.2           | 525.2       | Drinking Water | Dimethylphthalate                |
| 525.2           | 525.2       | Drinking Water | Di-n-butyl phthalate             |
| 525.2           | 525.2       | Drinking Water | Di-n-octyl phthalate             |
| 525.2           | 525.2       | Drinking Water | Endosulfan I (Alpha)             |
| 525.2           | 525.2       | Drinking Water | Endosulfan II (Beta)             |
| 525.2           | 525.2       | Drinking Water | Endosulfan sulfate               |
| 525.2           | 525.2       | Drinking Water | Endrin aldehyde                  |
| 525.2           | 525.2       | Drinking Water | EPTC                             |
| 525.2           | 525.2       | Drinking Water | Fluoranthene                     |
| 525.2           | 525.2       | Drinking Water | Fluorene                         |
| 525.2           | 525.2       | Drinking Water | gamma-Chlordane                  |
| 525.2           | 525.2       | Drinking Water | Indeno[1,2,3-cd]pyrene           |

# Accreditation/Certification Summary

Client: City & County of Honolulu  
 Project/Site: RED-HILL

Job ID: 380-32096-1

## Laboratory: Eurofins Eaton Monrovia (Continued)

Unless otherwise noted, all analytes for this laboratory were covered under each accreditation/certification below.

| Authority | Program | Identification Number | Expiration Date |
|-----------|---------|-----------------------|-----------------|
|-----------|---------|-----------------------|-----------------|

The following analytes are included in this report, but the laboratory is not certified by the governing authority. This list may include analytes for which the agency does not offer certification.

| Analysis Method | Prep Method | Matrix         | Analyte                          |
|-----------------|-------------|----------------|----------------------------------|
| 525.2           | 525.2       | Drinking Water | Isophorone                       |
| 525.2           | 525.2       | Drinking Water | Malathion                        |
| 525.2           | 525.2       | Drinking Water | Molinate                         |
| 525.2           | 525.2       | Drinking Water | Naphthalene                      |
| 525.2           | 525.2       | Drinking Water | Parathion                        |
| 525.2           | 525.2       | Drinking Water | Pendimethalin (Penoxaline)       |
| 525.2           | 525.2       | Drinking Water | Phenanthrene                     |
| 525.2           | 525.2       | Drinking Water | Pyrene                           |
| 525.2           | 525.2       | Drinking Water | Terbacil                         |
| 525.2           | 525.2       | Drinking Water | Terbutylazine                    |
| 525.2           | 525.2       | Drinking Water | Thiobencarb                      |
| 525.2           | 525.2       | Drinking Water | Total Permethrin (mixed isomers) |
| 525.2           | 525.2       | Drinking Water | trans-Nonachlor                  |
| 525.2           | 525.2       | Drinking Water | Trifluralin                      |



# Method Summary

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

| Method | Method Description                       | Protocol | Laboratory |
|--------|--|----------|------------|
| 525.2  | Semivolatile Organic Compounds (GC/MS)   | EPA      | EA MON     |
| 625    | EPA 625 Base/Neutral and Acid Organics i | EPA      |            |
| 8015   | 8015 - TPH DRO/ORO                       | EPA      |            |
| 8015B  | SW846 8015B Gasoline Range Organics      | SW846    |            |
| 525.2  | Extraction of Semivolatile Compounds     | EPA      | EA MON     |

**Protocol References:**

EPA = US Environmental Protection Agency

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

**Laboratory References:**

= Physis Environmental Laboratories, 1904 Wright Circle, Anaheim, CA 92806

EA MON = Eurofins Eaton Monrovia, 750 Royal Oaks Drive, Suite 100, Monrovia, CA 91016, TEL (626)386-1100



# Sample Summary

Client: City & County of Honolulu  
Project/Site: RED-HILL

Job ID: 380-32096-1

| Lab Sample ID | Client Sample ID            | Matrix         | Collected      | Received       | PWSID Number |
|---------------|-----------------------------|----------------|----------------|----------------|--------------|
| 380-32096-1   | Aiea Gulch Wells Pump 2     | Drinking Water | 12/19/22 10:00 | 12/21/22 10:00 | HI0000331    |
| 380-32096-2   | Aiea Wells Pump 2           | Drinking Water | 12/19/22 10:45 | 12/21/22 10:00 | HI0000331    |
| 380-32096-3   | Halawa Wells Pump 1         | Drinking Water | 12/19/22 11:18 | 12/21/22 10:00 | HI0000331    |
| 380-32096-4   | TB: Aiea Gulch Wells Pump 2 | Drinking Water | 12/19/22 10:00 | 12/21/22 10:00 |              |
| 380-32096-5   | TB: Aiea Wells Pump 2       | Drinking Water | 12/19/22 10:45 | 12/21/22 10:00 |              |
| 380-32096-6   | TB: Halawa Wells Pump 1     | Drinking Water | 12/19/22 11:18 | 12/21/22 10:00 |              |

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3051 Fujita Street  
 Torrance, CA 90505  
 Tel: (310)-618-8889

Date: 01-20-2023  
 EMAX Batch No.: 22L294

Attn: Jackie Contreras

Eurofins Eaton Analytical  
 750 Royal Oaks Dr., Suite 100  
 Monrovia, CA 91016-3629

Subject: Laboratory Report  
 Project: 380-32096

.....

Enclosed is the Laboratory report for samples received on 12/22/22.  
 The data reported relate only to samples listed below :

| Sample ID      | Control # | Col Date | Matrix | Analysis            |
|----------------|-----------|----------|--------|---------------------|
| 380-32096-1    | L294-01   | 12/19/22 | WATER  | TPH GASOLINE<br>TPH |
| 380-32096-2    | L294-02   | 12/19/22 | WATER  | TPH GASOLINE<br>TPH |
| 380-32096-3    | L294-03   | 12/19/22 | WATER  | TPH GASOLINE<br>TPH |
| 380-32096-4    | L294-04   | 12/19/22 | WATER  | TPH GASOLINE        |
| 380-32096-5    | L294-05   | 12/19/22 | WATER  | TPH GASOLINE        |
| 380-32096-6    | L294-06   | 12/19/22 | WATER  | TPH GASOLINE        |
| 380-32096-1MS  | L294-01M  | 12/19/22 | WATER  | TPH GASOLINE        |
| 380-32096-1MSD | L294-01S  | 12/19/22 | WATER  | TPH GASOLINE        |

The results are summarized on the following pages.

Please feel free to call if you have any questions concerning these results.

Sincerely yours,



Caspar J. Pang  
 Laboratory Director

This report is confidential and intended solely for the use of the individual or entity to whom it is addressed. This report shall not be reproduced except in full or without the written approval of EMAX.

EMAX certifies that results included in this report meets all TNI & DOD requirements unless noted in the Case Narrative.

NELAP Accredited Certificate Number CA002912022-22  
 ANAB Accredited DoD ELAP and ISO/IEC 17025 Certificate Number L2278 Testing  
 California ELAP Accredited Certificate Number 2672



**Monrovia, CA (Suite 100)**  
 750 Royal Oaks Drive Suite 100  
 Monrovia, CA 91016  
 Phone: 626-386-1100

# Chain of Custody Record

222294



eurofins

Environment Testing

|  |  |  |  |  |  |   |  |
|--|--|--|--|--|--|---|--|
| <b>Client Information (Sub Contract Lab)</b><br>Client Contact: Rachelle Arada<br>Shipping/Receiving: Rachelle Arada<br>Company: EMAX Laboratories Inc<br>Address: 3051 Fujita Street, Torrance, CA 90505<br>City: Torrance, State: CA, Zip: 90505<br>Phone: PO #, WO #, Project #: 38001111, SSON#: |  | Lab PM: Arada, Rachelle<br>E-Mail: Rachelle.Arada@et.eurofinsus.com<br>Accreditations Required (See note): State - Hawaii  |  | Camer Tracking No(s): 380-31670.1<br>State of Origin: Hawaii<br>Page: Page 1 of 1<br>Job #: 380-32096-1  |  | COC No: 380-31670.1<br>Preservation Codes:<br>A - HCL<br>B - NaOH<br>C - Zn Acetate<br>D - Nitric Acid<br>E - NaHSO4<br>F - MeOH<br>G - Amchlor<br>H - Ascorbic Acid<br>I - Ice<br>J - DI Water<br>K - EDTA<br>L - EDA<br>Z - other (specify)<br>Other: |  |
| <b>Sample Information</b><br>Sample Date: 12/19/22<br>Sample Time: 10:00 Hawaiian<br>Sample Type (C=comp, G=grab): Water<br>Matrix (W=water, S=solid, O=other): Water  |  | Due Date Requested: 1/9/2023<br>TAT Requested (days):  |  | Analysis Requested:  |  | Total Number of Containers: 6   |  |
| Sample ID (Lab ID): Aiea Gulch Wells Pump 2 (380-32096-1)<br>Aiea Wells Pump 2 (380-32096-2)<br>Halawa Wells Pump 1 (380-32096-3)<br>TB: Aiea Gulch Wells Pump 2 (380-32096-4)<br>TB: Aiea Wells Pump 2 (380-32096-5)<br>TB: Halawa Wells Pump 1 (380-32096-6)                                       |  | Field Filtered Sample (Yes or No):<br>Perform MS/MSD (Yes or No):<br>SUB (8015 Gas (Purgeable) LL (EAL)) 8015 Gas<br>SUB (Purgeable) LL (EAL)<br>SUB (8015 LL DROM/RO/JPS/JP8) 8015 LL DROM/RO/JPS/JP8 |  | Special Instructions/Note:<br>See Attached Instructions<br>See Attached Instructions<br>See Attached Instructions<br>See Attached Instructions<br>See Attached Instructions<br>See Attached Instructions |  | Special Instructions/Note:<br>See Attached Instructions<br>See Attached Instructions<br>See Attached Instructions<br>See Attached Instructions<br>See Attached Instructions   |  |

Note: Since laboratory accreditations are subject to change, Eurofins Eaton Analytical, LLC places the ownership of method, analyte & accreditation compliance upon our subcontract laboratories. This sample shipment is forwarded under chain-of-custody. If the laboratory does not currently maintain accreditation in the State of Origin listed above for analysis/tests/matrix being analyzed, the samples must be shipped back to the Eurofins Eaton Analytical, LLC laboratory or other instructions will be provided. Any changes to accreditation status should be brought to Eurofins Eaton Analytical, LLC attention immediately. If all requested accreditations are current to date, return the signed Chain of Custody attesting to Eurofins Eaton Analytical, LLC.

**Possible Hazard Identification**  
 Unconfirmed  
 Deliverable Requested: I, II, III, IV, Other (specify) Primary Deliverable Rank: 2  
 Date: 12/22/2022 09:48  
 Relinquished by: J. G. PEPPER  
 Relinquished by: J. G. PEPPER  
 Relinquished by: J. G. PEPPER  
 Date/Time: 12/22/2022 09:48  
 Date/Time: 12/22/2022 14:45  
 Date/Time: 12/22/2022 14:45  
 Received by: J. G. PEPPER  
 Received by: J. G. PEPPER  
 Received by: J. G. PEPPER  
 Company: EEA  
 Company: EEA  
 Company: EEA  
 Special Instructions/OC Requirements: Return To Client  Disposal By Lab  Archive For  Months  
 Method of Shipment: 12-22-22 12:15  
 Date/Time: 12/22/2022 14:45  
 Date/Time: 12/22/2022 14:45  
 Date/Time: 12/22/2022 14:45  
 Company: EEA  
 Company: EEA  
 Company: EEA  
 Custody Seals Intact: Custody Seal No.:  
 REPORT ID: 22L294  
 Cooling Temperature(s) °C and Other Remarks: 4.1 2.9 K: -0.2





|   |                           |  |
|---|---------------------------|--|
| Type of Delivery<br><input type="checkbox"/> Fedex <input type="checkbox"/> UPS <input type="checkbox"/> GSO <input type="checkbox"/> Others<br><input type="checkbox"/> EMAX Courier <input checked="" type="checkbox"/> Client Delivery | Airbill / Tracking Number | ECN <b>22L294</b><br>Recipient <b>Jocelyne Solis-Ramos</b><br>Date <b>12/22/22</b> Time <b>14:45</b> |
|---|---------------------------|--|

**COC INSPECTION**

|   |   |  |  |  |  |
|---|---|--|--|--|--|
| <input checked="" type="checkbox"/> Client Name | <input type="checkbox"/> Client PM/PC                 | <input type="checkbox"/> Sampler Name        | <input checked="" type="checkbox"/> Sampling Date/Time | <input checked="" type="checkbox"/> Sample ID  | <input checked="" type="checkbox"/> Matrix |
| <input checked="" type="checkbox"/> Address     | <input checked="" type="checkbox"/> Tel # / Fax #     | <input type="checkbox"/> Courier Signature   | <input checked="" type="checkbox"/> Analysis Required  | <input type="checkbox"/> Preservative (if any) | <input checked="" type="checkbox"/> TAT    |
| Safety Issues (if any)                          | <input type="checkbox"/> High concentrations expected | <input type="checkbox"/> From Superfund Site | <input type="checkbox"/> Rad screening required        |  |  |

Note: \_\_\_\_\_

**PACKAGING INSPECTION**

|  |  |   |  |
|--|--|---|--|
| Container                                    | <input checked="" type="checkbox"/> Cooler                     | <input type="checkbox"/> Box                | <input type="checkbox"/> Other             |
| Condition                                    | <input type="checkbox"/> Custody Seal                          | <input type="checkbox"/> Intact             | <input type="checkbox"/> Damaged           |
| Packaging                                    | <input checked="" type="checkbox"/> Bubble Pack                | <input type="checkbox"/> Styrofoam          | <input type="checkbox"/> Popcorn           |
| Temperatures<br>(Cool, ≤6 °C but not frozen) | <input checked="" type="checkbox"/> Cooler 1 <b>4.1/3.9 °C</b> | <input type="checkbox"/> Cooler 2 _____ °C  | <input type="checkbox"/> Cooler 3 _____ °C |
| Thermometer:                                 | <input type="checkbox"/> Cooler 6 _____ °C                     | <input type="checkbox"/> Cooler 7 _____ °C  | <input type="checkbox"/> Cooler 8 _____ °C |
|  | <input type="checkbox"/> Cooler 9 _____ °C                     | <input type="checkbox"/> Cooler 10 _____ °C |  |

Comments:  Temperature is out of range. PM was informed IMMEDIATELY.

Note: \_\_\_\_\_

**DISCREPANCIES**

| LabSampleID                                       | LabSampleContainerID | Code | ClientSample Label ID / Information        | Corrective Action |
|---|----------------------|------|--|-------------------|
| 1-3   | 5,6,4,12,7,10        | 02   | JPG/JPB is not indicated on label          | RS                |
| 4-6   | 14,20,21-24          | 07   | two dates on label - 12/13/22 and 12/19/22 | RI                |
| <i>[Large diagonal scribble across the table]</i> |                      |      |  |                   |

pH holding time requirement for water samples is 15 mins. Water samples for pH analysis are received beyond 15 minutes from sampling time.

**NOTES/OBSERVATIONS:**

SAMPLE MATRIX IS DRINKING WATER?  YES  NO

- LEGEND:**
- |   |   |   |
|---|---|---|
| Code Description-Sample Management              | Code Description-Sample Management            | Code Description-Sample Management  |
| D1 Analysis is not indicated in _____           | D13 Out of Holding Time                       | R1 Proceed as indicated in <input checked="" type="checkbox"/> COC <input type="checkbox"/> Label |
| <b>D2</b> Analysis mismatch COC vs label        | D14 Bubble is >6mm                            | R2 Refer to attached instruction  |
| D3 Sample ID mismatch COC vs label              | D15 No trip blank in cooler                   | R3 Cancel the analysis  |
| D4 Sample ID is not indicated in _____          | D16 Preservation not indicated in _____       | R4 Use vial with smallest bubble first  |
| D5 Container -[improper] [leaking] [broken]     | D17 Preservation mismatch COC vs label        | R5 Log-in with latest sampling date and time+1 min  |
| D6 Date/Time is not indicated in _____          | D18 Insufficient chemical preservative        | R6 Adjust pH as necessary   |
| <b>D7</b> Date/Time mismatch COC vs label       | D19 Insufficient Sample                       | R7 Filter and preserved as necessary  |
| D8 Sample listed in COC is not received         | D20 No filtration info for dissolved analysis | R8 <i>Informed client</i>   |
| D9 Sample received is not listed in COC         | D21 No sample for moisture determination      | R9 _____  |
| D10 No initial/date on corrections in COC/label | D22 _____                                     | R10 _____   |
| D11 Container count mismatch COC vs received    | D23 _____                                     | R11 _____   |
| D12 Container size mismatch COC vs received     | D24 _____                                     | R12 _____   |

**REVIEWS:**

Sample Labeling: *Jocelyne Solis-Ramos* / *[Signature]* SRF: *[Signature]*  
 Date: *12/22/22* Date: *12/22/22*

PM: *[Signature]*  
 Date: *12/28/22*



## REPORTING CONVENTIONS

### DATA QUALIFIERS:

| Lab Qualifier | AFCEE Qualifier | Description  |
|---------------|-----------------|--|
| J             | F               | Indicates that the analyte is positively identified and the result is less than RL but greater than MDL.       |
| N             |                 | Indicates presumptive evidence of a compound.  |
| B             | B               | Indicates that the analyte is found in the associated method blank as well as in the sample at above QC level. |
| E             | J               | Indicates that the result is above the maximum calibration range or estimated value.                           |
| *             | *               | Out of QC limit.   |

**Note:** The above qualifiers are used to flag the results unless the project requires a different set of qualification criteria.

### ACRONYMS AND ABBREVIATIONS:

|      |                                   |
|------|-----------------------------------|
| CRDL | Contract Required Detection Limit |
| RL   | Reporting Limit                   |
| MRL  | Method Reporting Limit            |
| PQL  | Practical Quantitation Limit      |
| MDL  | Method Detection Limit            |
| DO   | Diluted out                       |

### DATES

The date and time information for leaching and preparation reflect the beginning date and time of the procedure unless the method, protocol, or project specifically requires otherwise.

LABORATORY REPORT FOR

EUROFINS EATON ANALYTICAL

380-32096

METHOD 5030B/8015B  
TOTAL PETROLEUM HYDROCARBONS BY PURGE AND TRAP

SDG#: 22L294

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CASE NARRATIVE

Client : EUROFINS EATON ANALYTICAL

Project: 380-32096

SDG : 22L294

METHOD 5030B/8015B  
TOTAL PETROLEUM HYDROCARBONS BY PURGE AND TRAP

A total of six(6) water samples were received on 12/22/22 to be analyzed for Total Petroleum Hydrocarbons by Purge And Trap in accordance with Method 5030B/8015B and project specific requirements.

Holding Time

Samples were analyzed within the prescribed holding time.

Calibration

Multi-calibration points were generated to establish initial calibration (ICAL). ICAL was verified using a secondary source (ICV). Continuing calibration (CCV) verifications were carried out on a frequency specified by the project. All calibration requirements were within acceptance criteria. Refer to calibration summary forms of ICAL, ICV and CCV for details. MRL was analyzed as required by the project.

Method Blank

Method blank was prepared and analyzed at the frequency required by the project. For this SDG, one(1) method blank was analyzed. VG39L08B - result was compliant to project requirement. Refer to sample result summary form for details.

Lab Control Sample

Lab control sample was prepared and analyzed at a frequency required by the project. For this SDG, one(1) set of LCS/LCD was analyzed. VG39L08L/VG39L08C were within LCS limits. Refer to LCS summary form for details.

Matrix QC Sample

Matrix spike sample was prepared and analyzed at a frequency required by the project. For this SDG, one(1) set of MS/MSD was analyzed. Gasoline was within MS QC limits in L294-01M/L294-01S. Refer to Matrix QC summary form for details.

Surrogate

Surrogate was added on QC and field samples. All surrogate recoveries were within QC limits. Refer to sample result summary forms for details.

Sample Analysis

Samples were analyzed according to prescribed analytical procedures. Results were evaluated in accordance to project requirements. For this SDG, all quality control requirements were met.



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# SAMPLE RESULTS





METHOD 5030B/8015B  
TOTAL PETROLEUM HYDROCARBONS BY PURGE AND TRAP

|                                    |                                |
|------------------------------------|--------------------------------|
| Client : EUROFINS EATON ANALYTICAL | Date Collected: 12/19/22 11:18 |
| Project : 380-32096                | Date Received: 12/22/22        |
| Batch No. : 22L294                 | Date Extracted: 12/22/22 21:33 |
| Sample ID : 380-32096-3            | Date Analyzed: 12/22/22 21:33  |
| Lab Samp ID: L294-03               | Dilution Factor: 1             |
| Lab File ID: EL22021A              | Matrix: WATER                  |
| Ext Btch ID: 22VG39L08             | % Moisture: NA                 |
| Calib. Ref.: EL22015A              | Instrument ID: 39              |

| PARAMETERS                  | RESULTS<br>(mg/L) | RL<br>(mg/L) | MDL<br>(mg/L) |          |
|-----------------------------|-------------------|--------------|---------------|----------|
| -----<br>GASOLINE           | ND                | 0.020        | 0.010         |          |
| SURROGATE PARAMETERS        | RESULT            | SPK_AMT      | %RECOVERY     | QC LIMIT |
| -----<br>Bromofluorobenzene | 0.0322            | 0.0400       | 81            | 60-140   |

Notes:

Parameter      H-C Range  
Gasoline        C6-C10

Reported ND at RL quantitated per pattern recognition.

Detection limits are reported relative to sample result significant figures.

|                      |                      |
|----------------------|----------------------|
| Sample Amount : 5ml  | Final Volume : 5ml   |
| Prepared by : SCerva | Analyzed by : SCerva |







METHOD 5030B/8015B  
TOTAL PETROLEUM HYDROCARBONS BY PURGE AND TRAP

|              |                             |                                |
|--------------|-----------------------------|--------------------------------|
| Client       | : EUROFINS EATON ANALYTICAL | Date Collected: 12/19/22 11:18 |
| Project      | : 380-32096                 | Date Received: 12/22/22        |
| Batch No.    | : 22L294                    | Date Extracted: 12/22/22 23:20 |
| Sample ID    | : 380-32096-6               | Date Analyzed: 12/22/22 23:20  |
| Lab Samp ID: | L294-06                     | Dilution Factor: 1             |
| Lab File ID: | EL22024A                    | Matrix: WATER                  |
| Ext Btch ID: | 22VG39L08                   | % Moisture: NA                 |
| Calib. Ref.: | EL22015A                    | Instrument ID: 39              |

| PARAMETERS | RESULTS<br>(mg/L) | RL<br>(mg/L) | MDL<br>(mg/L) |
|------------|-------------------|--------------|---------------|
| GASOLINE   | ND                | 0.020        | 0.010         |

| SURROGATE PARAMETERS | RESULT | SPK_AMT | %RECOVERY | QC LIMIT |
|----------------------|--------|---------|-----------|----------|
| Bromofluorobenzene   | 0.0321 | 0.0400  | 80        | 60-140   |

Notes:

Parameter      H-C Range  
Gasoline        C6-C10

Reported ND at RL quantitated per pattern recognition.

Detection limits are reported relative to sample result significant figures.

|                      |                      |
|----------------------|----------------------|
| Sample Amount : 5ml  | Final Volume : 5ml   |
| Prepared by : SCerva | Analyzed by : SCerva |

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# QC SUMMARIES



EMAX QUALITY CONTROL DATA  
LAB CONTROL SAMPLE ANALYSIS

CLIENT : EUROFINS EATON ANALYTICAL  
PROJECT : 380-32096  
BATCH NO. : 22L294  
METHOD : 5030B/8015B

|                  |                  |                |                |
|------------------|------------------|----------------|----------------|
| MATRIX           | : WATER          |                | % MOISTURE:NA  |
| DILUTION FACTOR: | 1                | 1              | 1              |
| SAMPLE ID        | : MBLK1W         | LCS1W          | LCD1W          |
| LAB SAMPLE ID    | : VG39L08B       | VG39L08L       | VG39L08C       |
| LAB FILE ID      | : EL22005A       | EL22006A       | EL22007A       |
| DATE PREPARED    | : 12/22/22 11:58 | 12/22/22 12:34 | 12/22/22 13:09 |
| DATE ANALYZED    | : 12/22/22 11:58 | 12/22/22 12:34 | 12/22/22 13:09 |
| PREP BATCH       | : 22VG39L08      | 22VG39L08      | 22VG39L08      |
| CALIBRATION REF: | EL22004A         | EL22004A       | EL22004A       |

ACCESSION:

| PARAMETERS | MBResult<br>(mg/L) | SpikeAmt<br>(mg/L) | LCSResult<br>(mg/L) | LCSRec<br>(%) | SpikeAmt<br>(mg/L) | LCDResult<br>(mg/L) | LCDRec<br>(%) | RPD<br>(%) | QCLimit<br>(%) | MaxRPD<br>(%) |
|------------|--------------------|--------------------|---------------------|---------------|--------------------|---------------------|---------------|------------|----------------|---------------|
| Gasoline   | ND                 | 0.500              | 0.471               | 94            | 0.500              | 0.474               | 95            | 1          | 60-130         | 30            |

| SURROGATE PARAMETER | SpikeAmt<br>(mg/L) | LCSResult<br>(mg/L) | LCSRec<br>(%) | SpikeAmt<br>(mg/L) | LCDResult<br>(mg/L) | LCDRec<br>(%) | QCLimit<br>(%) |
|---------------------|--------------------|---------------------|---------------|--------------------|---------------------|---------------|----------------|
| Bromofluorobenzene  | 0.0400             | 0.0440              | 110           | 0.0400             | 0.0434              | 109           | 70-130         |

MB: Method Blank sample LCS: Lab Control Sample LCD: Lab Control Sample Duplicate

EMAX QUALITY CONTROL DATA  
MS/MSD ANALYSIS

CLIENT : EUROFINS EATON ANALYTICAL  
PROJECT : 380-32096  
BATCH NO. : 22L294  
METHOD : 5030B/8015B

|                                |                |                |
|--------------------------------|----------------|----------------|
| MATRIX : WATER                 |                | % MOISTURE:NA  |
| DILUTION FACTOR: 1             | 1              | 1              |
| SAMPLE ID : 380-32096-1        | 380-32096-1MS  | 380-32096-1MSD |
| LAB SAMPLE ID : L294-01        | L294-01M       | L294-01S       |
| LAB FILE ID : EL22017A         | EL22018A       | EL22019A       |
| DATE PREPARED : 12/22/22 19:09 | 12/22/22 19:45 | 12/22/22 20:21 |
| DATE ANALYZED : 12/22/22 19:09 | 12/22/22 19:45 | 12/22/22 20:21 |
| PREP BATCH : 22VG39L08         | 22VG39L08      | 22VG39L08      |
| CALIBRATION REF: EL22015A      | EL22015A       | EL22015A       |

ACCESSION:

| PARAMETERS | PSResult<br>(mg/L) | SpikeAmt<br>(mg/L) | MSResult<br>(mg/L) | MSRec<br>(%) | SpikeAmt<br>(mg/L) | MSDResult<br>(mg/L) | MSDRec<br>(%) | RPD<br>(%) | QCLimit<br>(%) | MaxRPD<br>(%) |
|------------|--------------------|--------------------|--------------------|--------------|--------------------|---------------------|---------------|------------|----------------|---------------|
| Gasoline   | ND                 | 0.500              | 0.494              | 99           | 0.500              | 0.484               | 97            | 2          | 50-130         | 30            |

| SURROGATE PARAMETER | SpikeAmt<br>(mg/L) | MSResult<br>(mg/L) | MSRec<br>(%) | SpikeAmt<br>(mg/L) | MSDResult<br>(mg/L) | MSDRec<br>(%) | QCLimit<br>(%) |
|---------------------|--------------------|--------------------|--------------|--------------------|---------------------|---------------|----------------|
| Bromofluorobenzene  | 0.0400             | 0.0432             | 108          | 0.0400             | 0.0441              | 110           | 60-140         |

PS: Parent Sample MS: Matrix Spike MSD: Matrix Spike Duplicate

LABORATORY REPORT FOR

EUROFINS EATON ANALYTICAL

380-32096

METHOD 3520C/8015B  
TOTAL PETROLEUM HYDROCARBONS BY EXTRACTION

SDG#: 22L294

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CASE NARRATIVE

Client : EUROFINS EATON ANALYTICAL

Project: 380-32096

SDG : 22L294

METHOD 3520C/8015B  
TOTAL PETROLEUM HYDROCARBONS BY EXTRACTION

A total of three(3) water samples were received on 12/22/22 to be analyzed for Total Petroleum Hydrocarbons by Extraction in accordance with Method 3520C/8015B and project specific requirements.

Holding Time

Samples were analyzed within the prescribed holding time.

Calibration

Multi-calibration points were generated to establish initial calibration (ICAL). ICAL was verified using a secondary source (ICV). Continuing calibration (CCV) verifications were carried out on a frequency specified by the project. All calibration requirements were within acceptance criteria. Refer to calibration summary forms of ICAL, ICV and CCV for details. MRL was analyzed as required by the project.

Method Blank

Method blank was prepared and analyzed at the frequency required by the project. For this SDG, one(1) method blank was analyzed. DSL040WB - result was compliant to project requirement. Refer to sample result summary form for details.

Lab Control Sample

Lab control sample was prepared and analyzed at a frequency required by the project. For this SDG, one(1) LCS was analyzed. Percent recovery for Diesel was within LCS QC limits in DSL040WL. Refer to LCS summary form for details.

Matrix QC Sample

Matrix spike sample was prepared and analyzed at a frequency required by the project. For this SDG, one(1) set of MS/MSD was analyzed. Diesel was within MS QC limits in 22L302-01M/22L302-01S. Refer to Matrix QC summary form for details.

Surrogate

Surrogates were added on QC and field samples. All surrogate recoveries were within QC limits. Refer to sample result summary forms for details.

Sample Analysis

Samples were analyzed according to prescribed analytical procedures. Results were evaluated in accordance to project requirements. For this SDG, all quality control requirements were met.

CASE NARRATIVE

Client : EUROFINS EATON ANALYTICAL

Project: 380-32096

SDG : 22L294

METHOD 3520C/8015B  
PETROLEUM HYDROCARBONS BY EXTRACTION

A total of three(3) water samples were received on 12/22/22 to be analyzed for Petroleum Hydrocarbons by Extraction in accordance with Method 3520C/8015B and project specific requirements.

Holding Time

Samples were analyzed within the prescribed holding time.

Calibration

Multi-calibration points were generated to establish initial calibration (ICAL). ICAL was verified using a secondary source (ICV). Continuing calibration (CCV) verifications were carried out on a frequency specified by the project. All calibration requirements were within acceptance criteria. Refer to calibration summary forms of ICAL, ICV and CCV for details. MRL was analyzed as required by the project.

Method Blank

Method blank was prepared and analyzed at the frequency required by the project. For this SDG, one(1) method blank was analyzed. DSL040WB - result was compliant to project requirement. Refer to sample result summary form for details.

Lab Control Sample

Lab control sample was prepared and analyzed at a frequency required by the project. For this SDG, one(1) LCS was analyzed. Percent recovery for JP5 was within LCS QC limits in J5L040WL. Refer to LCS summary form for details.

Matrix QC Sample

Matrix spike sample was prepared and analyzed at a frequency required by the project. For this SDG, one(1) set of MS/MSD was analyzed. JP5 was within MS QC limits in 22L302-01M/22L302-01S. Refer to Matrix QC summary form for details.

Surrogate

Surrogates were added on QC and field samples. All surrogate recoveries were within QC limits. Refer to sample result summary forms for details.

Sample Analysis

Samples were analyzed according to prescribed analytical procedures. Results were evaluated in accordance to project requirements. For this SDG, all quality control requirements were met.

CASE NARRATIVE

Client : EUROFINS EATON ANALYTICAL

Project: 380-32096

SDG : 22L294

METHOD 3520C/8015B  
PETROLEUM HYDROCARBONS BY EXTRACTION

A total of three(3) water samples were received on 12/22/22 to be analyzed for Petroleum Hydrocarbons by Extraction in accordance with Method 3520C/8015B and project specific requirements.

Holding Time

Samples were analyzed within the prescribed holding time.

Calibration

Multi-calibration points were generated to establish initial calibration (ICAL). ICAL was verified using a secondary source (ICV). Continuing calibration (CCV) verifications were carried out on a frequency specified by the project. All calibration requirements were within acceptance criteria. Refer to calibration summary forms of ICAL, ICV and CCV for details. MRL was analyzed as required by the project.

Method Blank

Method blank was prepared and analyzed at the frequency required by the project. For this SDG, one(1) method blank was analyzed. DSL040WB - result was compliant to project requirement. Refer to sample result summary form for details.

Lab Control Sample

Lab control sample was prepared and analyzed at a frequency required by the project. For this SDG, one(1) LCS was analyzed. Percent recovery for JP8 was within LCS QC limits in J8L040WL. Refer to LCS summary form for details.

Matrix QC Sample

Matrix spike sample was prepared and analyzed at a frequency required by the project. For this SDG, one(1) set of MS/MSD was analyzed. JP8 was within MS QC limits in 22L303-01M/22L303-01S. Refer to Matrix QC summary form for details.

Surrogate

Surrogates were added on QC and field samples. All surrogate recoveries were within QC limits. Refer to sample result summary forms for details.

Sample Analysis

Samples were analyzed according to prescribed analytical procedures. Results were evaluated in accordance to project requirements. For this SDG, all quality control requirements were met.







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# SAMPLE RESULTS

METHOD 3520C/8015B  
TOTAL PETROLEUM HYDROCARBONS BY EXTRACTION

|              |                             |                  |                |
|--------------|-----------------------------|------------------|----------------|
| Client       | : EUROFINS EATON ANALYTICAL | Date Collected:  | 12/19/22 10:00 |
| Project      | : 380-32096                 | Date Received:   | 12/22/22       |
| Batch No.    | : 22L294                    | Date Extracted:  | 12/28/22 13:15 |
| Sample ID    | : 380-32096-1               | Date Analyzed:   | 01/05/23 21:25 |
| Lab Samp ID: | 22L294-01                   | Dilution Factor: | 1              |
| Lab File ID: | LA05030A                    | Matrix:          | WATER          |
| Ext Btch ID: | 22DSL040W                   | % Moisture:      | NA             |
| Calib. Ref.: | LA05020A                    | Instrument ID:   | D5             |

| PARAMETERS | RESULTS<br>(mg/L) | RL<br>(mg/L) | MDL<br>(mg/L) |
|------------|-------------------|--------------|---------------|
| Diesel     | ND                | 0.026        | 0.013         |
| Motor Oil  | ND                | 0.052        | 0.026         |

| SURROGATE PARAMETERS | RESULT | SPK_AMT | %RECOVERY | QC LIMIT |
|----------------------|--------|---------|-----------|----------|
| Bromobenzene         | 0.490  | 0.525   | 93        | 60-130   |
| Hexacosane           | 0.150  | 0.131   | 115       | 60-130   |

Notes:

Parameter H-C Range  
Diesel C10-C24  
Motor Oil C24-C36

Reported ND at RL quantitated per pattern recognition.

Detection limits are reported relative to sample result significant figures.

Sample Amount : 950ml Final Volume : 5ml  
Prepared by : P0reto Analyzed by : SDeeso



METHOD 3520C/8015B  
 PETROLEUM HYDROCARBONS BY EXTRACTION

|              |                             |                  |                |
|--------------|-----------------------------|------------------|----------------|
| Client       | : EUROFINS EATON ANALYTICAL | Date Collected:  | 12/19/22 10:00 |
| Project      | : 380-32096                 | Date Received:   | 12/22/22       |
| Batch No.    | : 22L294                    | Date Extracted:  | 12/28/22 13:15 |
| Sample ID    | : 380-32096-1               | Date Analyzed:   | 01/05/23 21:25 |
| Lab Samp ID: | 22L294-01                   | Dilution Factor: | 1              |
| Lab File ID: | LA05030A                    | Matrix:          | WATER          |
| Ext Btch ID: | 22DSL040W                   | % Moisture:      | NA             |
| Calib. Ref.: | LA05021A                    | Instrument ID:   | D5             |

| PARAMETERS | RESULTS<br>(mg/L) | RL<br>(mg/L) | MDL<br>(mg/L) |
|------------|-------------------|--------------|---------------|
| JP5        | ND                | 0.052        | 0.026         |

| SURROGATE PARAMETERS | RESULT | SPK_AMT | %RECOVERY | QC LIMIT |
|----------------------|--------|---------|-----------|----------|
| Bromobenzene         | 0.490  | 0.525   | 93        | 60-130   |
| Hexacosane           | 0.150  | 0.131   | 115       | 60-130   |

Notes:

RL : Reporting Limit

Parameter H-C Range

JP5 C8-C18

Reported ND at RL quantitated per pattern recognition.

Detection limits are reported relative to sample result significant figures.

Sample Amount : 950ml

Final Volume : 5ml

Prepared by : P0reto

Analyzed by : SDeeso

METHOD 3520C/8015B  
 PETROLEUM HYDROCARBONS BY EXTRACTION

|              |                             |                  |                |
|--------------|-----------------------------|------------------|----------------|
| Client       | : EUROFINS EATON ANALYTICAL | Date Collected:  | 12/19/22 10:00 |
| Project      | : 380-32096                 | Date Received:   | 12/22/22       |
| Batch No.    | : 22L294                    | Date Extracted:  | 12/28/22 13:15 |
| Sample ID    | : 380-32096-1               | Date Analyzed:   | 01/05/23 21:25 |
| Lab Samp ID: | 22L294-01                   | Dilution Factor: | 1              |
| Lab File ID: | LA05030A                    | Matrix:          | WATER          |
| Ext Btch ID: | 22DSL040W                   | % Moisture:      | NA             |
| Calib. Ref.: | LA05022A                    | Instrument ID:   | D5             |

| PARAMETERS           | RESULTS<br>(mg/L) | RL<br>(mg/L) | MDL<br>(mg/L) |          |
|----------------------|-------------------|--------------|---------------|----------|
| JP8                  | ND                | 0.052        | 0.026         |          |
| SURROGATE PARAMETERS | RESULT            | SPK_AMT      | %RECOVERY     | QC LIMIT |
| Bromobenzene         | 0.490             | 0.525        | 93            | 60-130   |
| Hexacosane           | 0.150             | 0.131        | 115           | 60-130   |

Notes:

RL : Reporting Limit

Parameter H-C Range

JP8 C8-C18

Reported ND at RL quantitated per pattern recognition.

Detection limits are reported relative to sample result significant figures.

Sample Amount : 950ml

Final Volume : 5ml

Prepared by : P0reto

Analyzed by : SDeeso

METHOD 3520C/8015B  
TOTAL PETROLEUM HYDROCARBONS BY EXTRACTION

```

Client      : EUROFINS EATON ANALYTICAL   Date Collected: 12/19/22 10:45
Project     : 380-32096                   Date Received: 12/22/22
Batch No.   : 22L294                       Date Extracted: 12/28/22 13:15
Sample ID   : 380-32096-2                 Date Analyzed: 01/05/23 21:44
Lab Samp ID: 22L294-02                    Dilution Factor: 1
Lab File ID: LA05031A                     Matrix: WATER
Ext Btch ID: 22DSL040W                    % Moisture: NA
Calib. Ref.: LA05020A                     Instrument ID: D5
  
```

| PARAMETERS           | RESULTS<br>(mg/L) | RL<br>(mg/L) | MDL<br>(mg/L) |          |
|----------------------|-------------------|--------------|---------------|----------|
| Diesel               | ND                | 0.027        | 0.013         |          |
| Motor Oil            | ND                | 0.053        | 0.027         |          |
| SURROGATE PARAMETERS | RESULT            | SPK_AMT      | %RECOVERY     | QC LIMIT |
| Bromobenzene         | 0.443             | 0.530        | 84            | 60-130   |
| Hexacosane           | 0.147             | 0.132        | 111           | 60-130   |

Notes:

Parameter H-C Range  
Diesel C10-C24  
Motor Oil C24-C36

Reported ND at RL quantitated per pattern recognition.

Detection limits are reported relative to sample result significant figures.

Sample Amount : 940ml                      Final Volume : 5ml  
Prepared by : P0reto                         Analyzed by : SDeeso

METHOD 3520C/8015B  
 PETROLEUM HYDROCARBONS BY EXTRACTION

|              |                             |                  |                |
|--------------|-----------------------------|------------------|----------------|
| Client       | : EUROFINS EATON ANALYTICAL | Date Collected:  | 12/19/22 10:45 |
| Project      | : 380-32096                 | Date Received:   | 12/22/22       |
| Batch No.    | : 22L294                    | Date Extracted:  | 12/28/22 13:15 |
| Sample ID    | : 380-32096-2               | Date Analyzed:   | 01/05/23 21:44 |
| Lab Samp ID: | 22L294-02                   | Dilution Factor: | 1              |
| Lab File ID: | LA05031A                    | Matrix:          | WATER          |
| Ext Btch ID: | 22DSL040W                   | % Moisture:      | NA             |
| Calib. Ref.: | LA05021A                    | Instrument ID:   | D5             |

| PARAMETERS | RESULTS<br>(mg/L) | RL<br>(mg/L) | MDL<br>(mg/L) |
|------------|-------------------|--------------|---------------|
| JP5        | ND                | 0.053        | 0.027         |

| SURROGATE PARAMETERS | RESULT | SPK_AMT | %RECOVERY | QC LIMIT |
|----------------------|--------|---------|-----------|----------|
| Bromobenzene         | 0.443  | 0.530   | 84        | 60-130   |
| Hexacosane           | 0.147  | 0.132   | 111       | 60-130   |

Notes:

RL : Reporting Limit

Parameter H-C Range

JP5 C8-C18

Reported ND at RL quantitated per pattern recognition.

Detection limits are reported relative to sample result significant figures.

Sample Amount : 940ml

Final Volume : 5ml

Prepared by : P0reto

Analyzed by : SDeeso

METHOD 3520C/8015B  
 PETROLEUM HYDROCARBONS BY EXTRACTION

|              |                             |                  |                |
|--------------|-----------------------------|------------------|----------------|
| Client       | : EUROFINS EATON ANALYTICAL | Date Collected:  | 12/19/22 10:45 |
| Project      | : 380-32096                 | Date Received:   | 12/22/22       |
| Batch No.    | : 22L294                    | Date Extracted:  | 12/28/22 13:15 |
| Sample ID    | : 380-32096-2               | Date Analyzed:   | 01/05/23 21:44 |
| Lab Samp ID: | 22L294-02                   | Dilution Factor: | 1              |
| Lab File ID: | LA05031A                    | Matrix:          | WATER          |
| Ext Btch ID: | 22DSL040W                   | % Moisture:      | NA             |
| Calib. Ref.: | LA05022A                    | Instrument ID:   | D5             |

| PARAMETERS | RESULTS<br>(mg/L) | RL<br>(mg/L) | MDL<br>(mg/L) |  |
|------------|-------------------|--------------|---------------|--|
| JP8        | ND                | 0.053        | 0.027         |  |

| SURROGATE PARAMETERS | RESULT | SPK_AMT | %RECOVERY | QC LIMIT |
|----------------------|--------|---------|-----------|----------|
| Bromobenzene         | 0.443  | 0.530   | 84        | 60-130   |
| Hexacosane           | 0.147  | 0.132   | 111       | 60-130   |

Notes:

RL : Reporting Limit

Parameter H-C Range

JP8 C8-C18

Reported ND at RL quantitated per pattern recognition.

Detection limits are reported relative to sample result significant figures.

Sample Amount : 940ml

Final Volume : 5ml

Prepared by : P0reto

Analyzed by : SDeeso



METHOD 3520C/8015B  
 PETROLEUM HYDROCARBONS BY EXTRACTION

|              |                             |                  |                |
|--------------|-----------------------------|------------------|----------------|
| Client       | : EUROFINS EATON ANALYTICAL | Date Collected:  | 12/19/22 11:18 |
| Project      | : 380-32096                 | Date Received:   | 12/22/22       |
| Batch No.    | : 22L294                    | Date Extracted:  | 12/28/22 13:15 |
| Sample ID    | : 380-32096-3               | Date Analyzed:   | 01/05/23 22:02 |
| Lab Samp ID: | 22L294-03                   | Dilution Factor: | 1              |
| Lab File ID: | LA05032A                    | Matrix:          | WATER          |
| Ext Btch ID: | 22DSL040W                   | % Moisture:      | NA             |
| Calib. Ref.: | LA05021A                    | Instrument ID:   | D5             |

| PARAMETERS | RESULTS<br>(mg/L) | RL<br>(mg/L) | MDL<br>(mg/L) |
|------------|-------------------|--------------|---------------|
| JP5        | ND                | 0.053        | 0.027         |

| SURROGATE PARAMETERS | RESULT | SPK_AMT | %RECOVERY | QC LIMIT |
|----------------------|--------|---------|-----------|----------|
| Bromobenzene         | 0.440  | 0.530   | 83        | 60-130   |
| Hexacosane           | 0.128  | 0.132   | 97        | 60-130   |

Notes:

RL : Reporting Limit

Parameter H-C Range

JP5 C8-C18

Reported ND at RL quantitated per pattern recognition.

Detection limits are reported relative to sample result significant figures.

Sample Amount : 940ml

Final Volume : 5ml

Prepared by : P0reto

Analyzed by : SDeeso

METHOD 3520C/8015B  
 PETROLEUM HYDROCARBONS BY EXTRACTION

|              |                             |                  |                |
|--------------|-----------------------------|------------------|----------------|
| Client       | : EUROFINS EATON ANALYTICAL | Date Collected:  | 12/19/22 11:18 |
| Project      | : 380-32096                 | Date Received:   | 12/22/22       |
| Batch No.    | : 22L294                    | Date Extracted:  | 12/28/22 13:15 |
| Sample ID    | : 380-32096-3               | Date Analyzed:   | 01/05/23 22:02 |
| Lab Samp ID: | 22L294-03                   | Dilution Factor: | 1              |
| Lab File ID: | LA05032A                    | Matrix:          | WATER          |
| Ext Btch ID: | 22DSL040W                   | % Moisture:      | NA             |
| Calib. Ref.: | LA05022A                    | Instrument ID:   | D5             |

| PARAMETERS | RESULTS<br>(mg/L) | RL<br>(mg/L) | MDL<br>(mg/L) |  |
|------------|-------------------|--------------|---------------|--|
| JP8        | ND                | 0.053        | 0.027         |  |

| SURROGATE PARAMETERS | RESULT | SPK_AMT | %RECOVERY | QC LIMIT |
|----------------------|--------|---------|-----------|----------|
| Bromobenzene         | 0.440  | 0.530   | 83        | 60-130   |
| Hexacosane           | 0.128  | 0.132   | 97        | 60-130   |

Notes:

RL : Reporting Limit  
 Parameter H-C Range  
 JP8 C8-C18  
 Reported ND at RL quantitated per pattern recognition.

Detection limits are reported relative to sample result significant figures.

Sample Amount : 940ml Final Volume : 5ml  
 Prepared by : P0reto Analyzed by : SDeeso



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# QC SUMMARIES

METHOD 3520C/8015B  
TOTAL PETROLEUM HYDROCARBONS BY EXTRACTION

|              |                             |                  |                |
|--------------|-----------------------------|------------------|----------------|
| Client       | : EUROFINS EATON ANALYTICAL | Date Collected:  | 12/28/22 13:15 |
| Project      | : 380-32096                 | Date Received:   | 12/28/22       |
| Batch No.    | : 22L294                    | Date Extracted:  | 12/28/22 13:15 |
| Sample ID    | : MBLK1W                    | Date Analyzed:   | 01/05/23 20:12 |
| Lab Samp ID: | DSL040WB                    | Dilution Factor: | 1              |
| Lab File ID: | LA05026A                    | Matrix:          | WATER          |
| Ext Btch ID: | 22DSL040W                   | % Moisture:      | NA             |
| Calib. Ref.: | LA05020A                    | Instrument ID:   | D5             |

| PARAMETERS           | RESULTS<br>(mg/L) | RL<br>(mg/L) | MDL<br>(mg/L) |          |
|----------------------|-------------------|--------------|---------------|----------|
| Diesel               | ND                | 0.025        | 0.012         |          |
| Motor Oil            | ND                | 0.050        | 0.025         |          |
| SURROGATE PARAMETERS | RESULT            | SPK_AMT      | %RECOVERY     | QC LIMIT |
| Bromobenzene         | 0.406             | 0.500        | 81            | 60-130   |
| Hexacosane           | 0.131             | 0.125        | 105           | 60-130   |

Notes:

Parameter H-C Range  
Diesel C10-C24  
Motor Oil C24-C36

Reported ND at RL quantitated per pattern recognition.

Detection limits are reported relative to sample result significant figures.

Sample Amount : 1000ml Final Volume : 5ml  
Prepared by : P0reto Analyzed by : SDeeso

EMAX QUALITY CONTROL DATA  
LAB CONTROL SAMPLE ANALYSIS

CLIENT : EUROFINS EATON ANALYTICAL  
PROJECT : 380-32096  
BATCH NO. : 22L294  
METHOD : 3520C/8015B

MATRIX : WATER % MOISTURE:NA  
DILUTION FACTOR: 1 1  
SAMPLE ID : MBLK1W LCS1W  
LAB SAMPLE ID : DSL040WB DSL040WL  
LAB FILE ID : LA05026A LA05027A  
DATE PREPARED : 12/28/22 13:15 12/28/22 13:15  
DATE ANALYZED : 01/05/23 20:12 01/05/23 20:30  
PREP BATCH : 22DSL040W 22DSL040W  
CALIBRATION REF: LA05020A LA05020A

ACCESSION:

| PARAMETERS | MBResult<br>(mg/L) | SpikeAmt<br>(mg/L) | LCSResult<br>(mg/L) | LCSRec<br>(%) | QCLimit<br>(%) |
|------------|--------------------|--------------------|---------------------|---------------|----------------|
| Diesel     | ND                 | 2.50               | 2.83                | 113           | 50-130         |

| SURROGATE PARAMETERS | SpikeAmt<br>(mg/L) | LCSResult<br>(mg/L) | LCSRec<br>(%) | QCLimit<br>(%) |
|----------------------|--------------------|---------------------|---------------|----------------|
| Bromobenzene         | 0.500              | 0.510               | 102           | 60-130         |
| Hexacosane           | 0.125              | 0.132               | 106           | 60-130         |

MB: Method Blank sample LCS: Lab Control Sample

METHOD 3520C/8015B  
 PETROLEUM HYDROCARBONS BY EXTRACTION

|              |                             |                  |                |
|--------------|-----------------------------|------------------|----------------|
| Client       | : EUROFINS EATON ANALYTICAL | Date Collected:  | 12/28/22 13:15 |
| Project      | : 380-32096                 | Date Received:   | 12/28/22       |
| Batch No.    | : 22L294                    | Date Extracted:  | 12/28/22 13:15 |
| Sample ID    | : MBLK1W                    | Date Analyzed:   | 01/05/23 20:12 |
| Lab Samp ID: | DSL040WB                    | Dilution Factor: | 1              |
| Lab File ID: | LA05026A                    | Matrix:          | WATER          |
| Ext Btch ID: | 22DSL040W                   | % Moisture:      | NA             |
| Calib. Ref.: | LA05021A                    | Instrument ID:   | D5             |

| PARAMETERS           | RESULTS<br>(mg/L) | RL<br>(mg/L) | MDL<br>(mg/L) |          |
|----------------------|-------------------|--------------|---------------|----------|
| JP5                  | ND                | 0.050        | 0.025         |          |
| SURROGATE PARAMETERS | RESULT            | SPK_AMT      | %RECOVERY     | QC LIMIT |
| Bromobenzene         | 0.406             | 0.500        | 81            | 60-130   |
| Hexacosane           | 0.131             | 0.125        | 105           | 60-130   |

Notes:

RL : Reporting Limit

Parameter H-C Range

JP5 C8-C18

Reported ND at RL quantitated per pattern recognition.

Detection limits are reported relative to sample result significant figures.

Sample Amount : 1000ml

Final Volume : 5ml

Prepared by : P0reto

Analyzed by : SDeeso

EMAX QUALITY CONTROL DATA  
LAB CONTROL SAMPLE ANALYSIS

CLIENT : EUROFINS EATON ANALYTICAL  
PROJECT : 380-32096  
BATCH NO. : 22L294  
METHOD : 3520C/8015B

MATRIX : WATER % MOISTURE:NA  
DILUTION FACTOR: 1 1  
SAMPLE ID : MBLK1W LCS1W  
LAB SAMPLE ID : DSL040WB J5L040WL  
LAB FILE ID : LA05026A LA05028A  
DATE PREPARED : 12/28/22 13:15 12/28/22 13:15  
DATE ANALYZED : 01/05/23 20:12 01/05/23 20:49  
PREP BATCH : 22DSL040W 22DSL040W  
CALIBRATION REF: LA05021A LA05021A

ACCESSION:

| PARAMETERS | MBResult<br>(mg/L) | SpikeAmt<br>(mg/L) | LCSResult<br>(mg/L) | LCSRec<br>(%) | QCLimit<br>(%) |
|------------|--------------------|--------------------|---------------------|---------------|----------------|
| JP5        | ND                 | 2.50               | 2.43                | 97            | 30-160         |

| SURROGATE PARAMETERS | SpikeAmt<br>(mg/L) | LCSResult<br>(mg/L) | LCSRec<br>(%) | QCLimit<br>(%) |
|----------------------|--------------------|---------------------|---------------|----------------|
| Bromobenzene         | 0.500              | 0.523               | 105           | 60-130         |
| Hexacosane           | 0.125              | 0.121               | 97            | 60-130         |

MB: Method Blank sample LCS: Lab Control Sample

METHOD 3520C/8015B  
 PETROLEUM HYDROCARBONS BY EXTRACTION

|              |                             |                  |                |
|--------------|-----------------------------|------------------|----------------|
| Client       | : EUROFINS EATON ANALYTICAL | Date Collected:  | 12/28/22 13:15 |
| Project      | : 380-32096                 | Date Received:   | 12/28/22       |
| Batch No.    | : 22L294                    | Date Extracted:  | 12/28/22 13:15 |
| Sample ID    | : MBLK1W                    | Date Analyzed:   | 01/05/23 20:12 |
| Lab Samp ID: | DSL040WB                    | Dilution Factor: | 1              |
| Lab File ID: | LA05026A                    | Matrix:          | WATER          |
| Ext Btch ID: | 22DSL040W                   | % Moisture:      | NA             |
| Calib. Ref.: | LA05022A                    | Instrument ID:   | D5             |

| PARAMETERS | RESULTS<br>(mg/L) | RL<br>(mg/L) | MDL<br>(mg/L) |
|------------|-------------------|--------------|---------------|
| JP8        | ND                | 0.050        | 0.025         |

| SURROGATE PARAMETERS | RESULT | SPK_AMT | %RECOVERY | QC LIMIT |
|----------------------|--------|---------|-----------|----------|
| Bromobenzene         | 0.406  | 0.500   | 81        | 60-130   |
| Hexacosane           | 0.131  | 0.125   | 105       | 60-130   |

Notes:

RL : Reporting Limit  
 Parameter H-C Range  
 JP8 C8-C18  
 Reported ND at RL quantitated per pattern recognition.

Detection limits are reported relative to sample result significant figures.

|               |          |              |          |
|---------------|----------|--------------|----------|
| Sample Amount | : 1000ml | Final Volume | : 5ml    |
| Prepared by   | : P0reto | Analyzed by  | : SDeeso |

EMAX QUALITY CONTROL DATA  
LAB CONTROL SAMPLE ANALYSIS

CLIENT : EUROFINS EATON ANALYTICAL  
PROJECT : 380-32096  
BATCH NO. : 22L294  
METHOD : 3520C/8015B

MATRIX : WATER % MOISTURE:NA  
DILUTION FACTOR: 1 1  
SAMPLE ID : MBLK1W LCS1W  
LAB SAMPLE ID : DSL040WB J8L040WL  
LAB FILE ID : LA05026A LA05029A  
DATE PREPARED : 12/28/22 13:15 12/28/22 13:15  
DATE ANALYZED : 01/05/23 20:12 01/05/23 21:07  
PREP BATCH : 22DSL040W 22DSL040W  
CALIBRATION REF: LA05022A LA05022A

ACCESSION:

| PARAMETERS | MBResult<br>(mg/L) | SpikeAmt<br>(mg/L) | LCSResult<br>(mg/L) | LCSRec<br>(%) | QCLimit<br>(%) |
|------------|--------------------|--------------------|---------------------|---------------|----------------|
| JP8        | ND                 | 2.50               | 2.90                | 116           | 30-160         |

| SURROGATE PARAMETERS | SpikeAmt<br>(mg/L) | LCSResult<br>(mg/L) | LCSRec<br>(%) | QCLimit<br>(%) |
|----------------------|--------------------|---------------------|---------------|----------------|
| Bromobenzene         | 0.500              | 0.489               | 98            | 60-130         |
| Hexacosane           | 0.125              | 0.127               | 102           | 60-130         |

MB: Method Blank sample LCS: Lab Control Sample

EMAX QUALITY CONTROL DATA  
MS/MSD ANALYSIS

CLIENT : EUROFINS EATON ANALYTICAL  
PROJECT : 380-32386  
BATCH NO. : 22L302  
METHOD : 3520C/8015B

|                  |                  |                |                |
|------------------|------------------|----------------|----------------|
| MATRIX           | : WATER          |                | % MOISTURE:NA  |
| DILUTION FACTOR: | 1                | 1              | 1              |
| SAMPLE ID        | : 380-32386-1    | 380-32386-1MS  | 380-32386-1MSD |
| LAB SAMPLE ID    | : 22L302-01      | 22L302-01M     | 22L302-01S     |
| LAB FILE ID      | : LA05039A       | LA05040A       | LA05041A       |
| DATE PREPARED    | : 12/28/22 13:15 | 12/28/22 13:15 | 12/28/22 13:15 |
| DATE ANALYZED    | : 01/06/23 00:10 | 01/06/23 00:28 | 01/06/23 00:47 |
| PREP BATCH       | : 22DSL040W      | 22DSL040W      | 22DSL040W      |
| CALIBRATION REF: | LA05034A         | LA05034A       | LA05034A       |

ACCESSION:

| PARAMETERS | PSResult<br>(mg/L) | SpikeAmt<br>(mg/L) | MSResult<br>(mg/L) | MSRec<br>(%) | SpikeAmt<br>(mg/L) | MSDResult<br>(mg/L) | MSDRec<br>(%) | RPD<br>(%) | QCLimit<br>(%) | MaxRPD<br>(%) |
|------------|--------------------|--------------------|--------------------|--------------|--------------------|---------------------|---------------|------------|----------------|---------------|
| Diesel     | ND                 | 2.75               | 3.15               | 115          | 2.75               | 3.19                | 116           | 1          | 50-130         | 30            |

| SURROGATE PARAMETERS | SpikeAmt<br>(mg/L) | MSResult<br>(mg/L) | MSRec<br>(%) | SpikeAmt<br>(mg/L) | MSDResult<br>(mg/L) | MSDRec<br>(%) | QCLimit<br>(%) |
|----------------------|--------------------|--------------------|--------------|--------------------|---------------------|---------------|----------------|
| Bromobenzene         | 0.550              | 0.531              | 97           | 0.550              | 0.598               | 109           | 60-130         |
| Hexacosane           | 0.138              | 0.153              | 111          | 0.138              | 0.159               | 116           | 60-130         |

PS: Parent Sample MS: Matrix Spike MSD: Matrix Spike Duplicate



EMAX QUALITY CONTROL DATA  
MS/MSD ANALYSIS

CLIENT : EUROFINS EATON ANALYTICAL  
PROJECT : 380-32386  
BATCH NO. : 22L302  
METHOD : 3520C/8015B

|                  |                  |                |                |
|------------------|------------------|----------------|----------------|
| MATRIX           | : WATER          |                | % MOISTURE:NA  |
| DILUTION FACTOR: | 1                | 1              | 1              |
| SAMPLE ID        | : 380-32386-1    | 380-32386-1MS  | 380-32386-1MSD |
| LAB SAMPLE ID    | : 22L302-01      | 22L302-01M     | 22L302-01S     |
| LAB FILE ID      | : LA05039A       | LA05042A       | LA05043A       |
| DATE PREPARED    | : 12/28/22 13:15 | 12/28/22 13:15 | 12/28/22 13:15 |
| DATE ANALYZED    | : 01/06/23 00:10 | 01/06/23 01:05 | 01/06/23 01:23 |
| PREP BATCH       | : 22DSL040W      | 22DSL040W      | 22DSL040W      |
| CALIBRATION REF: | LA05035A         | LA05035A       | LA05035A       |

ACCESSION:

| PARAMETERS | PSResult<br>(mg/L) | SpikeAmt<br>(mg/L) | MSResult<br>(mg/L) | MSRec<br>(%) | SpikeAmt<br>(mg/L) | MSDResult<br>(mg/L) | MSDRec<br>(%) | RPD<br>(%) | QCLimit<br>(%) | MaxRPD<br>(%) |
|------------|--------------------|--------------------|--------------------|--------------|--------------------|---------------------|---------------|------------|----------------|---------------|
| JP5        | ND                 | 2.78               | 3.07               | 111          | 2.80               | 2.84                | 101           | 8          | 30-160         | 30            |

| SURROGATE PARAMETERS | SpikeAmt<br>(mg/L) | MSResult<br>(mg/L) | MSRec<br>(%) | SpikeAmt<br>(mg/L) | MSDResult<br>(mg/L) | MSDRec<br>(%) | QCLimit<br>(%) |
|----------------------|--------------------|--------------------|--------------|--------------------|---------------------|---------------|----------------|
| Bromobenzene         | 0.555              | 0.557              | 100          | 0.560              | 0.601               | 107           | 60-130         |
| Hexacosane           | 0.139              | 0.162              | 117          | 0.140              | 0.141               | 101           | 60-130         |

PS: Parent Sample MS: Matrix Spike MSD: Matrix Spike Duplicate

EMAX QUALITY CONTROL DATA  
MS/MSD ANALYSIS

CLIENT : EUROFINS EATON ANALYTICAL  
PROJECT : 380-32289  
BATCH NO. : 22L303  
METHOD : 3520C/8015B

|                  |                  |                |                |
|------------------|------------------|----------------|----------------|
| MATRIX           | : WATER          |                | % MOISTURE:NA  |
| DILUTION FACTOR: | 1                | 1              | 1              |
| SAMPLE ID        | : 380-32289-1    | 380-32289-1MS  | 380-32289-1MSD |
| LAB SAMPLE ID    | : 22L303-01      | 22L303-01M     | 22L303-01S     |
| LAB FILE ID      | : LA05044A       | LA05045A       | LA05046A       |
| DATE PREPARED    | : 12/28/22 13:15 | 12/28/22 13:15 | 12/28/22 13:15 |
| DATE ANALYZED    | : 01/06/23 01:42 | 01/06/23 02:00 | 01/06/23 02:18 |
| PREP BATCH       | : 22DSL040W      | 22DSL040W      | 22DSL040W      |
| CALIBRATION REF: | LA05036A         | LA05036A       | LA05036A       |

ACCESSION:

| PARAMETERS | PSResult<br>(mg/L) | SpikeAmt<br>(mg/L) | MSResult<br>(mg/L) | MSRec<br>(%) | SpikeAmt<br>(mg/L) | MSDResult<br>(mg/L) | MSDRec<br>(%) | RPD<br>(%) | QCLimit<br>(%) | MaxRPD<br>(%) |
|------------|--------------------|--------------------|--------------------|--------------|--------------------|---------------------|---------------|------------|----------------|---------------|
| JP8        | ND                 | 2.55               | 3.02               | 118          | 2.55               | 2.75                | 108           | 9          | 30-160         | 30            |

| SURROGATE PARAMETERS | SpikeAmt<br>(mg/L) | MSResult<br>(mg/L) | MSRec<br>(%) | SpikeAmt<br>(mg/L) | MSDResult<br>(mg/L) | MSDRec<br>(%) | QCLimit<br>(%) |
|----------------------|--------------------|--------------------|--------------|--------------------|---------------------|---------------|----------------|
| Bromobenzene         | 0.510              | 0.525              | 103          | 0.510              | 0.526               | 103           | 60-130         |
| Hexacosane           | 0.127              | 0.132              | 104          | 0.127              | 0.132               | 104           | 60-130         |

PS: Parent Sample MS: Matrix Spike MSD: Matrix Spike Duplicate

January 09, 2023

Rosalynn Dang  
 Eurofins Eaton Analytical  
 750 Royal Oaks Drive  
 Suite 100  
 Monrovia, CA 91016-

Project Name: RED-HILL Project # 38001111 Job # 380-32096-1  
 Physis Project ID: 1407003-351

Dear Rosalynn,


Enclosed are the analytical results for samples submitted to PHYSIS Environmental Laboratories, Inc. (PHYSIS) on 12/21/2022. A total of 3 samples were received for analysis in accordance with the attached chain of custody (COC). Per the COC, the samples were analyzed for:

| Organics                                       |
|--|
| Polynuclear Aromatic Hydrocarbons by EPA 625.1 |
| Disalicylidenepropanediamine by EPA 625.1      |
| Dibenzo [a,l] Pyrene w/ PAHs by EPA 625.1      |

Analytical results in this report apply only to samples submitted to PHYSIS in accordance with the COC and are intended to be considered in their entirety.

Please feel free to contact me at any time with any questions. PHYSIS appreciates the opportunity to provide you with our analytical and support services.

Regards,

  
 Misty Mercier  
 714 602-5320  
 Extension 202  
 mistymercier@physislabs.com



## PROJECT SAMPLE LIST

Eurofins Eaton Analytical

PHYSIS Project ID: 1407003-351

RED-HILL Project # 38001111 Job # 380-32096-1

Total Samples: 3

| PHYSIS ID | Sample ID               | Description | Date      | Time  | Matrix      | Sample Type   |
|-----------|-------------------------|-------------|-----------|-------|-------------|---------------|
| 102965    | Aiea Gulch Wells Pump 2 | 380-32096-1 | 12/19/202 | 10:00 | Samplewater | Not Specified |
| 102966    | Aiea Wells Pump 2       | 380-32096-2 | 12/19/202 | 10:45 | Samplewater | Not Specified |
| 102967    | Halawa Wells Pump 1     | 380-32096-3 | 12/19/202 | 11:18 | Samplewater | Not Specified |

## ABBREVIATIONS and ACRONYMS

|      |  |
|------|--|
| QM   | Quality Manual                         |
| QA   | Quality Assurance                      |
| QC   | Quality Control                        |
| MDL  | method detection limit                 |
| RL   | reporting limit                        |
| R1   | project sample                         |
| R2   | project sample replicate               |
| MS1  | matrix spike                           |
| MS2  | matrix spike replicate                 |
| B1   | procedural blank                       |
| B2   | procedural blank replicate             |
| BS1  | blank spike                            |
| BS2  | blank spike replicate                  |
| LCS1 | laboratory control spike               |
| LCS2 | laboratory control spike replicate     |
| LCM1 | laboratory control material            |
| LCM2 | laboratory control material replicate  |
| CRM1 | certified reference material           |
| CRM2 | certified reference material replicate |
| RPD  | relative percent difference            |
| LMW  | low molecular weight                   |
| HMW  | high molecular weight                  |

## QUALITY ASSURANCE SUMMARY

**LABORATORY BATCH:** Physis' QM defines a laboratory batch as a group of 20 or fewer project samples of similar matrix, processed together under the same conditions and with the same reagents. QC samples are associated with each batch and were used to assess the validity of the sample analyses.

**PROCEDURAL BLANK:** Laboratory contamination introduced during method use is assessed through the preparation and analysis of procedural blanks is provided at a minimum frequency of one per batch.

**ACCURACY:** Accuracy of analytical measurements is the degree of closeness based on percent recovery calculations between measured values and the actual or true value and includes a combination of reproducibility error and systematic bias due to sampling and analytical operations. Accuracy of the project data was indicated by analysis of MS, BS, LCS, LCM, CRM, and/or surrogate spikes on a minimum frequency of one per batch. Physis' QM requires that 95% of the target compounds greater than 10 times the MDL be within the specified acceptance limits.

**PRECISION:** Precision is the agreement among a set of replicate measurements without assumption of knowledge of the true value and is based on RPD calculations between repeated values. Precision of the project data was determined by analysis of replicate MS<sub>1</sub>/MS<sub>2</sub>, BS<sub>1</sub>/BS<sub>2</sub>, LCS<sub>1</sub>/LCS<sub>2</sub>, LCM<sub>1</sub>/LCM<sub>2</sub>, CRM<sub>1</sub>/CRM<sub>2</sub>, surrogate spikes and/or replicate project sample analysis (R<sub>1</sub>/R<sub>2</sub>) on a minimum frequency of one per batch. Physis' QM requires that for 95% of the compounds greater than 10 times the MDL, the percent RPD should be within the specified acceptance range.

**BLANK SPIKES:** BS is the introduction of a known concentration of analyte into the procedural blank. BS demonstrates performance of the preparation and analytical methods on a clean matrix void of potential matrix related interferences. The BS is performed in laboratory deionized water, making these recoveries a better indicator of the efficiency of the laboratory method per se.

**MATRIX SPIKES:** MS is the introduction of a known concentration of analyte into a sample. MS samples demonstrate the effect a particular project sample matrix has on the accuracy of a measurement. Individually, MS samples also indicate the bias of analytical measurements due to chemical interferences inherent in the in the specific project sample spiked. Intrinsic target analyte concentration in the specific project sample can also significantly impact MS recovery.

**CERTIFIED REFERENCE MATERIALS:** CRMs are materials of various matrices for which analytical information has been determined and certified by a recognized authority. These are used to provide a quantitative assessment of the accuracy of an analytical method. CRMs provide evidence that the laboratory preparation and analysis produces results that are comparable to those obtained by an independent organization.

**LABORATORY CONTROL MATERIAL:** LCM is provided because a suitable natural seawater CRM is not available and can be used to indicate accuracy of the method. Physis' internal LCM is seawater collected at ~800 meters in the Southern California San Pedro Basin and can be used as a reference for background concentrations in clean, natural seawater for comparison to project samples.

**LABORATORY CONTROL SPIKES:** LCS is the introduction of a known concentration of analyte into Physis' LCM. LCS samples were employed to assess the effect the seawater matrix has on the accuracy of a measurement. LCS also indicate the bias of this method due to chemical interferences inherent in the in the seawater matrix. Intrinsic LCM concentration can also significantly impact LCS recovery.

**SURROGATES:** A surrogate is a pure analyte unlikely to be found in any project sample, behaves similarly to

the target analyte and most often used with organic analytical procedures. Surrogates are added in known concentration to all samples and are measured to indicate overall efficiency of the method including processing and analyses.

**HOLDING TIME:** Method recommended holding times are the length of time a project sample can be stored under specific conditions after collection and prior to analysis without significantly affecting the analyte's concentration. Holding times can be extended if preservation techniques are employed to reduce biodegradation, volatilization, oxidation, sorption, precipitation, and other physical and chemical processes.

**SAMPLE STORAGE/RETENTION:** In order to maintain chemical integrity prior to analysis, all samples submitted to Physis are refrigerated (liquids) or frozen (solids) upon receipt unless otherwise recommended by applicable methods. Solid samples are retained for 1 year from collection while liquid samples are retained until method recommended holding times elapse.

**TOTAL/DISSOLVED FRACTION:** In some instances, the results for the dissolved fraction may be higher than the total fraction for a particular analyte (e.g. trace metals). This is typically caused by the analytical variation for each result and indicates that the target analyte is primarily in the dissolved phase, within the sample.

## PHYSIS QUALIFIER CODES

| CODE | DEFINITION  |
|------|---|
| #    | see Case Narrative  |
| ND   | analyte not detected at or above the MDL  |
| B    | analyte was detected in the procedural blank greater than 10 times the MDL  |
| E    | analyte concentration exceeds the upper limit of the linear calibration range, reported value is estimated  |
| H    | sample received and/or analyzed past the recommended holding time   |
| J    | analyte was detected at a concentration below the RL and above the MDL, reported value is estimated   |
| N    | insufficient sample, analysis could not be performed  |
| M    | analyte was outside the specified accuracy and/or precision acceptance limits due to matrix interference. The associated B/BS were within limits, therefore the sample data was reported without further clarification  |
| SH   | analyte concentration in the project sample exceeded the spike concentration, therefore accuracy and/or precision acceptance limits do not apply  |
| SL   | analyte results were lower than 10 times the MDL, therefore accuracy and/or precision acceptance limits do not apply  |
| NH   | project sample was heterogeneous and sample homogeneity could not be readily achieved using routine laboratory practices, therefore accuracy and/or precision acceptance limits do not apply  |
| Q    | analyte was outside the specified QAPP acceptance limits for precision and/or accuracy but within Physis derived acceptance limits, therefore the sample data was reported without further clarification  |
| R    | Physis' QM allows for 5% of the target compounds greater than 10 times the MDL to be outside the specified acceptance limits for precision and/or accuracy. This is often due to random error and does not indicate any significant problems with the analysis of these project samples |



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## CASE NARRATIVE

### QUALIFIER NOTES

In addition to the use of analyte specific Physis Qualifier Codes where applicable, the following were also noted.

#### **ND**

MDL is listed due to report format restrictions; it is not used in reporting. Analytical results reported are ND at the RL.

# ANALYTICALS

# REPORT

TERRA AURA  
ENVIRONMENTAL LABORATORIES, INC.

*Innovative Solutions for Nature*

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- 2
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### Base/Neutral Extractable Compounds

| ANALYTE  | Method    | Units | RESULT | DF | MDL  | RL  | Fraction | QA CODE | Batch ID | Date Processed | Date Analyzed |
|--|-----------|-------|--------|----|------|-----|----------|---------|----------|----------------|---------------|
| <b>Sample ID: 102965-R1 Aiea Gulch Wells Pump 2 380-3209 Matrix: Samplewater</b> |           |       |        |    |      |     |          |         |          |                |               |
| Disalicylideneopropanediamine  | EPA 625.1 | µg/L  | ND     | 1  | 0.05 | 0.1 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| <b>Sample ID: 102966-R1 Aiea Wells Pump 2 380-32096-2 Matrix: Samplewater</b>    |           |       |        |    |      |     |          |         |          |                |               |
| Disalicylideneopropanediamine  | EPA 625.1 | µg/L  | ND     | 1  | 0.05 | 0.1 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| <b>Sample ID: 102967-R1 Halawa Wells Pump 1 380-32096-3 Matrix: Samplewater</b>  |           |       |        |    |      |     |          |         |          |                |               |
| Disalicylideneopropanediamine  | EPA 625.1 | µg/L  | ND     | 1  | 0.05 | 0.1 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |

## Polynuclear Aromatic Hydrocarbons

| ANALYTE  | Method    | Units      | RESULT | DF | MDL   | RL    | Fraction                        | QA CODE | Batch ID                   | Date Processed | Date Analyzed |
|--|-----------|------------|--------|----|-------|-------|---------------------------------|---------|----------------------------|----------------|---------------|
| <b>Sample ID: 102965-R1 Aiea Gulch Wells Pump 2 380-3209 Matrix: Samplewater</b> |           |            |        |    |       |       | <b>Sampled: 19-Dec-22 10:00</b> |         | <b>Received: 21-Dec-22</b> |                |               |
| (d10-Acenaphthene)   | EPA 625.1 | % Recovery | 82     | 1  |       |       | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| (d10-Phenanthrene)   | EPA 625.1 | % Recovery | 87     | 1  |       |       | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| (d12-Chrysene)   | EPA 625.1 | % Recovery | 94     | 1  |       |       | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| (d12-Perylene)   | EPA 625.1 | % Recovery | 85     | 1  |       |       | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| (d8-Naphthalene)   | EPA 625.1 | % Recovery | 72     | 1  |       |       | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| 1-Methylnaphthalene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| 1-Methylphenanthrene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| 2,3,5-Trimethylnaphthalene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| 2,6-Dimethylnaphthalene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| 2-Methylnaphthalene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| Acenaphthene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| Acenaphthylene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| Anthracene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| Benz[a]anthracene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| Benzo[a]pyrene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| Benzo[b]fluoranthene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| Benzo[e]pyrene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| Benzo[g,h,i]perylene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| Benzo[k]fluoranthene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| Biphenyl   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| Chrysene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| D benz[a,h]anthracene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| D benzo[a,l]pyrene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |
| D benzothiophene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074                    | 26-Dec-22      | 08-Jan-23     |

### Polynuclear Aromatic Hydrocarbons

| ANALYTE                | Method    | Units | RESULT | DF | MDL   | RL    | Fraction | QA CODE | Batch ID | Date Processed | Date Analyzed |
|------------------------|-----------|-------|--------|----|-------|-------|----------|---------|----------|----------------|---------------|
| Fluoranthene           | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Fluorene               | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Indeno[1,2,3-cd]pyrene | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Naphthalene            | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Perylene               | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Phenanthrene           | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Pyrene                 | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |



## Polynuclear Aromatic Hydrocarbons

| ANALYTE                     | Method                               | Units                      | RESULT | DF | MDL   | RL    | Fraction                        | QA CODE | Batch ID | Date Processed             | Date Analyzed |
|-----------------------------|--------------------------------------|----------------------------|--------|----|-------|-------|---------------------------------|---------|----------|----------------------------|---------------|
| <b>Sample ID: 102966-R1</b> | <b>Aiea Wells Pump 2 380-32096-2</b> | <b>Matrix: Samplewater</b> |        |    |       |       | <b>Sampled: 19-Dec-22 10:45</b> |         |          | <b>Received: 21-Dec-22</b> |               |
| (d10-Acenaphthene)          | EPA 625.1                            | % Recovery                 | 86     | 1  |       |       | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| (d10-Phenanthrene)          | EPA 625.1                            | % Recovery                 | 91     | 1  |       |       | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| (d12-Chrysene)              | EPA 625.1                            | % Recovery                 | 98     | 1  |       |       | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| (d12-Perylene)              | EPA 625.1                            | % Recovery                 | 91     | 1  |       |       | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| (d8-Naphthalene)            | EPA 625.1                            | % Recovery                 | 78     | 1  |       |       | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| 1-Methylnaphthalene         | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| 1-Methylphenanthrene        | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| 2,3,5-Trimethylnaphthalene  | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| 2,6-Dimethylnaphthalene     | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| 2-Methylnaphthalene         | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| Acenaphthene                | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| Acenaphthylene              | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| Anthracene                  | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| Benz[a]anthracene           | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| Benzo[a]pyrene              | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| Benzo[b]fluoranthene        | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| Benzo[e]pyrene              | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| Benzo[g,h,i]perylene        | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| Benzo[k]fluoranthene        | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| Biphenyl                    | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| Chrysene                    | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| D benz[a,h]anthracene       | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| D benzo[a,l]pyrene          | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |
| D benzothiophene            | EPA 625.1                            | µg/L                       | ND     | 1  | 0.001 | 0.005 | Total                           |         | O-40074  | 26-Dec-22                  | 08-Jan-23     |

### Polynuclear Aromatic Hydrocarbons

| ANALYTE                | Method    | Units | RESULT | DF | MDL   | RL    | Fraction | QA CODE | Batch ID | Date Processed | Date Analyzed |
|------------------------|-----------|-------|--------|----|-------|-------|----------|---------|----------|----------------|---------------|
| Fluoranthene           | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Fluorene               | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Indeno[1,2,3-cd]pyrene | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Naphthalene            | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Perylene               | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Phenanthrene           | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Pyrene                 | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |



## Polynuclear Aromatic Hydrocarbons

| ANALYTE   | Method    | Units      | RESULT | DF | MDL   | RL    | Fraction | QA CODE | Batch ID | Date Processed | Date Analyzed |
|---|-----------|------------|--------|----|-------|-------|----------|---------|----------|----------------|---------------|
| <b>Sample ID: 102967-R1 Halawa Wells Pump 1 380-32096-3 Matrix: Samplewater</b> |           |            |        |    |       |       |          |         |          |                |               |
| <b>Sampled: 19-Dec-22 11:18 Received: 21-Dec-22</b>                             |           |            |        |    |       |       |          |         |          |                |               |
| (d10-Acenaphthene)  | EPA 625.1 | % Recovery | 90     | 1  |       |       | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| (d10-Phenanthrene)  | EPA 625.1 | % Recovery | 94     | 1  |       |       | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| (d12-Chrysene)  | EPA 625.1 | % Recovery | 99     | 1  |       |       | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| (d12-Perylene)  | EPA 625.1 | % Recovery | 86     | 1  |       |       | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| (d8-Naphthalene)  | EPA 625.1 | % Recovery | 80     | 1  |       |       | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| 1-Methylnaphthalene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| 1-Methylphenanthrene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| 2,3,5-Trimethylnaphthalene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| 2,6-Dimethylnaphthalene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| 2-Methylnaphthalene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Acenaphthene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Acenaphthylene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Anthracene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Benz[a]anthracene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Benzo[a]pyrene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Benzo[b]fluoranthene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Benzo[e]pyrene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Benzo[g,h,i]perylene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Benzo[k]fluoranthene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Biphenyl  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Chrysene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| D benz[a,h]anthracene   | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| D benzo[a,l]pyrene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| D benzothiophene  | EPA 625.1 | µg/L       | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |



## Polynuclear Aromatic Hydrocarbons

| ANALYTE                | Method    | Units | RESULT | DF | MDL   | RL    | Fraction | QA CODE | Batch ID | Date Processed | Date Analyzed |
|------------------------|-----------|-------|--------|----|-------|-------|----------|---------|----------|----------------|---------------|
| Fluoranthene           | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Fluorene               | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Indeno[1,2,3-cd]pyrene | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Naphthalene            | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Perylene               | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Phenanthrene           | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |
| Pyrene                 | EPA 625.1 | µg/L  | ND     | 1  | 0.001 | 0.005 | Total    |         | O-40074  | 26-Dec-22      | 08-Jan-23     |



# QUALITY CONTROL REPORT

TERRA AURA  
ENVIRONMENTAL LABORATORIES, INC.

*Innovative Solutions for Nature*

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## Base/Neutral Extractable Compounds

## QUALITY CONTROL REPORT

| ANALYTE                      | FRACTION | RESULT                       | DF | MDL  | RL                         | UNITS | SPIKE | SOURCE              | ACCURACY |                     | PRECISION |        | QA CODEc |
|------------------------------|----------|------------------------------|----|------|----------------------------|-------|-------|---------------------|----------|---------------------|-----------|--------|----------|
|                              |          |                              |    |      |                            |       | LEVEL | RESULT              | %        | LIMITS              | %         | LIMITS |          |
| <b>Sample ID: 102964-B1</b>  |          | <b>QAQC Procedural Blank</b> |    |      | <b>Matrix: BlankMatrix</b> |       |       | <b>Sampled:</b>     |          | <b>Received:</b>    |           |        |          |
|                              |          | Method: EPA 625.1            |    |      | Batch ID: O-40074          |       |       | Prepared: 26-Dec-22 |          | Analyzed: 08-Jan-23 |           |        |          |
| Disalicylideneprapanediamin  | Total    | ND                           | 1  | 0.05 | 0.1                        | µg/L  |       |                     |          |                     |           |        |          |
| <b>Sample ID: 102964-BS1</b> |          | <b>QAQC Procedural Blank</b> |    |      | <b>Matrix: BlankMatrix</b> |       |       | <b>Sampled:</b>     |          | <b>Received:</b>    |           |        |          |
|                              |          | Method: EPA 625.1            |    |      | Batch ID: O-40074          |       |       | Prepared: 26-Dec-22 |          | Analyzed: 08-Jan-23 |           |        |          |
| Disalicylideneprapanediamin  | Total    | 42.9                         | 1  | 0.05 | 0.1                        | µg/L  | 50    | 0                   | 86       | 50 - 150%           | PASS      |        |          |
| <b>Sample ID: 102964-BS2</b> |          | <b>QAQC Procedural Blank</b> |    |      | <b>Matrix: BlankMatrix</b> |       |       | <b>Sampled:</b>     |          | <b>Received:</b>    |           |        |          |
|                              |          | Method: EPA 625.1            |    |      | Batch ID: O-40074          |       |       | Prepared: 26-Dec-22 |          | Analyzed: 08-Jan-23 |           |        |          |
| Disalicylideneprapanediamin  | Total    | 45.8                         | 1  | 0.05 | 0.1                        | µg/L  | 50    | 0                   | 92       | 50 - 150%           | PASS      | 6      | 30 PASS  |

**Polynuclear Aromatic Hydrocarbons**

**QUALITY CONTROL REPORT**

| ANALYTE                     | FRACTION | RESULT                       | DF | MDL   | RL                         | UNITS | SPIKE      | SOURCE              | ACCURACY | PRECISION           | QA CODE |        |
|-----------------------------|----------|------------------------------|----|-------|----------------------------|-------|------------|---------------------|----------|---------------------|---------|--------|
|                             |          |                              |    |       |                            |       | LEVEL      | RESULT              | %        | LIMITS              | %       | LIMITS |
| <b>Sample ID: 102964-B1</b> |          | <b>QAQC Procedural Blank</b> |    |       | <b>Matrix: BlankMatrix</b> |       |            | <b>Sampled:</b>     |          | <b>Received:</b>    |         |        |
|                             |          | Method: EPA 625.1            |    |       | Batch ID: O-40074          |       |            | Prepared: 26-Dec-22 |          | Analyzed: 08-Jan-23 |         |        |
| (d10-Acenaphthene)          | Total    | 92                           | 1  |       |                            |       | % Recovery | 100                 | 92       | 27 - 133%           | PASS    |        |
| (d10-Phenanthrene)          | Total    | 94                           | 1  |       |                            |       | % Recovery | 100                 | 94       | 43 - 129%           | PASS    |        |
| (d12-Chrysene)              | Total    | 99                           | 1  |       |                            |       | % Recovery | 100                 | 99       | 52 - 144%           | PASS    |        |
| (d12-Perylene)              | Total    | 103                          | 1  |       |                            |       | % Recovery | 100                 | 103      | 36 - 161%           | PASS    |        |
| (d8-Naphthalene)            | Total    | 87                           | 1  |       |                            |       | % Recovery | 100                 | 87       | 25 - 125%           | PASS    |        |
| 1-Methylnaphthalene         | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| 1-Methylphenanthrene        | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| 2,3,5-Trimethylnaphthalene  | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| 2,6-Dimethylnaphthalene     | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| 2-Methylnaphthalene         | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| Acenaphthene                | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| Acenaphthylene              | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| Anthracene                  | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| Benz[a]anthracene           | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| Benzo[a]pyrene              | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| Benzo[b]fluoranthene        | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| Benzo[e]pyrene              | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| Benzo[g,h,i]perylene        | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| Benzo[k]fluoranthene        | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| Biphenyl                    | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| Chrysene                    | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| Dibenz[a,h]anthracene       | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| Dibenzo[a,l]pyrene          | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |
| Dibenzothiophene            | Total    | ND                           | 1  | 0.001 | 0.005                      | µg/L  |            |                     |          |                     |         |        |



## Polynuclear Aromatic Hydrocarbons

## QUALITY CONTROL REPORT

| ANALYTE                | FRACTION | RESULT | DF | MDL   | RL    | UNITS | SPIKE | SOURCE | ACCURACY |        | PRECISION |        | QA CODEc |
|------------------------|----------|--------|----|-------|-------|-------|-------|--------|----------|--------|-----------|--------|----------|
|                        |          |        |    |       |       |       | LEVEL | RESULT | %        | LIMITS | %         | LIMITS |          |
| Fluoranthene           | Total    | ND     | 1  | 0.001 | 0.005 | µg/L  |       |        |          |        |           |        |          |
| Fluorene               | Total    | ND     | 1  | 0.001 | 0.005 | µg/L  |       |        |          |        |           |        |          |
| Indeno[1,2,3-cd]pyrene | Total    | ND     | 1  | 0.001 | 0.005 | µg/L  |       |        |          |        |           |        |          |
| Naphthalene            | Total    | ND     | 1  | 0.001 | 0.005 | µg/L  |       |        |          |        |           |        |          |
| Perylene               | Total    | ND     | 1  | 0.001 | 0.005 | µg/L  |       |        |          |        |           |        |          |
| Phenanthrene           | Total    | ND     | 1  | 0.001 | 0.005 | µg/L  |       |        |          |        |           |        |          |
| Pyrene                 | Total    | ND     | 1  | 0.001 | 0.005 | µg/L  |       |        |          |        |           |        |          |



## Polynuclear Aromatic Hydrocarbons

## QUALITY CONTROL REPORT

| ANALYTE                      | FRACTION | RESULT                       | DF | MDL   | RL                         | UNITS      | SPIKE               | SOURCE          | ACCURACY |                  | PRECISION | QA CODE |
|------------------------------|----------|------------------------------|----|-------|----------------------------|------------|---------------------|-----------------|----------|------------------|-----------|---------|
|                              |          |                              |    |       |                            |            | LEVEL               | RESULT          | %        | LIMITS           | %         | LIMITS  |
| <b>Sample ID: 102964-BS1</b> |          | <b>QAQC Procedural Blank</b> |    |       | <b>Matrix: BlankMatrix</b> |            |                     | <b>Sampled:</b> |          | <b>Received:</b> |           |         |
| Method: EPA 625.1            |          | Batch ID: O-40074            |    |       | Prepared: 26-Dec-22        |            | Analyzed: 08-Jan-23 |                 |          |                  |           |         |
| (d10-Acenaphthene)           | Total    | 93                           | 1  |       |                            | % Recovery | 100                 | 0               | 93       | 27 - 133%        | PASS      |         |
| (d10-Phenanthrene)           | Total    | 96                           | 1  |       |                            | % Recovery | 100                 | 0               | 96       | 43 - 129%        | PASS      |         |
| (d12-Chrysene)               | Total    | 97                           | 1  |       |                            | % Recovery | 100                 | 0               | 97       | 52 - 144%        | PASS      |         |
| (d12-Perylene)               | Total    | 99                           | 1  |       |                            | % Recovery | 100                 | 0               | 99       | 36 - 161%        | PASS      |         |
| (d8-Naphthalene)             | Total    | 84                           | 1  |       |                            | % Recovery | 100                 | 0               | 84       | 25 - 125%        | PASS      |         |
| 1-Methylnaphthalene          | Total    | 0.477                        | 1  | 0.001 | 0.005                      | µg/L       | 0.5                 | 0               | 95       | 31 - 128%        | PASS      |         |
| 1-Methylphenanthrene         | Total    | 0.448                        | 1  | 0.001 | 0.005                      | µg/L       | 0.5                 | 0               | 90       | 66 - 127%        | PASS      |         |
| 2,3,5-Trimethylnaphthalene   | Total    | 0.498                        | 1  | 0.001 | 0.005                      | µg/L       | 0.5                 | 0               | 100      | 55 - 122%        | PASS      |         |
| 2,6-Dimethylnaphthalene      | Total    | 0.493                        | 1  | 0.001 | 0.005                      | µg/L       | 0.5                 | 0               | 99       | 48 - 120%        | PASS      |         |
| 2-Methylnaphthalene          | Total    | 1.44                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5                 | 0               | 96       | 47 - 130%        | PASS      |         |
| Acenaphthene                 | Total    | 1.49                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5                 | 0               | 99       | 53 - 131%        | PASS      |         |
| Acenaphthylene               | Total    | 1.52                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5                 | 0               | 101      | 43 - 140%        | PASS      |         |
| Anthracene                   | Total    | 1.52                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5                 | 0               | 101      | 58 - 135%        | PASS      |         |
| Benz[a]anthracene            | Total    | 1.56                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5                 | 0               | 104      | 55 - 145%        | PASS      |         |
| Benzo[a]pyrene               | Total    | 1.54                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5                 | 0               | 103      | 51 - 143%        | PASS      |         |
| Benzo[b]fluoranthene         | Total    | 1.63                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5                 | 0               | 109      | 46 - 165%        | PASS      |         |
| Benzo[e]pyrene               | Total    | 0.502                        | 1  | 0.001 | 0.005                      | µg/L       | 0.5                 | 0               | 100      | 42 - 152%        | PASS      |         |
| Benzo[g,h,i]perylene         | Total    | 1.59                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5                 | 0               | 106      | 63 - 133%        | PASS      |         |
| Benzo[k]fluoranthene         | Total    | 1.64                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5                 | 0               | 109      | 56 - 145%        | PASS      |         |
| Biphenyl                     | Total    | 0.523                        | 1  | 0.001 | 0.005                      | µg/L       | 0.5                 | 0               | 105      | 56 - 119%        | PASS      |         |
| Chrysene                     | Total    | 1.55                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5                 | 0               | 103      | 56 - 141%        | PASS      |         |
| Dibenz[a,h]anthracene        | Total    | 1.55                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5                 | 0               | 103      | 55 - 150%        | PASS      |         |
| Dibenzo[a,l]pyrene           | Total    | 0.539                        | 1  | 0.001 | 0.005                      | µg/L       | 0.5                 | 0               | 108      | 50 - 150%        | PASS      |         |
| Dibenzothiophene             | Total    | 0.459                        | 1  | 0.001 | 0.005                      | µg/L       | 0.5                 | 0               | 92       | 46 - 126%        | PASS      |         |

## Polynuclear Aromatic Hydrocarbons

## QUALITY CONTROL REPORT

| ANALYTE                | FRACTION | RESULT | DF | MDL   | RL    | UNITS | SPIKE | SOURCE | ACCURACY |           | PRECISION |        | QA CODE <sub>c</sub> |
|------------------------|----------|--------|----|-------|-------|-------|-------|--------|----------|-----------|-----------|--------|----------------------|
|                        |          |        |    |       |       |       | LEVEL | RESULT | %        | LIMITS    | %         | LIMITS |                      |
| Fluoranthene           | Total    | 1.68   | 1  | 0.001 | 0.005 | µg/L  | 1.5   | 0      | 112      | 60 - 146% | PASS      |        |                      |
| Fluorene               | Total    | 1.58   | 1  | 0.001 | 0.005 | µg/L  | 1.5   | 0      | 105      | 58 - 131% | PASS      |        |                      |
| Indeno[1,2,3-cd]pyrene | Total    | 1.61   | 1  | 0.001 | 0.005 | µg/L  | 1.5   | 0      | 107      | 50 - 151% | PASS      |        |                      |
| Naphthalene            | Total    | 1.38   | 1  | 0.001 | 0.005 | µg/L  | 1.5   | 0      | 92       | 41 - 126% | PASS      |        |                      |
| Perylene               | Total    | 0.478  | 1  | 0.001 | 0.005 | µg/L  | 0.5   | 0      | 96       | 48 - 141% | PASS      |        |                      |
| Phenanthrene           | Total    | 1.57   | 1  | 0.001 | 0.005 | µg/L  | 1.5   | 0      | 105      | 67 - 127% | PASS      |        |                      |
| Pyrene                 | Total    | 1.62   | 1  | 0.001 | 0.005 | µg/L  | 1.5   | 0      | 108      | 54 - 156% | PASS      |        |                      |

## Polynuclear Aromatic Hydrocarbons

## QUALITY CONTROL REPORT

| ANALYTE                      | FRACTION | RESULT                       | DF | MDL   | RL                         | UNITS      | SPIKE | SOURCE              | ACCURACY |           | PRECISION        |        | QA CODEc |      |
|------------------------------|----------|------------------------------|----|-------|----------------------------|------------|-------|---------------------|----------|-----------|------------------|--------|----------|------|
|                              |          |                              |    |       |                            |            | LEVEL | RESULT              | %        | LIMITS    | %                | LIMITS |          |      |
| <b>Sample ID: 102964-BS2</b> |          | <b>QAQC Procedural Blank</b> |    |       | <b>Matrix: BlankMatrix</b> |            |       | <b>Sampled:</b>     |          |           | <b>Received:</b> |        |          |      |
| Method: EPA 625.1            |          | Batch ID: O-40074            |    |       | Prepared: 26-Dec-22        |            |       | Analyzed: 08-Jan-23 |          |           |                  |        |          |      |
| (d10-Acenaphthene)           | Total    | 92                           | 1  |       |                            | % Recovery | 100   | 0                   | 92       | 27 - 133% | PASS             | 1      | 30       | PASS |
| (d10-Phenanthrene)           | Total    | 95                           | 1  |       |                            | % Recovery | 100   | 0                   | 95       | 43 - 129% | PASS             | 1      | 30       | PASS |
| (d12-Chrysene)               | Total    | 95                           | 1  |       |                            | % Recovery | 100   | 0                   | 95       | 52 - 144% | PASS             | 2      | 30       | PASS |
| (d12-Perylene)               | Total    | 94                           | 1  |       |                            | % Recovery | 100   | 0                   | 94       | 36 - 161% | PASS             | 5      | 30       | PASS |
| (d8-Naphthalene)             | Total    | 84                           | 1  |       |                            | % Recovery | 100   | 0                   | 84       | 25 - 125% | PASS             | 0      | 30       | PASS |
| 1-Methylnaphthalene          | Total    | 0.49                         | 1  | 0.001 | 0.005                      | µg/L       | 0.5   | 0                   | 98       | 31 - 128% | PASS             | 3      | 30       | PASS |
| 1-Methylphenanthrene         | Total    | 0.452                        | 1  | 0.001 | 0.005                      | µg/L       | 0.5   | 0                   | 90       | 66 - 127% | PASS             | 0      | 30       | PASS |
| 2,3,5-Trimethylnaphthalene   | Total    | 0.516                        | 1  | 0.001 | 0.005                      | µg/L       | 0.5   | 0                   | 103      | 55 - 122% | PASS             | 3      | 30       | PASS |
| 2,6-Dimethylnaphthalene      | Total    | 0.505                        | 1  | 0.001 | 0.005                      | µg/L       | 0.5   | 0                   | 101      | 48 - 120% | PASS             | 2      | 30       | PASS |
| 2-Methylnaphthalene          | Total    | 1.48                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5   | 0                   | 99       | 47 - 130% | PASS             | 3      | 30       | PASS |
| Acenaphthene                 | Total    | 1.54                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5   | 0                   | 103      | 53 - 131% | PASS             | 4      | 30       | PASS |
| Acenaphthylene               | Total    | 1.57                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5   | 0                   | 105      | 43 - 140% | PASS             | 4      | 30       | PASS |
| Anthracene                   | Total    | 1.56                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5   | 0                   | 104      | 58 - 135% | PASS             | 3      | 30       | PASS |
| Benz[a]anthracene            | Total    | 1.58                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5   | 0                   | 105      | 55 - 145% | PASS             | 1      | 30       | PASS |
| Benzo[a]pyrene               | Total    | 1.66                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5   | 0                   | 111      | 51 - 143% | PASS             | 7      | 30       | PASS |
| Benzo[b]fluoranthene         | Total    | 1.66                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5   | 0                   | 111      | 46 - 165% | PASS             | 2      | 30       | PASS |
| Benzo[e]pyrene               | Total    | 0.504                        | 1  | 0.001 | 0.005                      | µg/L       | 0.5   | 0                   | 101      | 42 - 152% | PASS             | 1      | 30       | PASS |
| Benzo[g,h,i]perylene         | Total    | 1.57                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5   | 0                   | 105      | 63 - 133% | PASS             | 1      | 30       | PASS |
| Benzo[k]fluoranthene         | Total    | 1.67                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5   | 0                   | 111      | 56 - 145% | PASS             | 2      | 30       | PASS |
| Biphenyl                     | Total    | 0.54                         | 1  | 0.001 | 0.005                      | µg/L       | 0.5   | 0                   | 108      | 56 - 119% | PASS             | 3      | 30       | PASS |
| Chrysene                     | Total    | 1.57                         | 1  | 0.001 | 0.005                      | µg/L       | 1.5   | 0                   | 105      | 56 - 141% | PASS             | 2      | 30       | PASS |
| Dibenz[a,h]anthracene        | Total    | 1.5                          | 1  | 0.001 | 0.005                      | µg/L       | 1.5   | 0                   | 100      | 55 - 150% | PASS             | 3      | 30       | PASS |
| Dibenzo[a,l]pyrene           | Total    | 0.52                         | 1  | 0.001 | 0.005                      | µg/L       | 0.5   | 0                   | 104      | 50 - 150% | PASS             | 4      | 30       | PASS |
| Dibenzothiophene             | Total    | 0.469                        | 1  | 0.001 | 0.005                      | µg/L       | 0.5   | 0                   | 94       | 46 - 126% | PASS             | 2      | 30       | PASS |



**Polynuclear Aromatic Hydrocarbons**

**QUALITY CONTROL REPORT**

| ANALYTE                | FRACTION | RESULT | DF | MDL   | RL    | UNITS | SPIKE | SOURCE | ACCURACY |           | PRECISION |        | QA CODE <sub>c</sub> |      |
|------------------------|----------|--------|----|-------|-------|-------|-------|--------|----------|-----------|-----------|--------|----------------------|------|
|                        |          |        |    |       |       |       | LEVEL | RESULT | %        | LIMITS    | %         | LIMITS |                      |      |
| Fluoranthene           | Total    | 1.72   | 1  | 0.001 | 0.005 | µg/L  | 1.5   | 0      | 115      | 60 - 146% | PASS      | 3      | 30                   | PASS |
| Fluorene               | Total    | 1.64   | 1  | 0.001 | 0.005 | µg/L  | 1.5   | 0      | 109      | 58 - 131% | PASS      | 4      | 30                   | PASS |
| Indeno[1,2,3-cd]pyrene | Total    | 1.67   | 1  | 0.001 | 0.005 | µg/L  | 1.5   | 0      | 111      | 50 - 151% | PASS      | 4      | 30                   | PASS |
| Naphthalene            | Total    | 1.42   | 1  | 0.001 | 0.005 | µg/L  | 1.5   | 0      | 95       | 41 - 126% | PASS      | 3      | 30                   | PASS |
| Perylene               | Total    | 0.506  | 1  | 0.001 | 0.005 | µg/L  | 0.5   | 0      | 101      | 48 - 141% | PASS      | 5      | 30                   | PASS |
| Phenanthrene           | Total    | 1.6    | 1  | 0.001 | 0.005 | µg/L  | 1.5   | 0      | 107      | 67 - 127% | PASS      | 2      | 30                   | PASS |
| Pyrene                 | Total    | 1.65   | 1  | 0.001 | 0.005 | µg/L  | 1.5   | 0      | 110      | 54 - 156% | PASS      | 2      | 30                   | PASS |



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**PHYSIS**

**TENTATIVELY**

**IDENTIFIED COMPOUNDS**

ENVIRONMENTAL LABORATORIES, INC.

*Innovative Solutions for Nature*

Sample ID: 102965

| RT      | Area Pct | Concentration (ng/L) | Library/ID                             | Cas Number | Match Qual |
|---------|----------|----------------------|--|------------|------------|
| 36.1323 | 6.1715   | 1111                 | Anthracene-D10                         | 1517-22-2  | 92         |
| 10.8913 | 2.0404   | 367                  | Cyclopropane, 2-bromo-1,1,3-trimethyl- | 36617-00-2 | 90         |

Concentration estimated using the response for Anthracene-d10

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**Sample ID: 102966**

| RT      | Area Pct | Concentration (ng/L) | Library/ID                             | Cas Number | Match Qual |
|---------|----------|----------------------|--|------------|------------|
| 36.1345 | 5.6305   | 1111                 | Anthracene-D10-                        | 1517-22-2  | 92         |
| 10.8918 | 1.9795   | 391                  | Cyclopropane, 2-bromo-1,1,3-trimethyl- | 36617-00-2 | 88         |
| 32.8247 | 0.5459   | 108                  | Benzoic acid, 2-ethylhexyl ester       | 5444-75-7  | 97         |

Concentration estimated using the response for Anthracene-d10

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Sample ID: 102967

| RT      | Area Pct | Concentration (ng/L) | Library/ID                             | Cas Number | Match Qual |
|---------|----------|----------------------|--|------------|------------|
| 36.1332 | 5.8000   | 1111                 | Anthracene-D10-                        | 1517-22-2  | 92         |
| 10.8930 | 2.0143   | 386                  | Cyclopropane, 2-bromo-1,1,3-trimethyl- | 36617-00-2 | 90         |
| 32.8231 | 0.5242   | 100                  | Benzoic acid, 2-ethylhexyl ester       | 5444-75-7  | 97         |

Concentration estimated using the response for Anthracene-d10

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Sample ID: Lab Blank B1\_40074

| RT      | Area Pct | Concentration (ng/L) | Library/ID                            | Cas Number | Match Qual |
|---------|----------|----------------------|---------------------------------------|------------|------------|
| 36.1336 | 5.7677   | 1111                 | Anthracene-D10-                       | 1517-22-2  | 92         |
| 10.8911 | 2.2211   | 428                  | 1,5-Heptadien-4-one, 3,3,6-trimethyl- | 546-49-6   | 89         |

Concentration estimated using the response for Anthracene-d10

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# PERFORMANCE CHAIN OF CUSTODY

TERRA ENVIRONMENTAL LABORATORIES, INC. AURA

*Innovative Solutions for Nature*

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**ICOC No:**  
380-31643

**Containers**  
Count 6      Container Type Amber Glass 1 liter - Sodium Thiosulfate

Preservative  
Sodium Thiosulfate

**Subcontract Method Instructions**

| Sample IDs | Method      | Method Description   | Method Comments   |
|------------|-------------|--|-------------------|
| 1, 2, 3    | SUBCONTRACT | SUB (625 PAH Physis LL (EAL) + TICS)/ 625 PAH Physis LL (EAL) + TICS | PAH Fraction TICS |

Project Iteration ID: 1407003-351  
 Client Name: Eurofins Eaton Analytical  
 Project Name: RED-HILL Project # 38001111 Job # 380-32096-1  
 COC Page Number: 3 of 3  
 Bottle Label Color: NA

**Sample Receipt Summary**

Receiving Info

1. Initials Received By: NW
2. Date Received: 12/21
3. Time Received: 1605
4. Client Name: Eurofins
5. Courier Information: (Please circle)
  - Client
    - UPS
    - **Area Fast**
    - DRS
  - FedEx
    - GSO/GLS
    - Ontrac
    - PAMS
  - PHYSIS Driver:
    - i. Start Time: \_\_\_\_\_
    - ii. End Time: \_\_\_\_\_
    - iii. Total Mileage: \_\_\_\_\_
    - iv. Number of Pickups: \_\_\_\_\_
6. Container Information: (Please put the # of containers or circle none)
  - 1 Cooler
  - \_\_\_ Styrofoam Cooler
  - \_\_\_ Boxes
  - None
  - \_\_\_ Carboy(s)
  - \_\_\_ Carboy Trash Can(s)
  - \_\_\_ Carboy Cap(s)
  - Other \_\_\_\_\_
7. What type of ice was used: (Please circle any that apply)
  - **Wet Ice**
  - Blue Ice
  - Dry Ice
  - Water
  - None
8. Randomly Selected Samples Temperature (°C): 0.7 Used I/R Thermometer # 1-2

Inspection Info

1. Initials Inspected By: RGH

Sample Integrity Upon Receipt:

1. COC(s) included and completely filled out..... Yes / No
2. All sample containers arrived intact..... Yes / No
3. All samples listed on COC(s) are present..... Yes / No
4. Information on containers consistent with information on COC(s)..... Yes / No
5. Correct containers and volume for all analyses indicated..... Yes / No
6. All samples received within method holding time..... Yes / No
7. Correct preservation used for all analyses indicated..... Yes / No
8. Name of sampler included on COC(s)..... Yes / No

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# Login Sample Receipt Checklist

Client: City & County of Honolulu

Job Number: 380-32096-1

**Login Number: 32096**  
**List Number: 1**  
**Creator: Elyas, Matthew**

**List Source: Eurofins Eaton Monrovia**

| Question   | Answer | Comment                                |
|--|--------|--|
| The cooler's custody seal, if present, is intact.                                | True   |  |
| Sample custody seals, if present, are intact.                                    | True   |  |
| Samples were received on ice.  | True   |  |
| Cooler Temperature is acceptable.  | True   |  |
| Cooler Temperature is recorded.  | True   |  |
| COC is present.  | True   |  |
| COC is filled out in ink and legible.  | True   |  |
| COC is filled out with all pertinent information.                                | True   |  |
| There are no discrepancies between the containers received and the COC.          | False  | The sample IDs don't completely match. |
| Samples are received within Holding Time (excluding tests with immediate HTs)    | True   |  |
| Sample containers have legible labels.   | True   |  |
| Containers are not broken or leaking.  | True   |  |
| Sample collection date/times are provided.                                       | True   |  |
| There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs | True   |  |
| Containers requiring zero headspace have no headspace or bubble is <6mm (1/4").  | True   |  |
| Samples do not require splitting or compositing.                                 | True   |  |
| Container provided by EEA  | True   |  |

