

# ANALYTICAL REPORT

## PREPARED FOR

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Honolulu, Hawaii 96843

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## JOB DESCRIPTION

RED-HILL  
RUSH Weekly Red Hill

## JOB NUMBER

380-57979-1

# Eurofins Eaton Analytical Pomona

## Job Notes

This report may not be reproduced except in full, and with written approval from the laboratory. The results relate only to the samples tested. For questions please contact the Project Manager at the e-mail address or telephone number listed on this page.

The test results in this report relate only to the samples as received by the laboratory and meet all requirements of the methodology, with any exceptions noted. This report shall not be reproduced except in full, without the express written approval of the laboratory. All questions should be directed to the Eurofins Eaton Analytical, LLC Project Manager.

## 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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# Definitions/Glossary

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

Job ID: 380-57979-1

## Qualifiers

### GC/MS Semi VOA

| Qualifier | Qualifier Description  |
|-----------|--|
| *+        | LCS and/or LCSD is outside acceptance limits, high biased.   |
| *1        | LCS/LCSD RPD exceeds control limits.   |
| ^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                                     |
| N         | Presumptive evidence of material.   |
| T         | Result is a tentatively identified compound (TIC) and an estimated value. |

### LCMS

| Qualifier | Qualifier Description   |
|-----------|---|
| *5-       | Isotope dilution analyte is outside acceptance limits, low biased.  |
| 4         | MS, MSD: The analyte present in the original sample is greater than 4 times the matrix spike concentration; therefore, control limits are not applicable. |
| J         | Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value.  |

## Glossary

| Abbreviation   | These commonly used abbreviations may or may not be present in this report.                                 |
|----------------|---|
| α              | 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: 380-57979-1

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**Job ID: 380-57979-1**

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**Laboratory: Eurofins Eaton Analytical Pomona**

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**Narrative**

**Job Narrative**  
**380-57979-1**

**Comments**

No additional comments.

**Receipt**

The samples were received on 8/8/2023 10:00 AM. Unless otherwise noted below, the samples arrived in good condition, and where required, properly preserved and on ice. The temperature of the cooler at receipt was 4.9° C.

**GC/MS Semi VOA**

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

**LCMS**

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

**Organic Prep**

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

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# Detection Summary

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

Job ID: 380-57979-1

**Client Sample ID: HALAWA WELLS UNITS 1 & 2**  
**PWSID Number: HI0000331**

**Lab Sample ID: 380-57979-1**

| Analyte                              | Result | Qualifier | RL  | Unit | Dil Fac | D | Method | Prep Type |
|--------------------------------------|--------|-----------|-----|------|---------|---|--------|-----------|
| Perfluorohexanesulfonic acid (PFHxS) | 2.3    |           | 2.0 | ng/L | 1       |   | 533    | Total/NA  |
| Perfluorohexanoic acid (PFHxA)       | 2.1    |           | 2.0 | ng/L | 1       |   | 533    | Total/NA  |
| Perfluorooctanesulfonic acid (PFOS)  | 2.3    |           | 2.0 | ng/L | 1       |   | 533    | Total/NA  |
| Perfluorooctanoic acid (PFOA)        | 2.1    |           | 2.0 | ng/L | 1       |   | 533    | Total/NA  |
| Perfluoropentanoic acid (PFPeA)      | 2.7    |           | 2.0 | ng/L | 1       |   | 533    | Total/NA  |
| Perfluorooctanesulfonic acid (PFOS)  | 2.1    |           | 2.0 | ng/L | 1       |   | 537.1  | Total/NA  |
| Perfluorohexanoic acid (PFHxA)       | 2.4    |           | 2.0 | ng/L | 1       |   | 537.1  | Total/NA  |
| Perfluorooctanoic acid (PFOA)        | 2.4    |           | 2.0 | ng/L | 1       |   | 537.1  | Total/NA  |
| Perfluorohexanesulfonic acid (PFHxS) | 2.5    |           | 2.0 | ng/L | 1       |   | 537.1  | Total/NA  |

**Client Sample ID: FB: HALAWA WELLS UNITS 1 & 2**  
**PWSID Number: HI0000331**

**Lab Sample ID: 380-57979-3**

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: 380-57979-1

**Client Sample ID: HALAWA WELLS UNITS 1 & 2**

**Lab Sample ID: 380-57979-1**

Date Collected: 08/02/23 10:00

Matrix: Drinking Water

Date Received: 08/08/23 10:00

PWSID Number: HI0000331

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

| Analyte                          | Result | Qualifier | RL    | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------------------------------|--------|-----------|-------|------|---|----------------|----------------|---------|
| 1-Methylnaphthalene              | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| 2,4'-DDD                         | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| 2,4'-DDE                         | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| 2,4'-DDT                         | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| 2,4-Dinitrotoluene               | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| 2,6-Dinitrotoluene               | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| 2-Methylnaphthalene              | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| 4,4'-DDD                         | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| 4,4'-DDE                         | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| 4,4'-DDT                         | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Acenaphthene                     | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Acenaphthylene                   | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Acetochlor                       | <0.098 | *+        | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Alachlor                         | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| alpha-BHC                        | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| alpha-Chlordane                  | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Anthracene                       | <0.020 |           | 0.020 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Atrazine                         | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Benz(a)anthracene                | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Benzo[a]pyrene                   | <0.020 |           | 0.020 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Benzo[b]fluoranthene             | <0.020 |           | 0.020 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Benzo[g,h,i]perylene             | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Benzo[k]fluoranthene             | <0.020 |           | 0.020 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| beta-BHC                         | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Bis(2-ethylhexyl) phthalate      | <0.59  | *1        | 0.59  | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Bromacil                         | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Butachlor                        | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Butylbenzylphthalate             | <0.49  |           | 0.49  | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Chlorobenzilate                  | <0.098 | ^3+       | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Chloroneb                        | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Chlorothalonil (Draconil, Bravo) | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Chlorpyrifos                     | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Chrysene                         | <0.020 |           | 0.020 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| delta-BHC                        | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Di(2-ethylhexyl)adipate          | <0.59  |           | 0.59  | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Dibenz(a,h)anthracene            | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Diclorvos (DDVP)                 | <0.049 | *+ ^3+    | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Dieldrin                         | <0.20  |           | 0.20  | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Diethylphthalate                 | <0.49  |           | 0.49  | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Dimethylphthalate                | <0.49  |           | 0.49  | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Di-n-butyl phthalate             | <0.98  |           | 0.98  | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Di-n-octyl phthalate             | <0.098 | *1        | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Endosulfan I (Alpha)             | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Endosulfan II (Beta)             | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Endosulfan sulfate               | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Endrin                           | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Endrin aldehyde                  | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| EPTC                             | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Fluoranthene                     | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |

Eurofins Eaton Analytical Pomona

# Client Sample Results

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

Job ID: 380-57979-1

**Client Sample ID: HALAWA WELLS UNITS 1 & 2**

**Lab Sample ID: 380-57979-1**

**Date Collected: 08/02/23 10:00**

**Matrix: Drinking Water**

**Date Received: 08/08/23 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 |
|----------------------------------|--------|-----------|-------|------|---|----------------|----------------|---------|
| Fluorene                         | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| gamma-Chlordane                  | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Heptachlor                       | <0.039 |           | 0.039 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Heptachlor epoxide (isomer B)    | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Hexachlorobenzene                | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Hexachlorocyclopentadiene        | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Indeno[1,2,3-cd]pyrene           | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Isophorone                       | <0.49  |           | 0.49  | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Lindane                          | <0.039 |           | 0.039 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Malathion                        | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Methoxychlor                     | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Metolachlor                      | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Molinate                         | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Naphthalene                      | <0.29  |           | 0.29  | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Parathion                        | <0.098 | *+        | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Pendimethalin (Penoxaline)       | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Phenanthrene                     | <0.039 |           | 0.039 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Propachlor                       | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Pyrene                           | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Simazine                         | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Terbacil                         | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Terbutylazine                    | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Thiobencarb                      | <0.20  |           | 0.20  | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Total Permethrin (mixed isomers) | <0.20  |           | 0.20  | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| trans-Nonachlor                  | <0.049 |           | 0.049 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Trifluralin                      | <0.098 |           | 0.098 | ug/L |   | 08/10/23 08:15 | 08/11/23 11:56 | 1       |

| Tentatively Identified Compound | Est. Result | Qualifier | Unit | D | RT | CAS No. | Prepared       | Analyzed       | Dil Fac |
|---------------------------------|-------------|-----------|------|---|----|---------|----------------|----------------|---------|
| Tentatively Identified Compound | None        |           | ug/L |   |    | N/A     | 08/10/23 08:15 | 08/11/23 11:56 | 1       |

| Surrogate          | %Recovery | Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|--------------------|-----------|-----------|----------|----------------|----------------|---------|
| 2-Nitro-m-xylene   | 101       |           | 70 - 130 | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Perylene-d12       | 88        |           | 70 - 130 | 08/10/23 08:15 | 08/11/23 11:56 | 1       |
| Triphenylphosphate | 113       |           | 70 - 130 | 08/10/23 08:15 | 08/11/23 11:56 | 1       |

**Method: EPA 533 - Perfluorinated and Polyfluorinated Alkyl Substances in Drinking Water**

| Analyte   | Result     | Qualifier | RL  | Unit | D | Prepared       | Analyzed       | Dil Fac |
|---|------------|-----------|-----|------|---|----------------|----------------|---------|
| 11-Chloroeicosafiuoro-3-oxaundecan e-1-sulfonic acid (11Cl-PF3OUdS) | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 9-Chlorohexadecafluoro-3-oxanonan e-1-sulfonic acid(9Cl-PF3ONS)     | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA)                         | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)                 | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| Perfluorobutanesulfonic acid (PFBS)                                 | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| Perfluorodecanoic acid (PFDA)                                       | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| Perfluorododecanoic acid (PFDoA)                                    | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| Perfluoroheptanoic acid (PFHpA)                                     | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| <b>Perfluorohexanesulfonic acid (PFHxS)</b>                         | <b>2.3</b> |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |

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# Client Sample Results

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

Job ID: 380-57979-1

**Client Sample ID: HALAWA WELLS UNITS 1 & 2**

**Lab Sample ID: 380-57979-1**

Date Collected: 08/02/23 10:00

Matrix: Drinking Water

Date Received: 08/08/23 10:00

PWSID Number: HI0000331

**Method: EPA 533 - Perfluorinated and Polyfluorinated Alkyl Substances in Drinking Water (Continued)**

| Analyte   | Result     | Qualifier | RL  | Unit | D | Prepared       | Analyzed       | Dil Fac |
|---|------------|-----------|-----|------|---|----------------|----------------|---------|
| <b>Perfluorohexanoic acid (PFHxA)</b>               | <b>2.1</b> |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| Perfluorononanoic acid (PFNA)                       | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| <b>Perfluorooctanesulfonic acid (PFOS)</b>          | <b>2.3</b> |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| <b>Perfluorooctanoic acid (PFOA)</b>                | <b>2.1</b> |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| Perfluoroundecanoic acid (PFUnA)                    | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| Perfluorobutanoic acid (PFBA)                       | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS) | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS) | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS) | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)          | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| Perfluoro (2-ethoxyethane) sulfonic acid (PFEEESA)  | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| Perfluoro-3-methoxypropanoic acid (PFMPA)           | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| Perfluoro-4-methoxybutanoic acid (PFMBA)            | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| <b>Perfluoropentanoic acid (PFPeA)</b>              | <b>2.7</b> |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| Perfluoroheptanesulfonic acid (PFHpS)               | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| Perfluoropentanesulfonic acid (PFPeS)               | <2.0       |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 06:48 | 1       |

| Isotope Dilution | %Recovery | Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|------------------|-----------|-----------|----------|----------------|----------------|---------|
| 13C3 HFPO-DA     | 75        |           | 50 - 200 | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 13C6 PFDA        | 97        |           | 50 - 200 | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 13C5 PFHxA       | 89        |           | 50 - 200 | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 13C4 PFHpA       | 92        |           | 50 - 200 | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 13C8 PFOA        | 94        |           | 50 - 200 | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 13C9 PFNA        | 97        |           | 50 - 200 | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 13C7 PFUnA       | 97        |           | 50 - 200 | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 13C2 PFDoA       | 98        |           | 50 - 200 | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 13C4 PFBA        | 89        |           | 50 - 200 | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 13C5 PFPeA       | 91        |           | 50 - 200 | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 13C3 PFBS        | 96        |           | 50 - 200 | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 13C3 PFHxS       | 98        |           | 50 - 200 | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 13C8 PFOS        | 98        |           | 50 - 200 | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 13C2-4:2-FTS     | 117       |           | 50 - 200 | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 13C2-6:2-FTS     | 101       |           | 50 - 200 | 08/24/23 16:20 | 08/27/23 06:48 | 1       |
| 13C2-8:2-FTS     | 97        |           | 50 - 200 | 08/24/23 16:20 | 08/27/23 06:48 | 1       |

**Method: EPA 537.1 - Perfluorinated Alkyl Acids (LC/MS)**

| Analyte  | Result     | Qualifier | RL  | Unit | D | Prepared       | Analyzed       | Dil Fac |
|--|------------|-----------|-----|------|---|----------------|----------------|---------|
| Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)      | <2.0       |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/12/23 11:03 | 1       |
| <b>Perfluorooctanesulfonic acid (PFOS)</b>               | <b>2.1</b> |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/12/23 11:03 | 1       |
| Perfluoroundecanoic acid (PFUnA)                         | <2.0       |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/12/23 11:03 | 1       |
| N-methylperfluorooctanesulfonamide cetic acid (NMeFOSAA) | <2.0       |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/12/23 11:03 | 1       |

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# Client Sample Results

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

Job ID: 380-57979-1

**Client Sample ID: HALAWA WELLS UNITS 1 & 2**

**Lab Sample ID: 380-57979-1**

Date Collected: 08/02/23 10:00

Matrix: Drinking Water

Date Received: 08/08/23 10:00

PWSID Number: HI0000331

**Method: EPA 537.1 - Perfluorinated Alkyl Acids (LC/MS) (Continued)**

| Analyte   | Result           | Qualifier        | RL            | Unit | D | Prepared        | Analyzed        | Dil Fac        |
|---|------------------|------------------|---------------|------|---|-----------------|-----------------|----------------|
| N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)           | <2.0             |                  | 2.0           | ng/L |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |
| <b>Perfluorohexanoic acid (PFHxA)</b>                             | <b>2.4</b>       |                  | 2.0           | ng/L |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |
| Perfluorododecanoic acid (PFDoA)                                  | <2.0             |                  | 2.0           | ng/L |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |
| <b>Perfluorooctanoic acid (PFOA)</b>                              | <b>2.4</b>       |                  | 2.0           | ng/L |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |
| Perfluorodecanoic acid (PFDA)                                     | <2.0             |                  | 2.0           | ng/L |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |
| <b>Perfluorohexanesulfonic acid (PFHxS)</b>                       | <b>2.5</b>       |                  | 2.0           | ng/L |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |
| Perfluorobutanesulfonic acid (PFBS)                               | <2.0             |                  | 2.0           | ng/L |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |
| Perfluoroheptanoic acid (PFHpA)                                   | <2.0             |                  | 2.0           | ng/L |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |
| Perfluorononanoic acid (PFNA)                                     | <2.0             |                  | 2.0           | ng/L |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |
| Perfluorotetradecanoic acid (PFTA)                                | <2.0             |                  | 2.0           | ng/L |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |
| Perfluorotridecanoic acid (PFTrDA)                                | <2.0             |                  | 2.0           | ng/L |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |
| 9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid(9Cl-PF3ONS)    | <2.0             |                  | 2.0           | ng/L |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |
| 11-Chloroeicosfluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS) | <2.0             |                  | 2.0           | ng/L |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA)                       | <2.0             |                  | 2.0           | ng/L |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |
| <b>Surrogate</b>  | <b>%Recovery</b> | <b>Qualifier</b> | <b>Limits</b> |      |   | <b>Prepared</b> | <b>Analyzed</b> | <b>Dil Fac</b> |
| d5-NEtFOSAA   | 100              |                  | 70 - 130      |      |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |
| 13C2 PFHxA  | 127              |                  | 70 - 130      |      |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |
| 13C2 PFDA   | 121              |                  | 70 - 130      |      |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |
| 13C3-GenX   | 129              |                  | 70 - 130      |      |   | 08/10/23 05:33  | 08/12/23 11:03  | 1              |

**Client Sample ID: FB: HALAWA WELLS UNITS 1 & 2**

**Lab Sample ID: 380-57979-3**

Date Collected: 08/02/23 10:00

Matrix: Water

Date Received: 08/08/23 10:00

PWSID Number: HI0000331

**Method: EPA 533 - Perfluorinated and Polyfluorinated Alkyl Substances in Drinking Water**

| Analyte   | Result | Qualifier | RL  | Unit | D | Prepared       | Analyzed       | Dil Fac |
|---|--------|-----------|-----|------|---|----------------|----------------|---------|
| 11-Chloroeicosfluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS) | <2.0   |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid(9Cl-PF3ONS)    | <2.0   |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA)                       | <2.0   |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)               | <2.0   |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Perfluorobutanesulfonic acid (PFBS)                               | <2.0   |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Perfluorodecanoic acid (PFDA)                                     | <2.0   |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Perfluorododecanoic acid (PFDoA)                                  | <2.0   |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Perfluoroheptanoic acid (PFHpA)                                   | <2.0   |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Perfluorohexanesulfonic acid (PFHxS)                              | <2.0   |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Perfluorohexanoic acid (PFHxA)                                    | <2.0   |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Perfluorononanoic acid (PFNA)                                     | <2.0   |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Perfluorooctanesulfonic acid (PFOS)                               | <2.0   |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Perfluorooctanoic acid (PFOA)                                     | <2.0   |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Perfluoroundecanoic acid (PFUnA)                                  | <2.0   |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Perfluorobutanoic acid (PFBA)                                     | <2.0   |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)               | <2.0   |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |

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# Client Sample Results

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

Job ID: 380-57979-1

**Client Sample ID: FB: HALAWA WELLS UNITS 1 & 2**

**Lab Sample ID: 380-57979-3**

**Date Collected: 08/02/23 10:00**

**Matrix: Water**

**Date Received: 08/08/23 10:00**

**PWSID Number: HI0000331**

**Method: EPA 533 - Perfluorinated and Polyfluorinated Alkyl Substances in Drinking Water (Continued)**

| Analyte   | Result    | Qualifier | RL       | Unit | D | Prepared       | Analyzed       | Dil Fac |
|---|-----------|-----------|----------|------|---|----------------|----------------|---------|
| 1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS) | <2.0      |           | 2.0      | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS) | <2.0      |           | 2.0      | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)          | <2.0      |           | 2.0      | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Perfluoro (2-ethoxyethane) sulfonic acid (PFEEESA)  | <2.0      |           | 2.0      | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Perfluoro-3-methoxypropanoic acid (PFMPA)           | <2.0      |           | 2.0      | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Perfluoro-4-methoxybutanoic acid (PFMBA)            | <2.0      |           | 2.0      | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Perfluoropentanoic acid (PFPeA)                     | <2.0      |           | 2.0      | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Perfluoroheptanesulfonic acid (PFHpS)               | <2.0      |           | 2.0      | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Perfluoropentanesulfonic acid (PFPeS)               | <2.0      |           | 2.0      | ng/L |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| Isotope Dilution                                    | %Recovery | Qualifier | Limits   |      |   | Prepared       | Analyzed       | Dil Fac |
| 13C3 HFPO-DA  | 63        |           | 50 - 200 |      |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 13C6 PFDA   | 88        |           | 50 - 200 |      |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 13C5 PFHxA  | 71        |           | 50 - 200 |      |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 13C4 PFHpA  | 77        |           | 50 - 200 |      |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 13C8 PFOA   | 83        |           | 50 - 200 |      |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 13C9 PFNA   | 88        |           | 50 - 200 |      |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 13C7 PFUnA  | 93        |           | 50 - 200 |      |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 13C2 PFDoA  | 89        |           | 50 - 200 |      |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 13C4 PFBA   | 68        |           | 50 - 200 |      |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 13C5 PFPeA  | 70        |           | 50 - 200 |      |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 13C3 PFBS   | 92        |           | 50 - 200 |      |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 13C3 PFHxS  | 89        |           | 50 - 200 |      |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 13C8 PFOS   | 92        |           | 50 - 200 |      |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 13C2-4:2-FTS  | 97        |           | 50 - 200 |      |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 13C2-6:2-FTS  | 90        |           | 50 - 200 |      |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |
| 13C2-8:2-FTS  | 87        |           | 50 - 200 |      |   | 08/24/23 16:20 | 08/27/23 07:07 | 1       |

**Method: EPA 537.1 - Perfluorinated Alkyl Acids (LC/MS)**

| Analyte   | Result | Qualifier | RL  | Unit | D | Prepared       | Analyzed       | Dil Fac |
|---|--------|-----------|-----|------|---|----------------|----------------|---------|
| Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)       | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| Perfluorooctanesulfonic acid (PFOS)                       | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| Perfluoroundecanoic acid (PFUnA)                          | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| N-methylperfluorooctanesulfonamide acetic acid (NMeFOSAA) | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| N-ethylperfluorooctanesulfonamide acetic acid (NEtFOSAA)  | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| Perfluorohexanoic acid (PFHxA)                            | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| Perfluorododecanoic acid (PFDoA)                          | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| Perfluorooctanoic acid (PFOA)                             | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| Perfluorodecanoic acid (PFDA)                             | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| Perfluorohexanesulfonic acid (PFHxS)                      | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| Perfluorobutanesulfonic acid (PFBS)                       | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| Perfluoroheptanoic acid (PFHpA)                           | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |

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# Client Sample Results

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

Job ID: 380-57979-1

**Client Sample ID: FB: HALAWA WELLS UNITS 1 & 2**

**Lab Sample ID: 380-57979-3**

**Date Collected: 08/02/23 10:00**

**Matrix: Water**

**Date Received: 08/08/23 10:00**

**PWSID Number: HI0000331**

**Method: EPA 537.1 - Perfluorinated Alkyl Acids (LC/MS) (Continued)**

| Analyte   | Result | Qualifier | RL  | Unit | D | Prepared       | Analyzed       | Dil Fac |
|---|--------|-----------|-----|------|---|----------------|----------------|---------|
| Perfluorononanoic acid (PFNA)   | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| Perfluorotetradecanoic acid (PFTA)                                    | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| Perfluorotridecanoic acid (PFTrDA)                                    | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| 9-Chlorohexadecafluoro-3-oxanonan<br>e-1-sulfonic acid(9Cl-PF3ONS)    | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| 11-Chloroeicosafuoro-3-oxaundecan<br>e-1-sulfonic acid (11Cl-PF3OUdS) | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| 4,8-Dioxa-3H-perfluorononanoic acid<br>(ADONA)                        | <2.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 12:14 | 1       |

| Surrogate   | %Recovery | Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|-------------|-----------|-----------|----------|----------------|----------------|---------|
| d5-NEtFOSAA | 116       |           | 70 - 130 | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| 13C2 PFHxA  | 124       |           | 70 - 130 | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| 13C2 PFDA   | 110       |           | 70 - 130 | 08/10/23 05:33 | 08/11/23 12:14 | 1       |
| 13C3-GenX   | 112       |           | 70 - 130 | 08/10/23 05:33 | 08/11/23 12:14 | 1       |

# Action Limit Summary

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

Job ID: 380-57979-1

**Client Sample ID: HALAWA WELLS UNITS 1 & 2**  
**PWSID Number: HI0000331**

**Lab Sample ID: 380-57979-1**

## 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                      | <0.049 |           | ug/L | 2      | 0.049 | 525.2  | Total/NA  |
| Atrazine                      | <0.049 |           | ug/L | 3      | 0.049 | 525.2  | Total/NA  |
| Benzo[a]pyrene                | <0.020 |           | ug/L | 0.2    | 0.020 | 525.2  | Total/NA  |
| Bis(2-ethylhexyl) phthalate   | <0.59  | *1        | ug/L | 6      | 0.59  | 525.2  | Total/NA  |
| Di(2-ethylhexyl)adipate       | <0.59  |           | ug/L | 400    | 0.59  | 525.2  | Total/NA  |
| Endrin                        | <0.098 |           | ug/L | 2      | 0.098 | 525.2  | Total/NA  |
| Heptachlor                    | <0.039 |           | ug/L | 0.4    | 0.039 | 525.2  | Total/NA  |
| Heptachlor epoxide (isomer B) | <0.049 |           | ug/L | 0.2    | 0.049 | 525.2  | Total/NA  |
| Hexachlorobenzene             | <0.049 |           | ug/L | 1      | 0.049 | 525.2  | Total/NA  |
| Hexachlorocyclopentadiene     | <0.049 |           | ug/L | 50     | 0.049 | 525.2  | Total/NA  |
| Lindane                       | <0.039 |           | ug/L | 0.2    | 0.039 | 525.2  | Total/NA  |
| Methoxychlor                  | <0.098 |           | ug/L | 40     | 0.098 | 525.2  | Total/NA  |
| Simazine                      | <0.049 |           | ug/L | 4      | 0.049 | 525.2  | Total/NA  |

# Surrogate Summary

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

Job ID: 380-57979-1

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

Matrix: Drinking Water

Prep Type: Total/NA

### Percent Surrogate Recovery (Acceptance Limits)

| Lab Sample ID | Client Sample ID         | Percent Surrogate Recovery (Acceptance Limits) |                 |                 |
|---------------|--------------------------|--|-----------------|-----------------|
|               |                          | 2NMX<br>(70-130)                               | PRY<br>(70-130) | TPP<br>(70-130) |
| 380-57979-1   | HALAWA WELLS UNITS 1 & 2 | 101  | 88              | 113             |

#### Surrogate Legend

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

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

Matrix: Water

Prep Type: Total/NA

### Percent Surrogate Recovery (Acceptance Limits)

| Lab Sample ID       | Client Sample ID       | Percent Surrogate Recovery (Acceptance Limits) |                 |                 |
|---------------------|------------------------|--|-----------------|-----------------|
|                     |                        | 2NMX<br>(70-130)                               | PRY<br>(70-130) | TPP<br>(70-130) |
| 380-57897-P-1-A MS  | Matrix Spike           | 98   | 92              | 115             |
| 380-57909-AR-1-A DU | Duplicate              | 97   | 92              | 110             |
| LCS 380-51109/3-A   | Lab Control Sample     | 102  | 96              | 113             |
| LCSD 380-51109/4-A  | Lab Control Sample Dup | 103  | 95              | 104             |
| MB 380-51109/1-A    | Method Blank           | 101  | 92              | 115             |
| MRL 380-51109/2-A   | Lab Control Sample     | 106  | 97              | 108             |

#### Surrogate Legend

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

## Method: 537.1 - Perfluorinated Alkyl Acids (LC/MS)

Matrix: Drinking Water

Prep Type: Total/NA

### Percent Surrogate Recovery (Acceptance Limits)

| Lab Sample ID | Client Sample ID         | Percent Surrogate Recovery (Acceptance Limits) |                   |                  |                  |
|---------------|--------------------------|--|-------------------|------------------|------------------|
|               |                          | d5NEFOS<br>(70-130)                            | PFHxA<br>(70-130) | PFDA<br>(70-130) | GenX<br>(70-130) |
| 380-57979-1   | HALAWA WELLS UNITS 1 & 2 | 100  | 127               | 121              | 129              |

#### Surrogate Legend

d5NEFOS = d5-NEtFOSAA  
PFHxA = 13C2 PFHxA  
PFDA = 13C2 PFDA  
GenX = 13C3-GenX

## Method: 537.1 - Perfluorinated Alkyl Acids (LC/MS)

Matrix: Water

Prep Type: Total/NA

### Percent Surrogate Recovery (Acceptance Limits)

| Lab Sample ID        | Client Sample ID             | Percent Surrogate Recovery (Acceptance Limits) |                   |                  |                  |
|----------------------|------------------------------|--|-------------------|------------------|------------------|
|                      |                              | d5NEFOS<br>(70-130)                            | PFHxA<br>(70-130) | PFDA<br>(70-130) | GenX<br>(70-130) |
| 380-57979-3          | FB: HALAWA WELLS UNITS 1 & 2 | 116  | 124               | 110              | 112              |
| 380-58087-B-1-C LMS  | Matrix Spike                 | 117  | 129               | 119              | 119              |
| 380-58087-C-1-C LMSD | Matrix Spike Duplicate       | 114  | 117               | 108              | 107              |
| LCS 380-51108/25-A   | Lab Control Sample           | 108  | 120               | 112              | 110              |
| LCSD 380-51108/26-A  | Lab Control Sample Dup       | 110  | 128               | 113              | 130              |
| MBL 380-51108/23-A   | Method Blank                 | 109  | 115               | 121              | 106              |
| MRL 380-51108/24-A   | Lab Control Sample           | 97   | 122               | 110              | 113              |

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# Surrogate Summary

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

Job ID: 380-57979-1

## Surrogate Legend

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d5NEFOS = d5-NEtFOSAA

PFHxA = 13C2 PFHxA

PFDA = 13C2 PFDA

GenX = 13C3-GenX

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# Isotope Dilution Summary

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

Job ID: 380-57979-1

## Method: 533 - Perfluorinated and Polyfluorinated Alkyl Substances in Drinking Water

Matrix: Drinking Water

Prep Type: Total/NA

### Percent Isotope Dilution Recovery (Acceptance Limits)

| Lab Sample ID | Client Sample ID         | HFPODA<br>(50-200) | C6PFDA<br>(50-200) | 13C5PHA<br>(50-200) | C4PFHA<br>(50-200) | C8PFOA<br>(50-200) | C9PFNA<br>(50-200) | 13C7PUA<br>(50-200) | PFDaA<br>(50-200) |
|---------------|--------------------------|--------------------|--------------------|---------------------|--------------------|--------------------|--------------------|---------------------|-------------------|
| 380-57979-1   | HALAWA WELLS UNITS 1 & 2 | 75                 | 97                 | 89                  | 92                 | 94                 | 97                 | 97                  | 98                |

### Percent Isotope Dilution Recovery (Acceptance Limits)

| Lab Sample ID | Client Sample ID         | PFBA<br>(50-200) | PFPeA<br>(50-200) | C3PFBS<br>(50-200) | C3PFHS<br>(50-200) | C8PFOS<br>(50-200) | 42FTS<br>(50-200) | 62FTS<br>(50-200) | 82FTS<br>(50-200) |
|---------------|--------------------------|------------------|-------------------|--------------------|--------------------|--------------------|-------------------|-------------------|-------------------|
| 380-57979-1   | HALAWA WELLS UNITS 1 & 2 | 89               | 91                | 96                 | 98                 | 98                 | 117               | 101               | 97                |

#### Surrogate Legend

- HFPODA = 13C3 HFPO-DA
- C6PFDA = 13C6 PFDA
- 13C5PHA = 13C5 PFHxA
- C4PFHA = 13C4 PFHpA
- C8PFOA = 13C8 PFOA
- C9PFNA = 13C9 PFNA
- 13C7PUA = 13C7 PFUnA
- PFDaA = 13C2 PFDaA
- PFBA = 13C4 PFBA
- PFPeA = 13C5 PFPeA
- C3PFBS = 13C3 PFBS
- C3PFHS = 13C3 PFHxS
- C8PFOS = 13C8 PFOS
- 42FTS = 13C2-4:2-FTS
- 62FTS = 13C2-6:2-FTS
- 82FTS = 13C2-8:2-FTS

## Method: 533 - Perfluorinated and Polyfluorinated Alkyl Substances in Drinking Water

Matrix: Water

Prep Type: Total/NA

### Percent Isotope Dilution Recovery (Acceptance Limits)

| Lab Sample ID        | Client Sample ID             | HFPODA<br>(50-200) | C6PFDA<br>(50-200) | 13C5PHA<br>(50-200) | C4PFHA<br>(50-200) | C8PFOA<br>(50-200) | C9PFNA<br>(50-200) | 13C7PUA<br>(50-200) | PFDaA<br>(50-200) |
|----------------------|------------------------------|--------------------|--------------------|---------------------|--------------------|--------------------|--------------------|---------------------|-------------------|
| 380-57432-O-1-D LMS  | Matrix Spike                 | 48 *5-             | 77                 | 58                  | 61                 | 69                 | 72                 | 77                  | 87                |
| 380-57432-P-1-C LMSD | Matrix Spike Duplicate       | 68                 | 84                 | 78                  | 78                 | 85                 | 85                 | 85                  | 90                |
| 380-57979-3          | FB: HALAWA WELLS UNITS 1 & 2 | 63                 | 88                 | 71                  | 77                 | 83                 | 88                 | 93                  | 89                |
| LCS 380-53035/23-A   | Lab Control Sample           | 91                 | 95                 | 94                  | 96                 | 97                 | 99                 | 96                  | 98                |
| LCSD 380-53035/24-A  | Lab Control Sample Dup       | 86                 | 98                 | 85                  | 93                 | 96                 | 97                 | 96                  | 97                |
| MBL 380-53035/21-A   | Method Blank                 | 79                 | 95                 | 88                  | 88                 | 93                 | 91                 | 88                  | 91                |
| MRL 380-53035/22-A   | Lab Control Sample           | 82                 | 93                 | 90                  | 90                 | 95                 | 93                 | 91                  | 93                |

### Percent Isotope Dilution Recovery (Acceptance Limits)

| Lab Sample ID        | Client Sample ID             | PFBA<br>(50-200) | PFPeA<br>(50-200) | C3PFBS<br>(50-200) | C3PFHS<br>(50-200) | C8PFOS<br>(50-200) | 42FTS<br>(50-200) | 62FTS<br>(50-200) | 82FTS<br>(50-200) |
|----------------------|------------------------------|------------------|-------------------|--------------------|--------------------|--------------------|-------------------|-------------------|-------------------|
| 380-57432-O-1-D LMS  | Matrix Spike                 | 63               | 57                | 92                 | 95                 | 93                 | 103               | 96                | 94                |
| 380-57432-P-1-C LMSD | Matrix Spike Duplicate       | 83               | 78                | 88                 | 98                 | 97                 | 112               | 97                | 99                |
| 380-57979-3          | FB: HALAWA WELLS UNITS 1 & 2 | 68               | 70                | 92                 | 89                 | 92                 | 97                | 90                | 87                |
| LCS 380-53035/23-A   | Lab Control Sample           | 93               | 94                | 92                 | 96                 | 95                 | 95                | 93                | 90                |
| LCSD 380-53035/24-A  | Lab Control Sample Dup       | 84               | 86                | 95                 | 97                 | 96                 | 106               | 94                | 94                |
| MBL 380-53035/21-A   | Method Blank                 | 92               | 90                | 90                 | 90                 | 91                 | 101               | 98                | 109               |
| MRL 380-53035/22-A   | Lab Control Sample           | 94               | 94                | 87                 | 88                 | 89                 | 94                | 87                | 88                |

#### Surrogate Legend

Eurofins Eaton Analytical Pomona



# Isotope Dilution Summary

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

Job ID: 380-57979-1

HFPODA = 13C3 HFPO-DA  
C6PFDA = 13C6 PFDA  
13C5PHA = 13C5 PFHxA  
C4PFHA = 13C4 PFHpA  
C8PFOA = 13C8 PFOA  
C9PFNA = 13C9 PFNA  
13C7PUA = 13C7 PFUnA  
PFDoA = 13C2 PFDoA  
PFBA = 13C4 PFBA  
PFPeA = 13C5 PFPeA  
C3PFBS = 13C3 PFBS  
C3PFHS = 13C3 PFHxS  
C8PFOS = 13C8 PFOS  
42FTS = 13C2-4:2-FTS  
62FTS = 13C2-6:2-FTS  
82FTS = 13C2-8:2-FTS

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- 3
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# QC Sample Results

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

Job ID: 380-57979-1

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

**Lab Sample ID: MB 380-51109/1-A**  
**Matrix: Water**  
**Analysis Batch: 51305**

**Client Sample ID: Method Blank**  
**Prep Type: Total/NA**  
**Prep Batch: 51109**

| Analyte                          | MB     | MB        | RL    | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------------------------------|--------|-----------|-------|------|---|----------------|----------------|---------|
|                                  | Result | Qualifier |       |      |   |                |                |         |
| 1-Methylnaphthalene              | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| 2,4'-DDD                         | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| 2,4'-DDE                         | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| 2,4'-DDT                         | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| 2,4-Dinitrotoluene               | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| 2,6-Dinitrotoluene               | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| 2-Methylnaphthalene              | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| 4,4'-DDD                         | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| 4,4'-DDE                         | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| 4,4'-DDT                         | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Acenaphthene                     | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Acenaphthylene                   | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Acetochlor                       | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Alachlor                         | <0.050 |           | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| alpha-BHC                        | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| alpha-Chlordane                  | <0.050 |           | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Anthracene                       | <0.020 |           | 0.020 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Atrazine                         | <0.050 |           | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Benz(a)anthracene                | <0.050 |           | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Benzo[a]pyrene                   | <0.020 |           | 0.020 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Benzo[b]fluoranthene             | <0.020 |           | 0.020 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Benzo[g,h,i]perylene             | <0.050 |           | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Benzo[k]fluoranthene             | <0.020 |           | 0.020 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| beta-BHC                         | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Bis(2-ethylhexyl) phthalate      | <0.60  |           | 0.60  | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Bromacil                         | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Butachlor                        | <0.050 |           | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Butylbenzylphthalate             | <0.50  |           | 0.50  | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Chlorobenzilate                  | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Chloroneb                        | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Chlorothalonil (Draconil, Bravo) | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Chlorpyrifos                     | <0.050 |           | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Chrysene                         | <0.020 |           | 0.020 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| delta-BHC                        | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Di(2-ethylhexyl)adipate          | <0.60  |           | 0.60  | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Dibenz(a,h)anthracene            | <0.050 |           | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Diclorvos (DDVP)                 | <0.050 |           | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Dieldrin                         | <0.20  |           | 0.20  | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Diethylphthalate                 | <0.50  |           | 0.50  | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Dimethylphthalate                | <0.50  |           | 0.50  | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Di-n-butyl phthalate             | <0.99  |           | 0.99  | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Di-n-octyl phthalate             | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Endosulfan I (Alpha)             | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Endosulfan II (Beta)             | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Endosulfan sulfate               | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Endrin                           | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Endrin aldehyde                  | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| EPTC                             | <0.099 |           | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |

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# QC Sample Results

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

Job ID: 380-57979-1

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

**Lab Sample ID: MB 380-51109/1-A**  
**Matrix: Water**  
**Analysis Batch: 51305**

**Client Sample ID: Method Blank**  
**Prep Type: Total/NA**  
**Prep Batch: 51109**

| Analyte                          | MB Result | MB Qualifier | RL    | Unit | D | Prepared       | Analyzed       | Dil Fac |
|----------------------------------|-----------|--------------|-------|------|---|----------------|----------------|---------|
| Fluoranthene                     | <0.099    |              | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Fluorene                         | <0.050    |              | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| gamma-Chlordane                  | <0.050    |              | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Heptachlor                       | <0.040    |              | 0.040 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Heptachlor epoxide (isomer B)    | <0.050    |              | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Hexachlorobenzene                | <0.050    |              | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Hexachlorocyclopentadiene        | <0.050    |              | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Indeno[1,2,3-cd]pyrene           | <0.050    |              | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Isophorone                       | <0.50     |              | 0.50  | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Lindane                          | <0.040    |              | 0.040 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Malathion                        | <0.099    |              | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Methoxychlor                     | <0.099    |              | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Metolachlor                      | <0.050    |              | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Molinate                         | <0.099    |              | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Naphthalene                      | <0.30     |              | 0.30  | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Parathion                        | <0.099    |              | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Pendimethalin (Penoxaline)       | <0.099    |              | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Phenanthrene                     | <0.040    |              | 0.040 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Propachlor                       | <0.050    |              | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Pyrene                           | <0.050    |              | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Simazine                         | <0.050    |              | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Terbacil                         | <0.099    |              | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Terbutylazine                    | <0.099    |              | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Thiobencarb                      | <0.20     |              | 0.20  | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Total Permethrin (mixed isomers) | <0.20     |              | 0.20  | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| trans-Nonachlor                  | <0.050    |              | 0.050 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Trifluralin                      | <0.099    |              | 0.099 | ug/L |   | 08/10/23 07:00 | 08/11/23 09:16 | 1       |

| Tentatively Identified Compound | MB Est. Result | MB Qualifier | Unit | D | RT   | CAS No. | Prepared       | Analyzed       | Dil Fac |
|---------------------------------|----------------|--------------|------|---|------|---------|----------------|----------------|---------|
| Phenol, 4-(1,1-dimethylpropyl)- | 1.30           | T J N        | ug/L |   | 3.87 | 80-46-6 | 08/10/23 07:00 | 08/11/23 09:16 | 1       |

| Surrogate          | MB %Recovery | MB Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|--------------------|--------------|--------------|----------|----------------|----------------|---------|
| 2-Nitro-m-xylene   | 101          |              | 70 - 130 | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Perylene-d12       | 92           |              | 70 - 130 | 08/10/23 07:00 | 08/11/23 09:16 | 1       |
| Triphenylphosphate | 115          |              | 70 - 130 | 08/10/23 07:00 | 08/11/23 09:16 | 1       |

**Lab Sample ID: LCS 380-51109/3-A**  
**Matrix: Water**  
**Analysis Batch: 51305**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 51109**

| Analyte             | Spike Added | LCS Result | LCS Qualifier | Unit | D | %Rec | %Rec Limits |
|---------------------|-------------|------------|---------------|------|---|------|-------------|
| 1-Methylnaphthalene | 1.99        | 1.99       |               | ug/L |   | 100  | 70 - 130    |
| 2,4'-DDD            | 1.99        | 2.13       |               | ug/L |   | 107  | 70 - 130    |
| 2,4'-DDE            | 1.99        | 1.99       |               | ug/L |   | 100  | 70 - 130    |
| 2,4'-DDT            | 1.99        | 2.18       |               | ug/L |   | 110  | 70 - 130    |
| 2,4-Dinitrotoluene  | 1.99        | 2.16       |               | ug/L |   | 109  | 70 - 130    |
| 2,6-Dinitrotoluene  | 1.99        | 2.09       |               | ug/L |   | 105  | 70 - 130    |
| 2-Methylnaphthalene | 1.99        | 1.99       |               | ug/L |   | 100  | 70 - 130    |

Eurofins Eaton Analytical Pomona

# QC Sample Results

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

Job ID: 380-57979-1

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

**Lab Sample ID: LCS 380-51109/3-A**  
**Matrix: Water**  
**Analysis Batch: 51305**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 51109**

| Analyte                          | Spike Added | LCS Result | LCS Qualifier | Unit | D | %Rec | %Rec Limits |
|----------------------------------|-------------|------------|---------------|------|---|------|-------------|
| 4,4'-DDD                         | 1.99        | 2.08       |               | ug/L |   | 105  | 70 - 130    |
| 4,4'-DDE                         | 1.99        | 2.01       |               | ug/L |   | 101  | 70 - 130    |
| 4,4'-DDT                         | 1.99        | 2.24       |               | ug/L |   | 113  | 70 - 130    |
| Acenaphthene                     | 1.99        | 1.87       |               | ug/L |   | 94   | 70 - 130    |
| Acenaphthylene                   | 1.99        | 2.05       |               | ug/L |   | 103  | 70 - 130    |
| Acetochlor                       | 1.99        | 2.63       | *+            | ug/L |   | 132  | 70 - 130    |
| Alachlor                         | 1.99        | 2.34       |               | ug/L |   | 118  | 70 - 130    |
| alpha-BHC                        | 1.99        | 2.09       |               | ug/L |   | 105  | 70 - 130    |
| alpha-Chlordane                  | 1.99        | 2.03       |               | ug/L |   | 102  | 70 - 130    |
| Anthracene                       | 1.99        | 1.93       |               | ug/L |   | 97   | 70 - 130    |
| Atrazine                         | 1.99        | 2.16       |               | ug/L |   | 109  | 70 - 130    |
| Benz(a)anthracene                | 1.99        | 2.13       |               | ug/L |   | 107  | 70 - 130    |
| Benzo[a]pyrene                   | 1.99        | 2.06       |               | ug/L |   | 104  | 70 - 130    |
| Benzo[b]fluoranthene             | 1.99        | 2.10       |               | ug/L |   | 105  | 70 - 130    |
| Benzo[g,h,i]perylene             | 1.99        | 2.25       |               | ug/L |   | 113  | 70 - 130    |
| Benzo[k]fluoranthene             | 1.99        | 2.20       |               | ug/L |   | 111  | 70 - 130    |
| beta-BHC                         | 1.99        | 2.14       |               | ug/L |   | 108  | 70 - 130    |
| Bis(2-ethylhexyl) phthalate      | 1.99        | 2.19       |               | ug/L |   | 110  | 70 - 130    |
| Bromacil                         | 1.99        | 2.34       |               | ug/L |   | 118  | 70 - 130    |
| Butachlor                        | 1.99        | 2.56       |               | ug/L |   | 129  | 70 - 130    |
| Butylbenzylphthalate             | 1.99        | 2.18       |               | ug/L |   | 110  | 70 - 130    |
| Chlorobenzilate                  | 1.99        | 2.53       |               | ug/L |   | 128  | 70 - 130    |
| Chloroneb                        | 1.99        | 2.06       |               | ug/L |   | 104  | 70 - 130    |
| Chlorothalonil (Draconil, Bravo) | 1.99        | 2.01       |               | ug/L |   | 101  | 70 - 130    |
| Chlorpyrifos                     | 1.99        | 2.27       |               | ug/L |   | 114  | 70 - 130    |
| Chrysene                         | 1.99        | 2.05       |               | ug/L |   | 103  | 70 - 130    |
| delta-BHC                        | 1.99        | 2.03       |               | ug/L |   | 102  | 70 - 130    |
| Di(2-ethylhexyl)adipate          | 1.99        | 2.18       |               | ug/L |   | 110  | 70 - 130    |
| Dibenz(a,h)anthracene            | 1.99        | 2.23       |               | ug/L |   | 112  | 70 - 130    |
| Diclorvos (DDVP)                 | 1.99        | 2.70       | *+            | ug/L |   | 136  | 70 - 130    |
| Dieldrin                         | 1.99        | 2.02       |               | ug/L |   | 102  | 70 - 130    |
| Diethylphthalate                 | 1.99        | 2.32       |               | ug/L |   | 117  | 70 - 130    |
| Dimethylphthalate                | 1.99        | 2.17       |               | ug/L |   | 109  | 70 - 130    |
| Di-n-butyl phthalate             | 3.97        | 4.36       |               | ug/L |   | 110  | 70 - 130    |
| Di-n-octyl phthalate             | 1.99        | 2.14       |               | ug/L |   | 107  | 70 - 130    |
| Endosulfan I (Alpha)             | 1.99        | 1.90       |               | ug/L |   | 95   | 70 - 130    |
| Endosulfan II (Beta)             | 1.99        | 2.17       |               | ug/L |   | 109  | 70 - 130    |
| Endosulfan sulfate               | 1.99        | 2.03       |               | ug/L |   | 102  | 70 - 130    |
| Endrin                           | 1.99        | 2.38       |               | ug/L |   | 120  | 70 - 130    |
| Endrin aldehyde                  | 1.99        | 2.15       |               | ug/L |   | 108  | 70 - 130    |
| EPTC                             | 1.99        | 2.29       |               | ug/L |   | 115  | 70 - 130    |
| Fluoranthene                     | 1.99        | 2.23       |               | ug/L |   | 112  | 70 - 130    |
| Fluorene                         | 1.99        | 2.15       |               | ug/L |   | 108  | 70 - 130    |
| gamma-Chlordane                  | 1.99        | 2.07       |               | ug/L |   | 104  | 70 - 130    |
| Heptachlor                       | 1.99        | 2.14       |               | ug/L |   | 108  | 70 - 130    |
| Heptachlor epoxide (isomer B)    | 1.99        | 2.17       |               | ug/L |   | 109  | 70 - 130    |
| Hexachlorobenzene                | 1.99        | 1.91       |               | ug/L |   | 96   | 70 - 130    |
| Hexachlorocyclopentadiene        | 1.99        | 1.91       |               | ug/L |   | 96   | 70 - 130    |
| Indeno[1,2,3-cd]pyrene           | 1.99        | 2.19       |               | ug/L |   | 110  | 70 - 130    |

Eurofins Eaton Analytical Pomona

# QC Sample Results

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

Job ID: 380-57979-1

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

**Lab Sample ID: LCS 380-51109/3-A**  
**Matrix: Water**  
**Analysis Batch: 51305**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 51109**

| Analyte                    | Spike Added | LCS Result | LCS Qualifier | Unit | D | %Rec | %Rec Limits |
|----------------------------|-------------|------------|---------------|------|---|------|-------------|
| Isophorone                 | 1.99        | 2.16       |               | ug/L |   | 109  | 70 - 130    |
| Lindane                    | 1.99        | 2.12       |               | ug/L |   | 107  | 70 - 130    |
| Malathion                  | 1.99        | 2.35       |               | ug/L |   | 119  | 70 - 130    |
| Methoxychlor               | 1.99        | 2.26       |               | ug/L |   | 114  | 70 - 130    |
| Metolachlor                | 1.99        | 2.45       |               | ug/L |   | 124  | 70 - 130    |
| Molinate                   | 1.99        | 2.27       |               | ug/L |   | 114  | 70 - 130    |
| Naphthalene                | 1.99        | 1.88       |               | ug/L |   | 95   | 70 - 130    |
| Parathion                  | 1.99        | 2.60       | *+            | ug/L |   | 131  | 70 - 130    |
| Pendimethalin (Penoxaline) | 1.99        | 2.21       |               | ug/L |   | 111  | 70 - 130    |
| Phenanthrene               | 1.99        | 1.92       |               | ug/L |   | 97   | 70 - 130    |
| Propachlor                 | 1.99        | 2.28       |               | ug/L |   | 115  | 70 - 130    |
| Pyrene                     | 1.99        | 2.20       |               | ug/L |   | 111  | 70 - 130    |
| Simazine                   | 1.99        | 2.27       |               | ug/L |   | 114  | 70 - 130    |
| Terbacil                   | 1.99        | 2.31       |               | ug/L |   | 116  | 70 - 130    |
| Terbutylazine              | 1.99        | 2.31       |               | ug/L |   | 116  | 70 - 130    |
| Thiobencarb                | 1.99        | 2.22       |               | ug/L |   | 112  | 70 - 130    |
| trans-Nonachlor            | 1.99        | 1.94       |               | ug/L |   | 97   | 70 - 130    |
| Trifluralin                | 1.99        | 2.08       |               | ug/L |   | 105  | 70 - 130    |

| Surrogate          | LCS %Recovery | LCS Qualifier | Limits   |
|--------------------|---------------|---------------|----------|
| 2-Nitro-m-xylene   | 102           |               | 70 - 130 |
| Perylene-d12       | 96            |               | 70 - 130 |
| Triphenylphosphate | 113           |               | 70 - 130 |

**Lab Sample ID: LCSD 380-51109/4-A**  
**Matrix: Water**  
**Analysis Batch: 51305**

**Client Sample ID: Lab Control Sample Dup**  
**Prep Type: Total/NA**  
**Prep Batch: 51109**

| Analyte             | Spike Added | LCSD Result | LCSD Qualifier | Unit | D | %Rec | %Rec Limits | RPD | RPD Limit |
|---------------------|-------------|-------------|----------------|------|---|------|-------------|-----|-----------|
| 1-Methylnaphthalene | 1.98        | 1.99        |                | ug/L |   | 100  | 70 - 130    | 0   | 20        |
| 2,4'-DDD            | 1.98        | 2.04        |                | ug/L |   | 103  | 70 - 130    | 4   | 20        |
| 2,4'-DDE            | 1.98        | 1.93        |                | ug/L |   | 97   | 70 - 130    | 3   | 20        |
| 2,4'-DDT            | 1.98        | 2.08        |                | ug/L |   | 105  | 70 - 130    | 5   | 20        |
| 2,4-Dinitrotoluene  | 1.98        | 2.22        |                | ug/L |   | 112  | 70 - 130    | 3   | 20        |
| 2,6-Dinitrotoluene  | 1.98        | 2.18        |                | ug/L |   | 110  | 70 - 130    | 4   | 20        |
| 2-Methylnaphthalene | 1.98        | 2.01        |                | ug/L |   | 101  | 70 - 130    | 1   | 20        |
| 4,4'-DDD            | 1.98        | 2.02        |                | ug/L |   | 102  | 70 - 130    | 3   | 20        |
| 4,4'-DDE            | 1.98        | 1.93        |                | ug/L |   | 98   | 70 - 130    | 4   | 20        |
| 4,4'-DDT            | 1.98        | 2.03        |                | ug/L |   | 103  | 70 - 130    | 10  | 20        |
| Acenaphthene        | 1.98        | 1.91        |                | ug/L |   | 96   | 70 - 130    | 2   | 20        |
| Acenaphthylene      | 1.98        | 2.05        |                | ug/L |   | 103  | 70 - 130    | 0   | 20        |
| Acetochlor          | 1.98        | 2.53        |                | ug/L |   | 128  | 70 - 130    | 4   | 20        |
| Alachlor            | 1.98        | 2.32        |                | ug/L |   | 117  | 70 - 130    | 1   | 20        |
| alpha-BHC           | 1.98        | 2.10        |                | ug/L |   | 106  | 70 - 130    | 0   | 20        |
| alpha-Chlordane     | 1.98        | 2.02        |                | ug/L |   | 102  | 70 - 130    | 1   | 20        |
| Anthracene          | 1.98        | 1.95        |                | ug/L |   | 98   | 70 - 130    | 1   | 20        |
| Atrazine            | 1.98        | 2.12        |                | ug/L |   | 107  | 70 - 130    | 2   | 20        |
| Benz(a)anthracene   | 1.98        | 2.01        |                | ug/L |   | 101  | 70 - 130    | 6   | 20        |

# QC Sample Results

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

Job ID: 380-57979-1

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

**Lab Sample ID: LCSD 380-51109/4-A**  
**Matrix: Water**  
**Analysis Batch: 51305**

**Client Sample ID: Lab Control Sample Dup**  
**Prep Type: Total/NA**  
**Prep Batch: 51109**

| Analyte                          | Spike Added | LCSD Result | LCSD Qualifier | Unit | D | %Rec | %Rec     |     | RPD | RPD Limit |
|----------------------------------|-------------|-------------|----------------|------|---|------|----------|-----|-----|-----------|
|                                  |             |             |                |      |   |      | Limits   | RPD |     |           |
| Benzo[a]pyrene                   | 1.98        | 2.15        |                | ug/L |   | 109  | 70 - 130 | 4   | 20  |           |
| Benzo[b]fluoranthene             | 1.98        | 2.17        |                | ug/L |   | 110  | 70 - 130 | 4   | 20  |           |
| Benzo[g,h,i]perylene             | 1.98        | 2.09        |                | ug/L |   | 105  | 70 - 130 | 7   | 20  |           |
| Benzo[k]fluoranthene             | 1.98        | 2.27        |                | ug/L |   | 115  | 70 - 130 | 3   | 20  |           |
| beta-BHC                         | 1.98        | 2.05        |                | ug/L |   | 104  | 70 - 130 | 4   | 20  |           |
| Bis(2-ethylhexyl) phthalate      | 1.98        | 1.62        | *1             | ug/L |   | 82   | 70 - 130 | 30  | 20  |           |
| Bromacil                         | 1.98        | 2.27        |                | ug/L |   | 114  | 70 - 130 | 3   | 20  |           |
| Butachlor                        | 1.98        | 2.47        |                | ug/L |   | 125  | 70 - 130 | 3   | 20  |           |
| Butylbenzylphthalate             | 1.98        | 2.07        |                | ug/L |   | 105  | 70 - 130 | 5   | 20  |           |
| Chlorobenzilate                  | 1.98        | 2.44        |                | ug/L |   | 123  | 70 - 130 | 4   | 20  |           |
| Chloroneb                        | 1.98        | 2.02        |                | ug/L |   | 102  | 70 - 130 | 2   | 20  |           |
| Chlorothalonil (Draconil, Bravo) | 1.98        | 1.98        |                | ug/L |   | 100  | 70 - 130 | 2   | 20  |           |
| Chlorpyrifos                     | 1.98        | 2.24        |                | ug/L |   | 113  | 70 - 130 | 1   | 20  |           |
| Chrysene                         | 1.98        | 1.97        |                | ug/L |   | 99   | 70 - 130 | 4   | 20  |           |
| delta-BHC                        | 1.98        | 1.97        |                | ug/L |   | 100  | 70 - 130 | 3   | 20  |           |
| Di(2-ethylhexyl)adipate          | 1.98        | 1.89        |                | ug/L |   | 96   | 70 - 130 | 14  | 20  |           |
| Dibenz(a,h)anthracene            | 1.98        | 2.05        |                | ug/L |   | 104  | 70 - 130 | 9   | 20  |           |
| Diclorvos (DDVP)                 | 1.98        | 2.73        | *+             | ug/L |   | 138  | 70 - 130 | 1   | 20  |           |
| Dieldrin                         | 1.98        | 1.97        |                | ug/L |   | 100  | 70 - 130 | 3   | 20  |           |
| Diethylphthalate                 | 1.98        | 2.26        |                | ug/L |   | 114  | 70 - 130 | 2   | 20  |           |
| Dimethylphthalate                | 1.98        | 2.21        |                | ug/L |   | 111  | 70 - 130 | 2   | 20  |           |
| Di-n-butyl phthalate             | 3.96        | 4.10        |                | ug/L |   | 103  | 70 - 130 | 6   | 20  |           |
| Di-n-octyl phthalate             | 1.98        | 1.49        | *1             | ug/L |   | 75   | 70 - 130 | 35  | 20  |           |
| Endosulfan I (Alpha)             | 1.98        | 1.90        |                | ug/L |   | 96   | 70 - 130 | 0   | 20  |           |
| Endosulfan II (Beta)             | 1.98        | 2.07        |                | ug/L |   | 104  | 70 - 130 | 5   | 20  |           |
| Endosulfan sulfate               | 1.98        | 2.13        |                | ug/L |   | 108  | 70 - 130 | 5   | 20  |           |
| Endrin                           | 1.98        | 2.35        |                | ug/L |   | 119  | 70 - 130 | 1   | 20  |           |
| Endrin aldehyde                  | 1.98        | 1.86        |                | ug/L |   | 94   | 70 - 130 | 15  | 20  |           |
| EPTC                             | 1.98        | 2.27        |                | ug/L |   | 115  | 70 - 130 | 1   | 20  |           |
| Fluoranthene                     | 1.98        | 2.18        |                | ug/L |   | 110  | 70 - 130 | 2   | 20  |           |
| Fluorene                         | 1.98        | 2.17        |                | ug/L |   | 109  | 70 - 130 | 1   | 20  |           |
| gamma-Chlordane                  | 1.98        | 2.09        |                | ug/L |   | 106  | 70 - 130 | 1   | 20  |           |
| Heptachlor                       | 1.98        | 2.21        |                | ug/L |   | 112  | 70 - 130 | 3   | 20  |           |
| Heptachlor epoxide (isomer B)    | 1.98        | 2.17        |                | ug/L |   | 109  | 70 - 130 | 0   | 20  |           |
| Hexachlorobenzene                | 1.98        | 1.99        |                | ug/L |   | 101  | 70 - 130 | 4   | 20  |           |
| Hexachlorocyclopentadiene        | 1.98        | 2.17        |                | ug/L |   | 110  | 70 - 130 | 13  | 20  |           |
| Indeno[1,2,3-cd]pyrene           | 1.98        | 2.03        |                | ug/L |   | 103  | 70 - 130 | 7   | 20  |           |
| Isophorone                       | 1.98        | 2.16        |                | ug/L |   | 109  | 70 - 130 | 0   | 20  |           |
| Lindane                          | 1.98        | 2.11        |                | ug/L |   | 107  | 70 - 130 | 0   | 20  |           |
| Malathion                        | 1.98        | 2.32        |                | ug/L |   | 117  | 70 - 130 | 1   | 20  |           |
| Methoxychlor                     | 1.98        | 2.13        |                | ug/L |   | 108  | 70 - 130 | 6   | 20  |           |
| Metolachlor                      | 1.98        | 2.43        |                | ug/L |   | 123  | 70 - 130 | 1   | 20  |           |
| Molinate                         | 1.98        | 2.34        |                | ug/L |   | 118  | 70 - 130 | 3   | 20  |           |
| Naphthalene                      | 1.98        | 1.91        |                | ug/L |   | 96   | 70 - 130 | 1   | 20  |           |
| Parathion                        | 1.98        | 2.54        |                | ug/L |   | 128  | 70 - 130 | 2   | 20  |           |
| Pendimethalin (Penoxaline)       | 1.98        | 2.17        |                | ug/L |   | 109  | 70 - 130 | 2   | 20  |           |
| Phenanthrene                     | 1.98        | 1.91        |                | ug/L |   | 97   | 70 - 130 | 0   | 20  |           |
| Propachlor                       | 1.98        | 2.32        |                | ug/L |   | 117  | 70 - 130 | 2   | 20  |           |
| Pyrene                           | 1.98        | 2.14        |                | ug/L |   | 108  | 70 - 130 | 3   | 20  |           |

Eurofins Eaton Analytical Pomona

# QC Sample Results

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

Job ID: 380-57979-1

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

**Lab Sample ID: LCSD 380-51109/4-A**  
**Matrix: Water**  
**Analysis Batch: 51305**

**Client Sample ID: Lab Control Sample Dup**  
**Prep Type: Total/NA**  
**Prep Batch: 51109**

| Analyte         | Spike Added | LCSD Result | LCSD Qualifier | Unit | D | %Rec | %Rec Limits | RPD | RPD Limit |
|-----------------|-------------|-------------|----------------|------|---|------|-------------|-----|-----------|
| Simazine        | 1.98        | 2.25        |                | ug/L |   | 114  | 70 - 130    | 1   | 20        |
| Terbacil        | 1.98        | 2.25        |                | ug/L |   | 114  | 70 - 130    | 2   | 20        |
| Terbuthylazine  | 1.98        | 2.27        |                | ug/L |   | 115  | 70 - 130    | 2   | 20        |
| Thiobencarb     | 1.98        | 2.22        |                | ug/L |   | 112  | 70 - 130    | 0   | 20        |
| trans-Nonachlor | 1.98        | 1.90        |                | ug/L |   | 96   | 70 - 130    | 2   | 20        |
| Trifluralin     | 1.98        | 2.13        |                | ug/L |   | 107  | 70 - 130    | 2   | 20        |

| Surrogate          | LCSD %Recovery | LCSD Qualifier | LCSD Limits |
|--------------------|----------------|----------------|-------------|
| 2-Nitro-m-xylene   | 103            |                | 70 - 130    |
| Perylene-d12       | 95             |                | 70 - 130    |
| Triphenylphosphate | 104            |                | 70 - 130    |

**Lab Sample ID: MRL 380-51109/2-A**  
**Matrix: Water**  
**Analysis Batch: 51305**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 51109**

| Analyte                          | Spike Added | MRL Result | MRL Qualifier | Unit | D | %Rec | %Rec Limits |
|----------------------------------|-------------|------------|---------------|------|---|------|-------------|
| 1-Methylnaphthalene              | 0.0993      | 0.119      |               | ug/L |   | 120  | 50 - 150    |
| 2,4'-DDD                         | 0.0993      | 0.118      |               | ug/L |   | 119  | 50 - 150    |
| 2,4'-DDE                         | 0.0993      | 0.103      |               | ug/L |   | 103  | 50 - 150    |
| 2,4'-DDT                         | 0.0993      | 0.115      |               | ug/L |   | 116  | 50 - 150    |
| 2,4-Dinitrotoluene               | 0.0993      | 0.119      |               | ug/L |   | 119  | 50 - 150    |
| 2,6-Dinitrotoluene               | 0.0993      | 0.116      |               | ug/L |   | 116  | 50 - 150    |
| 2-Methylnaphthalene              | 0.0993      | 0.109      |               | ug/L |   | 110  | 50 - 150    |
| 4,4'-DDD                         | 0.0993      | 0.111      |               | ug/L |   | 112  | 50 - 150    |
| 4,4'-DDE                         | 0.0993      | 0.0916     | J             | ug/L |   | 92   | 50 - 150    |
| 4,4'-DDT                         | 0.0993      | 0.127      |               | ug/L |   | 128  | 50 - 150    |
| Acenaphthene                     | 0.0993      | 0.100      |               | ug/L |   | 101  | 50 - 150    |
| Acenaphthylene                   | 0.0993      | 0.0947     | J             | ug/L |   | 95   | 50 - 150    |
| Acetochlor                       | 0.0496      | 0.0503     | J             | ug/L |   | 101  | 50 - 150    |
| Alachlor                         | 0.0496      | 0.0532     |               | ug/L |   | 107  | 50 - 150    |
| alpha-BHC                        | 0.0993      | 0.102      |               | ug/L |   | 103  | 50 - 150    |
| alpha-Chlordane                  | 0.0248      | <0.029     |               | ug/L |   | 94   | 50 - 150    |
| Anthracene                       | 0.0199      | 0.0196     | J             | ug/L |   | 99   | 50 - 150    |
| Atrazine                         | 0.0496      | 0.0683     |               | ug/L |   | 138  | 50 - 150    |
| Benz(a)anthracene                | 0.0496      | 0.0501     |               | ug/L |   | 101  | 50 - 150    |
| Benzo[a]pyrene                   | 0.0199      | 0.0166     | J             | ug/L |   | 84   | 50 - 150    |
| Benzo[b]fluoranthene             | 0.0199      | 0.0200     |               | ug/L |   | 101  | 50 - 150    |
| Benzo[g,h,i]perylene             | 0.0496      | 0.0512     |               | ug/L |   | 103  | 50 - 150    |
| Benzo[k]fluoranthene             | 0.0199      | 0.0188     | J             | ug/L |   | 95   | 50 - 150    |
| beta-BHC                         | 0.0993      | 0.0953     | J             | ug/L |   | 96   | 50 - 150    |
| Bis(2-ethylhexyl) phthalate      | 0.596       | 0.799      |               | ug/L |   | 134  | 50 - 150    |
| Bromacil                         | 0.0993      | 0.136      |               | ug/L |   | 137  | 50 - 150    |
| Butachlor                        | 0.0496      | 0.0642     |               | ug/L |   | 129  | 50 - 150    |
| Butylbenzylphthalate             | 0.149       | 0.202      | J             | ug/L |   | 136  | 50 - 150    |
| Chlorobenzilate                  | 0.0993      | 0.152      | ^3+           | ug/L |   | 153  | 50 - 150    |
| Chloroneb                        | 0.0993      | 0.106      |               | ug/L |   | 107  | 50 - 150    |
| Chlorothalonil (Draconil, Bravo) | 0.0993      | 0.121      |               | ug/L |   | 122  | 50 - 150    |

# QC Sample Results

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

Job ID: 380-57979-1

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

**Lab Sample ID: MRL 380-51109/2-A**  
**Matrix: Water**  
**Analysis Batch: 51305**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 51109**

| Analyte                       | Spike<br>Added | MRL<br>Result | MRL<br>Qualifier | Unit | D | %Rec | %Rec<br>Limits |
|-------------------------------|----------------|---------------|------------------|------|---|------|----------------|
| Chlorpyrifos                  | 0.0496         | 0.0505        |                  | ug/L |   | 102  | 50 - 150       |
| Chrysene                      | 0.0199         | 0.0205        |                  | ug/L |   | 103  | 50 - 150       |
| delta-BHC                     | 0.0993         | 0.110         |                  | ug/L |   | 111  | 50 - 150       |
| Di(2-ethylhexyl)adipate       | 0.298          | 0.395         | J                | ug/L |   | 133  | 50 - 150       |
| Dibenz(a,h)anthracene         | 0.0496         | 0.0513        |                  | ug/L |   | 103  | 50 - 150       |
| Diclorvos (DDVP)              | 0.0496         | 0.0921        | ^3+              | ug/L |   | 186  | 50 - 150       |
| Dieldrin                      | 0.0993         | 0.100         | J                | ug/L |   | 101  | 50 - 150       |
| Diethylphthalate              | 0.149          | 0.215         | J                | ug/L |   | 144  | 50 - 150       |
| Dimethylphthalate             | 0.298          | 0.337         | J                | ug/L |   | 113  | 50 - 150       |
| Di-n-butyl phthalate          | 0.298          | 0.369         | J                | ug/L |   | 124  | 49 - 243       |
| Di-n-octyl phthalate          | 0.0993         | 0.120         |                  | ug/L |   | 121  | 50 - 150       |
| Endosulfan I (Alpha)          | 0.0993         | 0.0815        | J                | ug/L |   | 82   | 50 - 150       |
| Endosulfan II (Beta)          | 0.0993         | 0.122         |                  | ug/L |   | 123  | 50 - 150       |
| Endosulfan sulfate            | 0.0993         | 0.112         |                  | ug/L |   | 113  | 50 - 150       |
| Endrin                        | 0.0993         | 0.118         |                  | ug/L |   | 118  | 50 - 150       |
| Endrin aldehyde               | 0.0993         | 0.136         |                  | ug/L |   | 137  | 50 - 150       |
| EPTC                          | 0.0993         | 0.126         |                  | ug/L |   | 127  | 50 - 150       |
| Fluoranthene                  | 0.0496         | 0.0519        | J                | ug/L |   | 105  | 50 - 150       |
| Fluorene                      | 0.0496         | 0.0600        |                  | ug/L |   | 121  | 50 - 150       |
| gamma-Chlordane               | 0.0248         | 0.0235        | J                | ug/L |   | 95   | 50 - 150       |
| Heptachlor                    | 0.0397         | 0.0455        |                  | ug/L |   | 115  | 50 - 150       |
| Heptachlor epoxide (isomer B) | 0.0496         | 0.0476        | J                | ug/L |   | 96   | 50 - 150       |
| Hexachlorobenzene             | 0.0496         | 0.0434        | J                | ug/L |   | 87   | 50 - 150       |
| Hexachlorocyclopentadiene     | 0.0496         | 0.0440        | J                | ug/L |   | 89   | 50 - 150       |
| Indeno[1,2,3-cd]pyrene        | 0.0496         | 0.0524        |                  | ug/L |   | 106  | 50 - 150       |
| Isophorone                    | 0.0993         | 0.113         | J                | ug/L |   | 114  | 50 - 150       |
| Lindane                       | 0.0397         | 0.0414        |                  | ug/L |   | 104  | 50 - 150       |
| Malathion                     | 0.0993         | 0.123         |                  | ug/L |   | 123  | 50 - 150       |
| Methoxychlor                  | 0.0993         | 0.135         |                  | ug/L |   | 136  | 50 - 150       |
| Metolachlor                   | 0.0496         | 0.0611        |                  | ug/L |   | 123  | 50 - 150       |
| Molinate                      | 0.0993         | 0.113         |                  | ug/L |   | 113  | 50 - 150       |
| Naphthalene                   | 0.0993         | 0.112         | J                | ug/L |   | 113  | 50 - 150       |
| Parathion                     | 0.0993         | 0.122         |                  | ug/L |   | 123  | 50 - 150       |
| Pendimethalin (Penoxaline)    | 0.0993         | 0.113         |                  | ug/L |   | 114  | 50 - 150       |
| Phenanthrene                  | 0.0199         | 0.0239        | J                | ug/L |   | 121  | 50 - 150       |
| Propachlor                    | 0.0496         | 0.0580        |                  | ug/L |   | 117  | 50 - 150       |
| Pyrene                        | 0.0496         | 0.0501        |                  | ug/L |   | 101  | 50 - 150       |
| Simazine                      | 0.0496         | 0.0569        |                  | ug/L |   | 115  | 50 - 150       |
| Terbacil                      | 0.0993         | 0.124         |                  | ug/L |   | 125  | 50 - 150       |
| Terbutylazine                 | 0.0993         | 0.106         |                  | ug/L |   | 106  | 50 - 150       |
| Thiobencarb                   | 0.0993         | 0.121         | J                | ug/L |   | 122  | 50 - 150       |
| trans-Nonachlor               | 0.0248         | <0.026        |                  | ug/L |   | 95   | 50 - 150       |
| Trifluralin                   | 0.0993         | 0.108         |                  | ug/L |   | 109  | 50 - 150       |

| Surrogate          | MRL<br>%Recovery | MRL<br>Qualifier | Limits   |
|--------------------|------------------|------------------|----------|
| 2-Nitro-m-xylene   | 106              |                  | 70 - 130 |
| Perylene-d12       | 97               |                  | 70 - 130 |
| Triphenylphosphate | 108              |                  | 70 - 130 |



# QC Sample Results

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

Job ID: 380-57979-1

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

**Lab Sample ID: 380-57897-P-1-A MS**  
**Matrix: Water**  
**Analysis Batch: 51305**

**Client Sample ID: Matrix Spike**  
**Prep Type: Total/NA**  
**Prep Batch: 51109**

| Analyte                          | Sample Result | Sample Qualifier | Spike Added | MS Result | MS Qualifier | Unit | D | %Rec | %Rec Limits |
|----------------------------------|---------------|------------------|-------------|-----------|--------------|------|---|------|-------------|
| 1-Methylnaphthalene              | <0.098        |                  | 1.96        | 1.95      |              | ug/L |   | 99   | 70 - 130    |
| 2,4'-DDD                         | <0.098        |                  | 1.96        | 2.09      |              | ug/L |   | 107  | 70 - 130    |
| 2,4'-DDE                         | <0.098        |                  | 1.96        | 1.99      |              | ug/L |   | 102  | 70 - 130    |
| 2,4'-DDT                         | <0.098        |                  | 1.96        | 2.19      |              | ug/L |   | 112  | 70 - 130    |
| 2,4-Dinitrotoluene               | <0.098        |                  | 1.96        | 2.33      |              | ug/L |   | 119  | 70 - 130    |
| 2,6-Dinitrotoluene               | <0.098        |                  | 1.96        | 2.24      |              | ug/L |   | 114  | 70 - 130    |
| 2-Methylnaphthalene              | <0.098        |                  | 1.96        | 1.97      |              | ug/L |   | 101  | 70 - 130    |
| 4,4'-DDD                         | <0.098        |                  | 1.96        | 2.11      |              | ug/L |   | 108  | 70 - 130    |
| 4,4'-DDE                         | <0.098        |                  | 1.96        | 1.90      |              | ug/L |   | 97   | 70 - 130    |
| 4,4'-DDT                         | <0.098        |                  | 1.96        | 2.20      |              | ug/L |   | 113  | 70 - 130    |
| Acenaphthene                     | <0.098        |                  | 1.96        | 1.83      |              | ug/L |   | 93   | 70 - 130    |
| Acenaphthylene                   | <0.098        |                  | 1.96        | 2.03      |              | ug/L |   | 104  | 70 - 130    |
| Acetochlor                       | <0.098        | *+               | 1.96        | 2.27      |              | ug/L |   | 116  | 70 - 130    |
| Alachlor                         | <0.049        |                  | 1.96        | 2.07      |              | ug/L |   | 106  | 70 - 130    |
| alpha-BHC                        | <0.098        |                  | 1.96        | 1.92      |              | ug/L |   | 98   | 70 - 130    |
| alpha-Chlordane                  | <0.049        |                  | 1.96        | 1.99      |              | ug/L |   | 102  | 70 - 130    |
| Anthracene                       | <0.020        | F1               | 1.96        | 1.17      | F1           | ug/L |   | 60   | 70 - 130    |
| Atrazine                         | <0.049        |                  | 1.96        | 2.18      |              | ug/L |   | 112  | 70 - 130    |
| Benz(a)anthracene                | <0.049        |                  | 1.96        | 2.03      |              | ug/L |   | 104  | 70 - 130    |
| Benzo[a]pyrene                   | <0.020        |                  | 1.96        | 1.71      |              | ug/L |   | 88   | 70 - 130    |
| Benzo[b]fluoranthene             | <0.020        |                  | 1.96        | 2.12      |              | ug/L |   | 109  | 70 - 130    |
| Benzo[g,h,i]perylene             | <0.049        |                  | 1.96        | 1.93      |              | ug/L |   | 99   | 70 - 130    |
| Benzo[k]fluoranthene             | <0.020        |                  | 1.96        | 2.19      |              | ug/L |   | 112  | 70 - 130    |
| beta-BHC                         | <0.098        |                  | 1.96        | 1.93      |              | ug/L |   | 99   | 70 - 130    |
| Bis(2-ethylhexyl) phthalate      | <0.59         | *1               | 1.96        | 1.58      |              | ug/L |   | 81   | 70 - 130    |
| Bromacil                         | <0.098        |                  | 1.96        | 2.34      |              | ug/L |   | 119  | 70 - 130    |
| Butachlor                        | <0.049        |                  | 1.96        | 2.21      |              | ug/L |   | 113  | 70 - 130    |
| Butylbenzylphthalate             | <0.49         |                  | 1.96        | 1.99      |              | ug/L |   | 102  | 70 - 130    |
| Chlorobenzilate                  | <0.098        | ^3+              | 1.96        | 2.23      |              | ug/L |   | 114  | 70 - 130    |
| Chloroneb                        | <0.098        |                  | 1.96        | 1.94      |              | ug/L |   | 99   | 70 - 130    |
| Chlorothalonil (Draconil, Bravo) | <0.098        |                  | 1.96        | 2.13      |              | ug/L |   | 109  | 70 - 130    |
| Chlorpyrifos                     | <0.049        |                  | 1.96        | 2.19      |              | ug/L |   | 112  | 70 - 130    |
| Chrysene                         | <0.020        |                  | 1.96        | 1.93      |              | ug/L |   | 99   | 70 - 130    |
| delta-BHC                        | <0.098        |                  | 1.96        | 1.79      |              | ug/L |   | 92   | 70 - 130    |
| Di(2-ethylhexyl)adipate          | <0.59         |                  | 1.96        | 1.94      |              | ug/L |   | 95   | 70 - 130    |
| Dibenz(a,h)anthracene            | <0.049        |                  | 1.96        | 1.98      |              | ug/L |   | 101  | 70 - 130    |
| Diclorvos (DDVP)                 | <0.049        | *+ ^3+ F1        | 1.96        | 2.77      | F1           | ug/L |   | 142  | 70 - 130    |
| Dieldrin                         | <0.20         |                  | 1.96        | 2.06      |              | ug/L |   | 105  | 70 - 130    |
| Diethylphthalate                 | <0.49         |                  | 1.96        | 2.13      |              | ug/L |   | 109  | 70 - 130    |
| Dimethylphthalate                | <0.49         |                  | 1.96        | 2.14      |              | ug/L |   | 109  | 70 - 130    |
| Di-n-butyl phthalate             | <0.98         |                  | 3.91        | 4.01      |              | ug/L |   | 102  | 70 - 130    |
| Di-n-octyl phthalate             | <0.098        | *1               | 1.96        | 1.49      |              | ug/L |   | 76   | 70 - 130    |
| Endosulfan I (Alpha)             | <0.098        |                  | 1.96        | 1.69      |              | ug/L |   | 86   | 70 - 130    |
| Endosulfan II (Beta)             | <0.098        |                  | 1.96        | 2.01      |              | ug/L |   | 103  | 70 - 130    |
| Endosulfan sulfate               | <0.098        |                  | 1.96        | 2.23      |              | ug/L |   | 114  | 70 - 130    |
| Endrin                           | <0.098        |                  | 1.96        | 1.94      |              | ug/L |   | 99   | 70 - 130    |
| Endrin aldehyde                  | <0.098        |                  | 1.96        | 1.40      |              | ug/L |   | 71   | 70 - 130    |
| EPTC                             | <0.098        |                  | 1.96        | 2.20      |              | ug/L |   | 112  | 70 - 130    |

# QC Sample Results

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

Job ID: 380-57979-1

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

**Lab Sample ID: 380-57897-P-1-A MS**  
**Matrix: Water**  
**Analysis Batch: 51305**

**Client Sample ID: Matrix Spike**  
**Prep Type: Total/NA**  
**Prep Batch: 51109**

| Analyte                       | Sample | Sample Qualifier | Spike Added | MS     | MS Qualifier | Unit | D | %Rec | %Rec Limits |
|-------------------------------|--------|------------------|-------------|--------|--------------|------|---|------|-------------|
|                               | Result |                  |             | Result |              |      |   |      |             |
| Fluoranthene                  | <0.098 |                  | 1.96        | 2.14   |              | ug/L |   | 109  | 70 - 130    |
| Fluorene                      | <0.049 |                  | 1.96        | 2.10   |              | ug/L |   | 107  | 70 - 130    |
| gamma-Chlordane               | <0.049 |                  | 1.96        | 1.98   |              | ug/L |   | 101  | 70 - 130    |
| Heptachlor                    | <0.039 |                  | 1.96        | 2.11   |              | ug/L |   | 108  | 70 - 130    |
| Heptachlor epoxide (isomer B) | <0.049 |                  | 1.96        | 2.06   |              | ug/L |   | 105  | 70 - 130    |
| Hexachlorobenzene             | <0.049 |                  | 1.96        | 2.04   |              | ug/L |   | 104  | 70 - 130    |
| Hexachlorocyclopentadiene     | <0.049 |                  | 1.96        | 2.34   |              | ug/L |   | 120  | 70 - 130    |
| Indeno[1,2,3-cd]pyrene        | <0.049 |                  | 1.96        | 1.99   |              | ug/L |   | 102  | 70 - 130    |
| Isophorone                    | <0.49  |                  | 1.96        | 2.14   |              | ug/L |   | 109  | 70 - 130    |
| Lindane                       | <0.039 |                  | 1.96        | 1.95   |              | ug/L |   | 100  | 70 - 130    |
| Malathion                     | <0.098 |                  | 1.96        | 2.18   |              | ug/L |   | 112  | 70 - 130    |
| Methoxychlor                  | <0.098 |                  | 1.96        | 2.20   |              | ug/L |   | 112  | 70 - 130    |
| Metolachlor                   | <0.049 |                  | 1.96        | 2.24   |              | ug/L |   | 115  | 70 - 130    |
| Molinate                      | <0.098 |                  | 1.96        | 2.26   |              | ug/L |   | 116  | 70 - 130    |
| Naphthalene                   | <0.29  |                  | 1.96        | 1.87   |              | ug/L |   | 96   | 70 - 130    |
| Parathion                     | <0.098 | *+               | 1.96        | 2.53   |              | ug/L |   | 129  | 70 - 130    |
| Pendimethalin (Penoxaline)    | <0.098 |                  | 1.96        | 2.43   |              | ug/L |   | 124  | 70 - 130    |
| Phenanthrene                  | <0.039 |                  | 1.96        | 1.84   |              | ug/L |   | 94   | 70 - 130    |
| Propachlor                    | <0.049 |                  | 1.96        | 2.27   |              | ug/L |   | 116  | 70 - 130    |
| Pyrene                        | <0.049 |                  | 1.96        | 2.16   |              | ug/L |   | 111  | 70 - 130    |
| Simazine                      | <0.049 |                  | 1.96        | 2.21   |              | ug/L |   | 113  | 70 - 130    |
| Terbacil                      | <0.098 |                  | 1.96        | 2.15   |              | ug/L |   | 110  | 70 - 130    |
| Terbutylazine                 | <0.098 |                  | 1.96        | 2.24   |              | ug/L |   | 115  | 70 - 130    |
| Thiobencarb                   | <0.20  |                  | 1.96        | 2.11   |              | ug/L |   | 108  | 70 - 130    |
| trans-Nonachlor               | <0.049 |                  | 1.96        | 2.10   |              | ug/L |   | 107  | 70 - 130    |
| Trifluralin                   | <0.098 |                  | 1.96        | 2.32   |              | ug/L |   | 119  | 70 - 130    |

| Surrogate          | MS MS     |           | Limits   |
|--------------------|-----------|-----------|----------|
|                    | %Recovery | Qualifier |          |
| 2-Nitro-m-xylene   | 98        |           | 70 - 130 |
| Perylene-d12       | 92        |           | 70 - 130 |
| Triphenylphosphate | 115       |           | 70 - 130 |

**Lab Sample ID: 380-57909-AR-1-A DU**  
**Matrix: Water**  
**Analysis Batch: 51305**

**Client Sample ID: Duplicate**  
**Prep Type: Total/NA**  
**Prep Batch: 51109**

| Analyte             | Sample Result | Sample Qualifier | DU Result | DU Qualifier | Unit | D | RPD |       |
|---------------------|---------------|------------------|-----------|--------------|------|---|-----|-------|
|                     |               |                  |           |              |      |   | RPD | Limit |
| 1-Methylnaphthalene | <0.099        |                  | <0.098    |              | ug/L |   | NC  | 20    |
| 2,4'-DDD            | <0.099        |                  | <0.098    |              | ug/L |   | NC  | 20    |
| 2,4'-DDE            | <0.099        |                  | <0.098    |              | ug/L |   | NC  | 20    |
| 2,4'-DDT            | <0.099        |                  | <0.098    |              | ug/L |   | NC  | 20    |
| 2,4-Dinitrotoluene  | <0.099        |                  | <0.098    |              | ug/L |   | NC  | 20    |
| 2,6-Dinitrotoluene  | <0.099        |                  | <0.098    |              | ug/L |   | NC  | 20    |
| 2-Methylnaphthalene | <0.099        |                  | <0.098    |              | ug/L |   | NC  | 20    |
| 4,4'-DDD            | <0.099        |                  | <0.098    |              | ug/L |   | NC  | 20    |
| 4,4'-DDE            | <0.099        |                  | <0.098    |              | ug/L |   | NC  | 20    |
| 4,4'-DDT            | <0.099        |                  | <0.098    |              | ug/L |   | NC  | 20    |
| Acenaphthene        | <0.099        |                  | <0.098    |              | ug/L |   | NC  | 20    |

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# QC Sample Results

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

Job ID: 380-57979-1

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

**Lab Sample ID: 380-57909-AR-1-A DU**  
**Matrix: Water**  
**Analysis Batch: 51305**

**Client Sample ID: Duplicate**  
**Prep Type: Total/NA**  
**Prep Batch: 51109**

| Analyte                          | Sample | Sample    | DU     | DU        | Unit | D | RPD | Limit |
|----------------------------------|--------|-----------|--------|-----------|------|---|-----|-------|
|                                  | Result | Qualifier | Result | Qualifier |      |   |     |       |
| Acenaphthylene                   | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| Acetochlor                       | <0.099 | *+        | <0.098 | *+        | ug/L |   | NC  | 20    |
| Alachlor                         | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| alpha-BHC                        | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| alpha-Chlordane                  | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| Anthracene                       | <0.020 |           | <0.020 |           | ug/L |   | NC  | 20    |
| Atrazine                         | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| Benz(a)anthracene                | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| Benzo[a]pyrene                   | <0.020 |           | <0.020 |           | ug/L |   | NC  | 20    |
| Benzo[b]fluoranthene             | <0.020 |           | <0.020 |           | ug/L |   | NC  | 20    |
| Benzo[g,h,i]perylene             | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| Benzo[k]fluoranthene             | <0.020 |           | <0.020 |           | ug/L |   | NC  | 20    |
| beta-BHC                         | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| Bis(2-ethylhexyl) phthalate      | <0.59  | *1        | <0.59  | *1        | ug/L |   | NC  | 20    |
| Bromacil                         | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| Butachlor                        | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| Butylbenzylphthalate             | <0.49  |           | <0.49  |           | ug/L |   | NC  | 20    |
| Chlorobenzilate                  | <0.099 | ^3+       | <0.098 |           | ug/L |   | NC  | 20    |
| Chloroneb                        | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| Chlorothalonil (Draconil, Bravo) | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| Chlorpyrifos                     | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| Chrysene                         | <0.020 |           | <0.020 |           | ug/L |   | NC  | 20    |
| delta-BHC                        | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| Di(2-ethylhexyl)adipate          | <0.59  |           | <0.59  |           | ug/L |   | NC  | 20    |
| Dibenz(a,h)anthracene            | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| Diclorvos (DDVP)                 | <0.049 | *+ ^3+    | <0.049 | *+        | ug/L |   | NC  | 20    |
| Dieldrin                         | <0.20  |           | <0.20  |           | ug/L |   | NC  | 20    |
| Diethylphthalate                 | <0.49  |           | <0.49  |           | ug/L |   | NC  | 20    |
| Dimethylphthalate                | <0.49  |           | <0.49  |           | ug/L |   | NC  | 20    |
| Di-n-butyl phthalate             | <0.99  |           | <0.98  |           | ug/L |   | NC  | 20    |
| Di-n-octyl phthalate             | <0.099 | *1        | <0.098 | *1        | ug/L |   | NC  | 20    |
| Endosulfan I (Alpha)             | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| Endosulfan II (Beta)             | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| Endosulfan sulfate               | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| Endrin                           | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| Endrin aldehyde                  | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| EPTC                             | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| Fluoranthene                     | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| Fluorene                         | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| gamma-Chlordane                  | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| Heptachlor                       | <0.040 |           | <0.039 |           | ug/L |   | NC  | 20    |
| Heptachlor epoxide (isomer B)    | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| Hexachlorobenzene                | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| Hexachlorocyclopentadiene        | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| Indeno[1,2,3-cd]pyrene           | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| Isophorone                       | <0.49  |           | <0.49  |           | ug/L |   | NC  | 20    |
| Lindane                          | <0.040 |           | <0.039 |           | ug/L |   | NC  | 20    |
| Malathion                        | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| Methoxychlor                     | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |

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# QC Sample Results

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

Job ID: 380-57979-1

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

**Lab Sample ID: 380-57909-AR-1-A DU**  
**Matrix: Water**  
**Analysis Batch: 51305**

**Client Sample ID: Duplicate**  
**Prep Type: Total/NA**  
**Prep Batch: 51109**

| Analyte                          | Sample | Sample    | DU     | DU        | Unit | D | RPD | Limit |
|----------------------------------|--------|-----------|--------|-----------|------|---|-----|-------|
|                                  | Result | Qualifier | Result | Qualifier |      |   |     |       |
| Metolachlor                      | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| Molinate                         | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| Naphthalene                      | <0.30  |           | <0.29  |           | ug/L |   | NC  | 20    |
| Parathion                        | <0.099 | *+        | <0.098 | *+        | ug/L |   | NC  | 20    |
| Pendimethalin (Penoxaline)       | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| Phenanthrene                     | <0.040 |           | <0.039 |           | ug/L |   | NC  | 20    |
| Propachlor                       | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| Pyrene                           | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| Simazine                         | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| Terbacil                         | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| Terbutylazine                    | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |
| Thiobencarb                      | <0.20  |           | <0.20  |           | ug/L |   | NC  | 20    |
| Total Permethrin (mixed isomers) | <0.20  |           | <0.20  |           | ug/L |   | NC  | 20    |
| trans-Nonachlor                  | <0.049 |           | <0.049 |           | ug/L |   | NC  | 20    |
| Trifluralin                      | <0.099 |           | <0.098 |           | ug/L |   | NC  | 20    |

| Surrogate          | DU        | DU        | Limits   |
|--------------------|-----------|-----------|----------|
|                    | %Recovery | Qualifier |          |
| 2-Nitro-m-xylene   | 97        |           | 70 - 130 |
| Perylene-d12       | 92        |           | 70 - 130 |
| Triphenylphosphate | 110       |           | 70 - 130 |

## Method: 533 - Perfluorinated and Polyfluorinated Alkyl Substances in Drinking Water

**Lab Sample ID: MBL 380-53035/21-A**  
**Matrix: Water**  
**Analysis Batch: 53385**

**Client Sample ID: Method Blank**  
**Prep Type: Total/NA**  
**Prep Batch: 53035**

| Analyte   | MBL    | MBL       | RL  | Unit | D | Prepared       | Analyzed       | Dil Fac |
|---|--------|-----------|-----|------|---|----------------|----------------|---------|
|   | Result | Qualifier |     |      |   |                |                |         |
| 11-Chloroeicosafuoro-3-oxaundecan<br>e-1-sulfonic acid (11Cl-PF3OUdS) | <0.30  |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 9-Chlorohexadecafluoro-3-oxanonan<br>e-1-sulfonic acid(9Cl-PF3ONS)    | <0.30  |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 4,8-Dioxa-3H-perfluorononanoic acid<br>(ADONA)                        | <0.60  |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Hexafluoropropylene Oxide Dimer<br>Acid (HFPO-DA/GenX)                | <1.0   |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Perfluorobutanesulfonic acid (PFBS)                                   | <0.37  |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Perfluorodecanoic acid (PFDA)   | <0.31  |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Perfluorododecanoic acid (PFDoA)                                      | <0.54  |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Perfluoroheptanoic acid (PFHpA)                                       | <0.39  |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Perfluorohexanesulfonic acid (PFHxS)                                  | <0.32  |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Perfluorohexanoic acid (PFHxA)  | <0.46  |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Perfluorononanoic acid (PFNA)   | <0.40  |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Perfluorooctanesulfonic acid (PFOS)                                   | <0.43  |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Perfluorooctanoic acid (PFOA)   | <0.38  |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Perfluoroundecanoic acid (PFUnA)                                      | <0.42  |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Perfluorobutanoic acid (PFBA)   | <0.69  |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 1H,1H,2H,2H-Perfluorodecane<br>sulfonic acid (8:2 FTS)                | <0.38  |           | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |

Eurofins Eaton Analytical Pomona

# QC Sample Results

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

Job ID: 380-57979-1

## Method: 533 - Perfluorinated and Polyfluorinated Alkyl Substances in Drinking Water (Continued)

**Lab Sample ID: MBL 380-53035/21-A**  
**Matrix: Water**  
**Analysis Batch: 53385**

**Client Sample ID: Method Blank**  
**Prep Type: Total/NA**  
**Prep Batch: 53035**

| Analyte   | MBL Result | MBL Qualifier | RL  | Unit | D | Prepared       | Analyzed       | Dil Fac |
|---|------------|---------------|-----|------|---|----------------|----------------|---------|
| 1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS) | <0.37      |               | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS) | <0.48      |               | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)          | <0.47      |               | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Perfluoro (2-ethoxyethane) sulfonic acid (PFEEESA)  | <0.25      |               | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Perfluoro-3-methoxypropanoic acid (PFMPA)           | <0.46      |               | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Perfluoro-4-methoxybutanoic acid (PFMBA)            | <0.15      |               | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Perfluoropentanoic acid (PFPeA)                     | <0.38      |               | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Perfluoroheptanesulfonic acid (PFHpS)               | <0.36      |               | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| Perfluoropentanesulfonic acid (PFPeS)               | <0.39      |               | 2.0 | ng/L |   | 08/24/23 16:20 | 08/27/23 03:42 | 1       |

| Isotope Dilution | MBL %Recovery | MBL Qualifier | Limits   | Prepared       | Analyzed       | Dil Fac |
|------------------|---------------|---------------|----------|----------------|----------------|---------|
| 13C3 HFPO-DA     | 79            |               | 50 - 200 | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 13C6 PFDA        | 95            |               | 50 - 200 | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 13C5 PFHxA       | 88            |               | 50 - 200 | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 13C4 PFHpA       | 88            |               | 50 - 200 | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 13C8 PFOA        | 93            |               | 50 - 200 | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 13C9 PFNA        | 91            |               | 50 - 200 | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 13C7 PFUnA       | 88            |               | 50 - 200 | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 13C2 PFDoA       | 91            |               | 50 - 200 | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 13C4 PFBA        | 92            |               | 50 - 200 | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 13C5 PFPeA       | 90            |               | 50 - 200 | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 13C3 PFBS        | 90            |               | 50 - 200 | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 13C3 PFHxS       | 90            |               | 50 - 200 | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 13C8 PFOS        | 91            |               | 50 - 200 | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 13C2-4:2-FTS     | 101           |               | 50 - 200 | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 13C2-6:2-FTS     | 98            |               | 50 - 200 | 08/24/23 16:20 | 08/27/23 03:42 | 1       |
| 13C2-8:2-FTS     | 109           |               | 50 - 200 | 08/24/23 16:20 | 08/27/23 03:42 | 1       |

**Lab Sample ID: LCS 380-53035/23-A**  
**Matrix: Water**  
**Analysis Batch: 53385**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 53035**

| Analyte   | Spike Added | LCS Result | LCS Qualifier | Unit | D | %Rec | %Rec Limits |
|---|-------------|------------|---------------|------|---|------|-------------|
| 11-Chloroeicosfluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS) | 60.2        | 50.3       |               | ng/L |   | 84   | 70 - 130    |
| 9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid(9Cl-PF3ONS)    | 60.2        | 52.4       |               | ng/L |   | 87   | 70 - 130    |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA)                       | 60.2        | 56.2       |               | ng/L |   | 93   | 70 - 130    |
| Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)               | 60.2        | 50.5       |               | ng/L |   | 84   | 70 - 130    |

Eurofins Eaton Analytical Pomona

# QC Sample Results

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

Job ID: 380-57979-1

## Method: 533 - Perfluorinated and Polyfluorinated Alkyl Substances in Drinking Water (Continued)

**Lab Sample ID: LCS 380-53035/23-A**  
**Matrix: Water**  
**Analysis Batch: 53385**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 53035**

| Analyte   | Spike Added | LCS Result | LCS Qualifier | Unit | D | %Rec | %Rec Limits |
|---|-------------|------------|---------------|------|---|------|-------------|
| Perfluorobutanesulfonic acid (PFBS)                 | 60.2        | 57.4       |               | ng/L |   | 95   | 70 - 130    |
| Perfluorodecanoic acid (PFDA)                       | 60.2        | 55.1       |               | ng/L |   | 92   | 70 - 130    |
| Perfluorododecanoic acid (PFDoA)                    | 60.2        | 54.4       |               | ng/L |   | 90   | 70 - 130    |
| Perfluoroheptanoic acid (PFHpA)                     | 60.2        | 54.2       |               | ng/L |   | 90   | 70 - 130    |
| Perfluorohexanesulfonic acid (PFHxS)                | 60.2        | 53.4       |               | ng/L |   | 89   | 70 - 130    |
| Perfluorohexanoic acid (PFHxA)                      | 60.2        | 56.1       |               | ng/L |   | 93   | 70 - 130    |
| Perfluorononanoic acid (PFNA)                       | 60.2        | 54.9       |               | ng/L |   | 91   | 70 - 130    |
| Perfluorooctanesulfonic acid (PFOS)                 | 60.2        | 53.3       |               | ng/L |   | 88   | 70 - 130    |
| Perfluorooctanoic acid (PFOA)                       | 60.2        | 54.5       |               | ng/L |   | 91   | 70 - 130    |
| Perfluoroundecanoic acid (PFUnA)                    | 60.2        | 56.7       |               | ng/L |   | 94   | 70 - 130    |
| Perfluorobutanoic acid (PFBA)                       | 60.2        | 57.2       |               | ng/L |   | 95   | 70 - 130    |
| 1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS) | 60.2        | 56.6       |               | ng/L |   | 94   | 70 - 130    |
| 1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS) | 60.2        | 55.0       |               | ng/L |   | 91   | 70 - 130    |
| 1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS) | 60.2        | 54.3       |               | ng/L |   | 90   | 70 - 130    |
| Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)          | 60.2        | 51.3       |               | ng/L |   | 85   | 70 - 130    |
| Perfluoro (2-ethoxyethane) sulfonic acid (PFEEESA)  | 60.2        | 56.1       |               | ng/L |   | 93   | 70 - 130    |
| Perfluoro-3-methoxypropanoic acid (PFMPA)           | 60.2        | 54.9       |               | ng/L |   | 91   | 70 - 130    |
| Perfluoro-4-methoxybutanoic acid (PFMBA)            | 60.2        | 49.9       |               | ng/L |   | 83   | 70 - 130    |
| Perfluoropentanoic acid (PFPeA)                     | 60.2        | 57.9       |               | ng/L |   | 96   | 70 - 130    |
| Perfluoroheptanesulfonic acid (PFHpS)               | 60.2        | 53.7       |               | ng/L |   | 89   | 70 - 130    |
| Perfluoropentanesulfonic acid (PFPeS)               | 60.2        | 53.4       |               | ng/L |   | 89   | 70 - 130    |

| Isotope Dilution | LCS %Recovery | LCS Qualifier | Limits   |
|------------------|---------------|---------------|----------|
| 13C3 HFPO-DA     | 91            |               | 50 - 200 |
| 13C6 PFDA        | 95            |               | 50 - 200 |
| 13C5 PFHxA       | 94            |               | 50 - 200 |
| 13C4 PFHpA       | 96            |               | 50 - 200 |
| 13C8 PFOA        | 97            |               | 50 - 200 |
| 13C9 PFNA        | 99            |               | 50 - 200 |
| 13C7 PFUnA       | 96            |               | 50 - 200 |
| 13C2 PFDoA       | 98            |               | 50 - 200 |
| 13C4 PFBA        | 93            |               | 50 - 200 |
| 13C5 PFPeA       | 94            |               | 50 - 200 |
| 13C3 PFBS        | 92            |               | 50 - 200 |
| 13C3 PFHxS       | 96            |               | 50 - 200 |
| 13C8 PFOS        | 95            |               | 50 - 200 |
| 13C2-4:2-FTS     | 95            |               | 50 - 200 |
| 13C2-6:2-FTS     | 93            |               | 50 - 200 |

# QC Sample Results

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

Job ID: 380-57979-1

## Method: 533 - Perfluorinated and Polyfluorinated Alkyl Substances in Drinking Water (Continued)

**Lab Sample ID: LCS 380-53035/23-A**  
**Matrix: Water**  
**Analysis Batch: 53385**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 53035**

| Isotope Dilution | LCS LCS   |           | Limits   |
|------------------|-----------|-----------|----------|
|                  | %Recovery | Qualifier |          |
| 13C2-8:2-FTS     | 90        |           | 50 - 200 |

**Lab Sample ID: LCSD 380-53035/24-A**  
**Matrix: Water**  
**Analysis Batch: 53385**

**Client Sample ID: Lab Control Sample Dup**  
**Prep Type: Total/NA**  
**Prep Batch: 53035**

| Analyte   | Spike Added | LCSD Result | LCSD Qualifier | Unit | D | %Rec | %Rec     |     | RPD | Limit |
|---|-------------|-------------|----------------|------|---|------|----------|-----|-----|-------|
|   |             |             |                |      |   |      | Limits   | RPD |     |       |
| 11-Chloroeicosfluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS) | 60.2        | 55.1        |                | ng/L |   | 92   | 70 - 130 | 9   | 30  |       |
| 9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid(9Cl-PF3ONS)    | 60.2        | 56.2        |                | ng/L |   | 93   | 70 - 130 | 7   | 30  |       |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA)                       | 60.2        | 55.4        |                | ng/L |   | 92   | 70 - 130 | 1   | 30  |       |
| Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)               | 60.2        | 50.6        |                | ng/L |   | 84   | 70 - 130 | 0   | 30  |       |
| Perfluorobutanesulfonic acid (PFBS)                               | 60.2        | 57.4        |                | ng/L |   | 95   | 70 - 130 | 0   | 30  |       |
| Perfluorodecanoic acid (PFDA)                                     | 60.2        | 56.5        |                | ng/L |   | 94   | 70 - 130 | 2   | 30  |       |
| Perfluorododecanoic acid (PFDoA)                                  | 60.2        | 55.9        |                | ng/L |   | 93   | 70 - 130 | 3   | 30  |       |
| Perfluoroheptanoic acid (PFHpA)                                   | 60.2        | 56.4        |                | ng/L |   | 94   | 70 - 130 | 4   | 30  |       |
| Perfluorohexanesulfonic acid (PFHxS)                              | 60.2        | 55.0        |                | ng/L |   | 91   | 70 - 130 | 3   | 30  |       |
| Perfluorohexanoic acid (PFHxA)                                    | 60.2        | 64.0        |                | ng/L |   | 106  | 70 - 130 | 13  | 30  |       |
| Perfluorononanoic acid (PFNA)                                     | 60.2        | 59.0        |                | ng/L |   | 98   | 70 - 130 | 7   | 30  |       |
| Perfluorooctanesulfonic acid (PFOS)                               | 60.2        | 57.2        |                | ng/L |   | 95   | 70 - 130 | 7   | 30  |       |
| Perfluorooctanoic acid (PFOA)                                     | 60.2        | 57.6        |                | ng/L |   | 96   | 70 - 130 | 5   | 30  |       |
| Perfluoroundecanoic acid (PFUnA)                                  | 60.2        | 59.2        |                | ng/L |   | 98   | 70 - 130 | 4   | 30  |       |
| Perfluorobutanoic acid (PFBA)                                     | 60.2        | 58.2        |                | ng/L |   | 97   | 70 - 130 | 2   | 30  |       |
| 1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)               | 60.2        | 61.1        |                | ng/L |   | 101  | 70 - 130 | 8   | 30  |       |
| 1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)               | 60.2        | 55.9        |                | ng/L |   | 93   | 70 - 130 | 2   | 30  |       |
| 1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)               | 60.2        | 60.3        |                | ng/L |   | 100  | 70 - 130 | 10  | 30  |       |
| Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)                        | 60.2        | 52.3        |                | ng/L |   | 87   | 70 - 130 | 2   | 30  |       |
| Perfluoro (2-ethoxyethane) sulfonic acid (PFEEESA)                | 60.2        | 60.1        |                | ng/L |   | 100  | 70 - 130 | 7   | 30  |       |
| Perfluoro-3-methoxypropanoic acid (PFMPA)                         | 60.2        | 57.1        |                | ng/L |   | 95   | 70 - 130 | 4   | 30  |       |
| Perfluoro-4-methoxybutanoic acid (PFMBA)                          | 60.2        | 58.4        |                | ng/L |   | 97   | 70 - 130 | 16  | 30  |       |
| Perfluoropentanoic acid (PFPeA)                                   | 60.2        | 61.6        |                | ng/L |   | 102  | 70 - 130 | 6   | 30  |       |
| Perfluoroheptanesulfonic acid (PFHpS)                             | 60.2        | 58.3        |                | ng/L |   | 97   | 70 - 130 | 8   | 30  |       |
| Perfluoropentanesulfonic acid (PFPeS)                             | 60.2        | 56.2        |                | ng/L |   | 93   | 70 - 130 | 5   | 30  |       |

# QC Sample Results

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

Job ID: 380-57979-1

## Method: 533 - Perfluorinated and Polyfluorinated Alkyl Substances in Drinking Water (Continued)

| Isotope Dilution | LCSD      |           | Limits   |
|------------------|-----------|-----------|----------|
|                  | %Recovery | Qualifier |          |
| 13C3 HFPO-DA     | 86        |           | 50 - 200 |
| 13C6 PFDA        | 98        |           | 50 - 200 |
| 13C5 PFHxA       | 85        |           | 50 - 200 |
| 13C4 PFHpA       | 93        |           | 50 - 200 |
| 13C8 PFOA        | 96        |           | 50 - 200 |
| 13C9 PFNA        | 97        |           | 50 - 200 |
| 13C7 PFUnA       | 96        |           | 50 - 200 |
| 13C2 PFDoA       | 97        |           | 50 - 200 |
| 13C4 PFBA        | 84        |           | 50 - 200 |
| 13C5 PFPeA       | 86        |           | 50 - 200 |
| 13C3 PFBS        | 95        |           | 50 - 200 |
| 13C3 PFHxS       | 97        |           | 50 - 200 |
| 13C8 PFOS        | 96        |           | 50 - 200 |
| 13C2-4:2-FTS     | 106       |           | 50 - 200 |
| 13C2-6:2-FTS     | 94        |           | 50 - 200 |
| 13C2-8:2-FTS     | 94        |           | 50 - 200 |

**Lab Sample ID: MRL 380-53035/22-A**  
**Matrix: Water**  
**Analysis Batch: 53385**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 53035**

| Analyte   | Spike Added | MRL Result | MRL Qualifier | Unit | D | %Rec | %Rec Limits |
|---|-------------|------------|---------------|------|---|------|-------------|
| 11-Chloroeicosfluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS) | 2.01        | 1.85       | J             | ng/L |   | 92   | 50 - 150    |
| 9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid(9Cl-PF3ONS)    | 2.01        | 1.92       | J             | ng/L |   | 95   | 50 - 150    |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA)                       | 2.01        | 2.18       | J             | ng/L |   | 108  | 50 - 150    |
| Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)               | 2.01        | 1.73       | J             | ng/L |   | 86   | 50 - 150    |
| Perfluorobutanesulfonic acid (PFBS)                               | 2.01        | 2.08       | J             | ng/L |   | 104  | 50 - 150    |
| Perfluorodecanoic acid (PFDA)                                     | 2.01        | 1.98       | J             | ng/L |   | 99   | 50 - 150    |
| Perfluorododecanoic acid (PFDoA)                                  | 2.01        | 1.96       | J             | ng/L |   | 98   | 50 - 150    |
| Perfluoroheptanoic acid (PFHpA)                                   | 2.01        | 2.04       | J             | ng/L |   | 102  | 50 - 150    |
| Perfluorohexanesulfonic acid (PFHxS)                              | 2.01        | 2.08       | J             | ng/L |   | 104  | 50 - 150    |
| Perfluorohexanoic acid (PFHxA)                                    | 2.01        | 2.32       | J             | ng/L |   | 115  | 50 - 150    |
| Perfluorononanoic acid (PFNA)                                     | 2.01        | 2.16       | J             | ng/L |   | 108  | 50 - 150    |
| Perfluorooctanesulfonic acid (PFOS)                               | 2.01        | 2.04       | J             | ng/L |   | 102  | 50 - 150    |
| Perfluorooctanoic acid (PFOA)                                     | 2.01        | 2.14       | J             | ng/L |   | 107  | 50 - 150    |
| Perfluoroundecanoic acid (PFUnA)                                  | 2.01        | 1.99       | J             | ng/L |   | 99   | 50 - 150    |
| Perfluorobutanoic acid (PFBA)                                     | 2.01        | 2.13       | J             | ng/L |   | 106  | 50 - 150    |
| 1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)               | 2.01        | 2.21       | J             | ng/L |   | 110  | 50 - 150    |
| 1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)               | 2.01        | 2.06       | J             | ng/L |   | 103  | 50 - 150    |
| 1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)               | 2.01        | 2.16       | J             | ng/L |   | 108  | 50 - 150    |



# QC Sample Results

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

Job ID: 380-57979-1

## Method: 533 - Perfluorinated and Polyfluorinated Alkyl Substances in Drinking Water (Continued)

**Lab Sample ID: MRL 380-53035/22-A**  
**Matrix: Water**  
**Analysis Batch: 53385**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 53035**

| Analyte  | Spike Added | MRL Result | MRL Qualifier | Unit | D | %Rec | %Rec Limits |
|--|-------------|------------|---------------|------|---|------|-------------|
| Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)         | 2.01        | 2.15       | J             | ng/L |   | 107  | 50 - 150    |
| Perfluoro (2-ethoxyethane) sulfonic acid (PFEEESA) | 2.01        | 1.99       | J             | ng/L |   | 99   | 50 - 150    |
| Perfluoro-3-methoxypropanoic acid (PFMPA)          | 2.01        | 2.02       | J             | ng/L |   | 101  | 50 - 150    |
| Perfluoro-4-methoxybutanoic acid (PFMBA)           | 2.01        | 2.07       | J             | ng/L |   | 103  | 50 - 150    |
| Perfluoropentanoic acid (PFPeA)                    | 2.01        | 2.13       | J             | ng/L |   | 106  | 50 - 150    |
| Perfluoroheptanesulfonic acid (PFHpS)              | 2.01        | 2.15       | J             | ng/L |   | 107  | 50 - 150    |
| Perfluoropentanesulfonic acid (PFPeS)              | 2.01        | 1.90       | J             | ng/L |   | 95   | 50 - 150    |

| Isotope Dilution | MRL %Recovery | MRL Qualifier | MRL Limits |
|------------------|---------------|---------------|------------|
| 13C3 HFPO-DA     | 82            |               | 50 - 200   |
| 13C6 PFDA        | 93            |               | 50 - 200   |
| 13C5 PFHxA       | 90            |               | 50 - 200   |
| 13C4 PFHpA       | 90            |               | 50 - 200   |
| 13C8 PFOA        | 95            |               | 50 - 200   |
| 13C9 PFNA        | 93            |               | 50 - 200   |
| 13C7 PFUnA       | 91            |               | 50 - 200   |
| 13C2 PFDoA       | 93            |               | 50 - 200   |
| 13C4 PFBA        | 94            |               | 50 - 200   |
| 13C5 PFPeA       | 94            |               | 50 - 200   |
| 13C3 PFBS        | 87            |               | 50 - 200   |
| 13C3 PFHxS       | 88            |               | 50 - 200   |
| 13C8 PFOS        | 89            |               | 50 - 200   |
| 13C2-4:2-FTS     | 94            |               | 50 - 200   |
| 13C2-6:2-FTS     | 87            |               | 50 - 200   |
| 13C2-8:2-FTS     | 88            |               | 50 - 200   |

**Lab Sample ID: 380-57432-O-1-D LMS**  
**Matrix: Water**  
**Analysis Batch: 53385**

**Client Sample ID: Matrix Spike**  
**Prep Type: Total/NA**  
**Prep Batch: 53035**

| Analyte   | Sample Result | Sample Qualifier | Spike Added | LMS Result | LMS Qualifier | Unit | D | %Rec | %Rec Limits |
|---|---------------|------------------|-------------|------------|---------------|------|---|------|-------------|
| 11-Chloroeicosfluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS) | <2.0          |                  | 2.01        | 1.90       | J             | ng/L |   | 94   | 50 - 150    |
| 9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid(9Cl-PF3ONS)    | <2.0          |                  | 2.01        | 1.91       | J             | ng/L |   | 95   | 50 - 150    |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA)                       | <2.0          |                  | 2.01        | 1.90       | J             | ng/L |   | 95   | 50 - 150    |
| Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)               | <2.0          | *5-              | 2.01        | 1.85       | J *5-         | ng/L |   | 92   | 50 - 150    |
| Perfluorobutanesulfonic acid (PFBS)                               | <2.0          |                  | 2.01        | 2.73       |               | ng/L |   | 109  | 50 - 150    |
| Perfluorodecanoic acid (PFDA)                                     | <2.0          |                  | 2.01        | 2.13       |               | ng/L |   | 106  | 50 - 150    |
| Perfluorododecanoic acid (PFDoA)                                  | <2.0          |                  | 2.01        | 1.99       | J             | ng/L |   | 99   | 50 - 150    |

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# QC Sample Results

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

Job ID: 380-57979-1

## Method: 533 - Perfluorinated and Polyfluorinated Alkyl Substances in Drinking Water (Continued)

Lab Sample ID: 380-57432-O-1-D LMS

Matrix: Water

Analysis Batch: 53385

Client Sample ID: Matrix Spike

Prep Type: Total/NA

Prep Batch: 53035

| Analyte   | Sample Result | Sample Qualifier | Spike Added | LMS Result | LMS Qualifier | Unit | D | %Rec | %Rec Limits |
|---|---------------|------------------|-------------|------------|---------------|------|---|------|-------------|
| Perfluoroheptanoic acid (PFHpA)                     | <2.0          |                  | 2.01        | 2.72       |               | ng/L |   | 107  | 50 - 150    |
| Perfluorohexanesulfonic acid (PFHxS)                | <2.0          |                  | 2.01        | 3.16       |               | ng/L |   | 94   | 50 - 150    |
| Perfluorohexanoic acid (PFHxA)                      | <2.0          |                  | 2.01        | 3.38       |               | ng/L |   | 110  | 50 - 150    |
| Perfluorononanoic acid (PFNA)                       | <2.0          |                  | 2.01        | 2.28       |               | ng/L |   | 114  | 50 - 150    |
| Perfluorooctanesulfonic acid (PFOS)                 | <2.0          |                  | 2.01        | 3.66       |               | ng/L |   | 95   | 50 - 150    |
| Perfluorooctanoic acid (PFOA)                       | <2.0          |                  | 2.01        | 2.74       |               | ng/L |   | 94   | 50 - 150    |
| Perfluoroundecanoic acid (PFUnA)                    | <2.0          |                  | 2.01        | 2.17       |               | ng/L |   | 108  | 50 - 150    |
| Perfluorobutanoic acid (PFBA)                       | 2.2           |                  | 2.01        | 4.19       |               | ng/L |   | 101  | 50 - 150    |
| 1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS) | <2.0          |                  | 2.01        | 2.07       |               | ng/L |   | 103  | 50 - 150    |
| 1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS) | <2.0          |                  | 2.01        | 2.10       |               | ng/L |   | 104  | 50 - 150    |
| 1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS) | <2.0          |                  | 2.01        | 2.06       |               | ng/L |   | 103  | 50 - 150    |
| Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)          | <2.0          |                  | 2.01        | 1.86       | J             | ng/L |   | 93   | 50 - 150    |
| Perfluoro (2-ethoxyethane) sulfonic acid (PFEEESA)  | <2.0          |                  | 2.01        | 2.08       |               | ng/L |   | 104  | 50 - 150    |
| Perfluoro-3-methoxypropanoic acid (PFMPA)           | <2.0          |                  | 2.01        | 1.80       | J             | ng/L |   | 90   | 50 - 150    |
| Perfluoro-4-methoxybutanoic acid (PFMBA)            | <2.0          |                  | 2.01        | 2.23       |               | ng/L |   | 111  | 50 - 150    |
| Perfluoropentanoic acid (PFPeA)                     | <2.0          |                  | 2.01        | 3.35       |               | ng/L |   | 118  | 50 - 150    |
| Perfluoroheptanesulfonic acid (PFHpS)               | <2.0          |                  | 2.01        | 2.27       |               | ng/L |   | 113  | 50 - 150    |
| Perfluoropentanesulfonic acid (PFPeS)               | <2.0          |                  | 2.01        | 2.28       |               | ng/L |   | 113  | 50 - 150    |

| Isotope Dilution | LMS %Recovery | LMS Qualifier | Limits   |
|------------------|---------------|---------------|----------|
| 13C3 HFPO-DA     | 48            | *5-           | 50 - 200 |
| 13C6 PFDA        | 77            |               | 50 - 200 |
| 13C5 PFHxA       | 58            |               | 50 - 200 |
| 13C4 PFHpA       | 61            |               | 50 - 200 |
| 13C8 PFOA        | 69            |               | 50 - 200 |
| 13C9 PFNA        | 72            |               | 50 - 200 |
| 13C7 PFUnA       | 77            |               | 50 - 200 |
| 13C2 PFDoA       | 87            |               | 50 - 200 |
| 13C4 PFBA        | 63            |               | 50 - 200 |
| 13C5 PFPeA       | 57            |               | 50 - 200 |
| 13C3 PFBS        | 92            |               | 50 - 200 |
| 13C3 PFHxS       | 95            |               | 50 - 200 |
| 13C8 PFOS        | 93            |               | 50 - 200 |
| 13C2-4:2-FTS     | 103           |               | 50 - 200 |
| 13C2-6:2-FTS     | 96            |               | 50 - 200 |
| 13C2-8:2-FTS     | 94            |               | 50 - 200 |

# QC Sample Results

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

Job ID: 380-57979-1

## Method: 533 - Perfluorinated and Polyfluorinated Alkyl Substances in Drinking Water (Continued)

**Lab Sample ID: 380-57432-P-1-C LMSD**

**Matrix: Water**

**Analysis Batch: 53385**

**Client Sample ID: Matrix Spike Duplicate**

**Prep Type: Total/NA**

**Prep Batch: 53035**

| Analyte   | Sample | Sample    | Spike | LMSD   | LMSD      | Unit | D | %Rec | %Rec     | RPD | RPD |
|---|--------|-----------|-------|--------|-----------|------|---|------|----------|-----|-----|
|   | Result | Qualifier | Added | Result | Qualifier |      |   |      | Limits   |     |     |
| 11-Chloroeicosfluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS) | <2.0   |           | 2.00  | 1.85   | J         | ng/L |   | 93   | 50 - 150 | 2   | 50  |
| 9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid(9Cl-PF3ONS)    | <2.0   |           | 2.00  | 1.96   | J         | ng/L |   | 98   | 50 - 150 | 3   | 50  |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA)                       | <2.0   |           | 2.00  | 2.20   |           | ng/L |   | 110  | 50 - 150 | 14  | 50  |
| Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)               | <2.0   | *5-       | 2.00  | 1.78   | J         | ng/L |   | 89   | 50 - 150 | 4   | 50  |
| Perfluorobutanesulfonic acid (PFBS)                               | <2.0   |           | 2.00  | 2.82   |           | ng/L |   | 114  | 50 - 150 | 3   | 50  |
| Perfluorodecanoic acid (PFDA)                                     | <2.0   |           | 2.00  | 2.13   |           | ng/L |   | 106  | 50 - 150 | 0   | 50  |
| Perfluorododecanoic acid (PFDoA)                                  | <2.0   |           | 2.00  | 2.07   |           | ng/L |   | 103  | 50 - 150 | 4   | 50  |
| Perfluoroheptanoic acid (PFHpA)                                   | <2.0   |           | 2.00  | 2.87   |           | ng/L |   | 115  | 50 - 150 | 5   | 50  |
| Perfluorohexanesulfonic acid (PFHxS)                              | <2.0   |           | 2.00  | 3.19   |           | ng/L |   | 96   | 50 - 150 | 1   | 50  |
| Perfluorohexanoic acid (PFHxA)                                    | <2.0   |           | 2.00  | 3.47   |           | ng/L |   | 114  | 50 - 150 | 2   | 50  |
| Perfluorononanoic acid (PFNA)                                     | <2.0   |           | 2.00  | 2.40   |           | ng/L |   | 120  | 50 - 150 | 5   | 50  |
| Perfluorooctanesulfonic acid (PFOS)                               | <2.0   |           | 2.00  | 3.63   |           | ng/L |   | 94   | 50 - 150 | 1   | 50  |
| Perfluorooctanoic acid (PFOA)                                     | <2.0   |           | 2.00  | 2.88   |           | ng/L |   | 101  | 50 - 150 | 5   | 50  |
| Perfluoroundecanoic acid (PFUnA)                                  | <2.0   |           | 2.00  | 2.08   |           | ng/L |   | 104  | 50 - 150 | 4   | 50  |
| Perfluorobutanoic acid (PFBA)                                     | 2.2    |           | 2.00  | 4.33   |           | ng/L |   | 108  | 50 - 150 | 3   | 50  |
| 1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)               | <2.0   |           | 2.00  | 2.17   |           | ng/L |   | 108  | 50 - 150 | 4   | 50  |
| 1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)               | <2.0   |           | 2.00  | 2.01   |           | ng/L |   | 100  | 50 - 150 | 4   | 50  |
| 1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)               | <2.0   |           | 2.00  | 2.37   |           | ng/L |   | 118  | 50 - 150 | 14  | 50  |
| Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)                        | <2.0   |           | 2.00  | 1.96   | J         | ng/L |   | 98   | 50 - 150 | 5   | 50  |
| Perfluoro (2-ethoxyethane) sulfonic acid (PFEEESA)                | <2.0   |           | 2.00  | 2.22   |           | ng/L |   | 111  | 50 - 150 | 6   | 50  |
| Perfluoro-3-methoxypropanoic acid (PFMPA)                         | <2.0   |           | 2.00  | 1.80   | J         | ng/L |   | 90   | 50 - 150 | 0   | 50  |
| Perfluoro-4-methoxybutanoic acid (PFMBA)                          | <2.0   |           | 2.00  | 2.31   |           | ng/L |   | 115  | 50 - 150 | 3   | 50  |
| Perfluoropentanoic acid (PFPeA)                                   | <2.0   |           | 2.00  | 3.53   |           | ng/L |   | 127  | 50 - 150 | 5   | 50  |
| Perfluoroheptanesulfonic acid (PFHpS)                             | <2.0   |           | 2.00  | 2.23   |           | ng/L |   | 111  | 50 - 150 | 2   | 50  |
| Perfluoropentanesulfonic acid (PFPeS)                             | <2.0   |           | 2.00  | 2.29   |           | ng/L |   | 114  | 50 - 150 | 0   | 50  |

| Isotope Dilution | LMSD      |           | Limits   |
|------------------|-----------|-----------|----------|
|                  | %Recovery | Qualifier |          |
| 13C3 HFPO-DA     | 68        |           | 50 - 200 |
| 13C6 PFDA        | 84        |           | 50 - 200 |
| 13C5 PFHxA       | 78        |           | 50 - 200 |
| 13C4 PFHpA       | 78        |           | 50 - 200 |
| 13C8 PFOA        | 85        |           | 50 - 200 |
| 13C9 PFNA        | 85        |           | 50 - 200 |

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# QC Sample Results

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

Job ID: 380-57979-1

## Method: 533 - Perfluorinated and Polyfluorinated Alkyl Substances in Drinking Water (Continued)

**Lab Sample ID: 380-57432-P-1-C LMSD**  
**Matrix: Water**  
**Analysis Batch: 53385**

**Client Sample ID: Matrix Spike Duplicate**  
**Prep Type: Total/NA**  
**Prep Batch: 53035**

| Isotope Dilution | LMSD LMSD |           | Limits   |
|------------------|-----------|-----------|----------|
|                  | %Recovery | Qualifier |          |
| 13C7 PFUnA       | 85        |           | 50 - 200 |
| 13C2 PFDoA       | 90        |           | 50 - 200 |
| 13C4 PFBA        | 83        |           | 50 - 200 |
| 13C5 PFPeA       | 78        |           | 50 - 200 |
| 13C3 PFBS        | 88        |           | 50 - 200 |
| 13C3 PFHxS       | 98        |           | 50 - 200 |
| 13C8 PFOS        | 97        |           | 50 - 200 |
| 13C2-4:2-FTS     | 112       |           | 50 - 200 |
| 13C2-6:2-FTS     | 97        |           | 50 - 200 |
| 13C2-8:2-FTS     | 99        |           | 50 - 200 |

## Method: 537.1 - Perfluorinated Alkyl Acids (LC/MS)

**Lab Sample ID: MBL 380-51108/23-A**  
**Matrix: Water**  
**Analysis Batch: 51333**

**Client Sample ID: Method Blank**  
**Prep Type: Total/NA**  
**Prep Batch: 51108**

| Analyte  | MBL    | MBL       | RL  | Unit | D | Prepared       | Analyzed       | Dil Fac |
|--|--------|-----------|-----|------|---|----------------|----------------|---------|
|  | Result | Qualifier |     |      |   |                |                |         |
| Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)                | <1.0   |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| Perfluorooctanesulfonic acid (PFOS)                                | <0.43  |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| Perfluoroundecanoic acid (PFUnA)                                   | <0.42  |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)           | <0.58  |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)            | <0.42  |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| Perfluorohexanoic acid (PFHxA)                                     | <0.46  |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| Perfluorododecanoic acid (PFDoA)                                   | <0.54  |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| Perfluorooctanoic acid (PFOA)                                      | <0.38  |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| Perfluorodecanoic acid (PFDA)                                      | <0.31  |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| Perfluorohexanesulfonic acid (PFHxS)                               | <0.32  |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| Perfluorobutanesulfonic acid (PFBS)                                | <0.37  |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| Perfluoroheptanoic acid (PFHpA)                                    | <0.39  |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| Perfluorononanoic acid (PFNA)                                      | <0.40  |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| Perfluorotetradecanoic acid (PFTA)                                 | <0.54  |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| Perfluorotridecanoic acid (PFTrDA)                                 | <0.36  |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| 9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid(9Cl-PF3ONS)     | <0.30  |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| 11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS) | <0.30  |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA)                        | <0.60  |           | 2.0 | ng/L |   | 08/10/23 05:33 | 08/11/23 10:57 | 1       |

| Surrogate   | MBL MBL   |           | Limits   | Prepared       | Analyzed       | Dil Fac |
|-------------|-----------|-----------|----------|----------------|----------------|---------|
|             | %Recovery | Qualifier |          |                |                |         |
| d5-NEtFOSAA | 109       |           | 70 - 130 | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| 13C2 PFHxA  | 115       |           | 70 - 130 | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| 13C2 PFDA   | 121       |           | 70 - 130 | 08/10/23 05:33 | 08/11/23 10:57 | 1       |
| 13C3-GenX   | 106       |           | 70 - 130 | 08/10/23 05:33 | 08/11/23 10:57 | 1       |

# QC Sample Results

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

Job ID: 380-57979-1

## Method: 537.1 - Perfluorinated Alkyl Acids (LC/MS) (Continued)

**Lab Sample ID: LCS 380-51108/25-A**  
**Matrix: Water**  
**Analysis Batch: 51333**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 51108**

| Analyte  | Spike Added | LCS Result | LCS Qualifier | Unit | D | %Rec | %Rec Limits |
|--|-------------|------------|---------------|------|---|------|-------------|
| Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)                | 25.1        | 25.8       |               | ng/L |   | 103  | 70 - 130    |
| Perfluorooctanesulfonic acid (PFOS)                                | 23.2        | 24.4       |               | ng/L |   | 105  | 70 - 130    |
| Perfluoroundecanoic acid (PFUnA)                                   | 25.1        | 25.2       |               | ng/L |   | 101  | 70 - 130    |
| N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)           | 25.1        | 26.2       |               | ng/L |   | 104  | 70 - 130    |
| N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)            | 25.1        | 25.0       |               | ng/L |   | 100  | 70 - 130    |
| Perfluorohexanoic acid (PFHxA)                                     | 25.1        | 28.2       |               | ng/L |   | 113  | 70 - 130    |
| Perfluorododecanoic acid (PFDoA)                                   | 25.1        | 25.0       |               | ng/L |   | 100  | 70 - 130    |
| Perfluorooctanoic acid (PFOA)                                      | 25.1        | 27.0       |               | ng/L |   | 108  | 70 - 130    |
| Perfluorodecanoic acid (PFDA)                                      | 25.1        | 26.8       |               | ng/L |   | 107  | 70 - 130    |
| Perfluorohexanesulfonic acid (PFHxS)                               | 22.9        | 23.9       |               | ng/L |   | 105  | 70 - 130    |
| Perfluorobutanesulfonic acid (PFBS)                                | 22.2        | 24.2       |               | ng/L |   | 109  | 70 - 130    |
| Perfluoroheptanoic acid (PFHpA)                                    | 25.1        | 25.0       |               | ng/L |   | 100  | 70 - 130    |
| Perfluorononanoic acid (PFNA)                                      | 25.1        | 29.5       |               | ng/L |   | 118  | 70 - 130    |
| Perfluorotetradecanoic acid (PFTA)                                 | 25.1        | 21.6       |               | ng/L |   | 86   | 70 - 130    |
| Perfluorotridecanoic acid (PFTrDA)                                 | 25.1        | 27.9       |               | ng/L |   | 111  | 70 - 130    |
| 9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid(9Cl-PF3ONS)     | 23.4        | 26.3       |               | ng/L |   | 112  | 70 - 130    |
| 11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS) | 23.7        | 24.5       |               | ng/L |   | 103  | 70 - 130    |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA)                        | 23.7        | 26.1       |               | ng/L |   | 110  | 70 - 130    |

| Surrogate   | LCS %Recovery | LCS Qualifier | Limits   |
|-------------|---------------|---------------|----------|
| d5-NEtFOSAA | 108           |               | 70 - 130 |
| 13C2 PFHxA  | 120           |               | 70 - 130 |
| 13C2 PFDA   | 112           |               | 70 - 130 |
| 13C3-GenX   | 110           |               | 70 - 130 |

**Lab Sample ID: LCSD 380-51108/26-A**  
**Matrix: Water**  
**Analysis Batch: 51333**

**Client Sample ID: Lab Control Sample Dup**  
**Prep Type: Total/NA**  
**Prep Batch: 51108**

| Analyte  | Spike Added | LCSD Result | LCSD Qualifier | Unit | D | %Rec | %Rec Limits | RPD | RPD Limit |
|--|-------------|-------------|----------------|------|---|------|-------------|-----|-----------|
| Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)      | 25.1        | 25.8        |                | ng/L |   | 103  | 70 - 130    | 0   | 30        |
| Perfluorooctanesulfonic acid (PFOS)                      | 23.2        | 26.0        |                | ng/L |   | 112  | 70 - 130    | 6   | 30        |
| Perfluoroundecanoic acid (PFUnA)                         | 25.1        | 26.8        |                | ng/L |   | 107  | 70 - 130    | 6   | 30        |
| N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA) | 25.1        | 27.7        |                | ng/L |   | 110  | 70 - 130    | 6   | 30        |

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# QC Sample Results

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

Job ID: 380-57979-1

## Method: 537.1 - Perfluorinated Alkyl Acids (LC/MS) (Continued)

**Lab Sample ID: LCSD 380-51108/26-A**  
**Matrix: Water**  
**Analysis Batch: 51333**

**Client Sample ID: Lab Control Sample Dup**  
**Prep Type: Total/NA**  
**Prep Batch: 51108**

| Analyte   | Spike Added | LCSD Result | LCSD Qualifier | Unit | D | %Rec | %Rec Limits | RPD | RPD Limit |
|---|-------------|-------------|----------------|------|---|------|-------------|-----|-----------|
| N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)           | 25.1        | 25.7        |                | ng/L |   | 102  | 70 - 130    | 3   | 30        |
| Perfluorohexanoic acid (PFHxA)                                    | 25.1        | 30.6        |                | ng/L |   | 122  | 70 - 130    | 8   | 30        |
| Perfluorododecanoic acid (PFDoA)                                  | 25.1        | 26.8        |                | ng/L |   | 107  | 70 - 130    | 7   | 30        |
| Perfluorooctanoic acid (PFOA)                                     | 25.1        | 28.2        |                | ng/L |   | 112  | 70 - 130    | 4   | 30        |
| Perfluorodecanoic acid (PFDA)                                     | 25.1        | 27.3        |                | ng/L |   | 109  | 70 - 130    | 2   | 30        |
| Perfluorohexanesulfonic acid (PFHxS)                              | 22.9        | 25.2        |                | ng/L |   | 110  | 70 - 130    | 5   | 30        |
| Perfluorobutanesulfonic acid (PFBS)                               | 22.2        | 26.4        |                | ng/L |   | 119  | 70 - 130    | 9   | 30        |
| Perfluoroheptanoic acid (PFHpA)                                   | 25.1        | 26.0        |                | ng/L |   | 104  | 70 - 130    | 4   | 30        |
| Perfluorononanoic acid (PFNA)                                     | 25.1        | 31.4        |                | ng/L |   | 125  | 70 - 130    | 6   | 30        |
| Perfluorotetradecanoic acid (PFTA)                                | 25.1        | 23.5        |                | ng/L |   | 94   | 70 - 130    | 8   | 30        |
| Perfluorotridecanoic acid (PFTrDA)                                | 25.1        | 29.5        |                | ng/L |   | 117  | 70 - 130    | 6   | 30        |
| 9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid(9Cl-PF3ONS)    | 23.5        | 27.7        |                | ng/L |   | 118  | 70 - 130    | 5   | 30        |
| 11-Chloroeicosafuoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS) | 23.7        | 26.7        |                | ng/L |   | 112  | 70 - 130    | 9   | 30        |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA)                       | 23.7        | 26.8        |                | ng/L |   | 113  | 70 - 130    | 2   | 30        |

| Surrogate   | LCSD %Recovery | LCSD Qualifier | LCSD Limits |
|-------------|----------------|----------------|-------------|
| d5-NEtFOSAA | 110            |                | 70 - 130    |
| 13C2 PFHxA  | 128            |                | 70 - 130    |
| 13C2 PFDA   | 113            |                | 70 - 130    |
| 13C3-GenX   | 130            |                | 70 - 130    |

**Lab Sample ID: MRL 380-51108/24-A**  
**Matrix: Water**  
**Analysis Batch: 51440**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 51108**

| Analyte  | Spike Added | MRL Result | MRL Qualifier | Unit | D | %Rec | %Rec Limits |
|--|-------------|------------|---------------|------|---|------|-------------|
| Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)      | 2.00        | 1.94       | J             | ng/L |   | 97   | 50 - 150    |
| Perfluorooctanesulfonic acid (PFOS)                      | 1.86        | 1.93       | J             | ng/L |   | 104  | 50 - 150    |
| Perfluoroundecanoic acid (PFUnA)                         | 2.00        | 2.09       | J             | ng/L |   | 104  | 50 - 150    |
| N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA) | 2.00        | 2.01       | J             | ng/L |   | 100  | 50 - 150    |
| N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)  | 2.00        | 1.77       | J             | ng/L |   | 88   | 50 - 150    |
| Perfluorohexanoic acid (PFHxA)                           | 2.00        | 2.34       | J             | ng/L |   | 117  | 50 - 150    |
| Perfluorododecanoic acid (PFDoA)                         | 2.00        | 2.10       | J             | ng/L |   | 105  | 50 - 150    |
| Perfluorooctanoic acid (PFOA)                            | 2.00        | 2.40       | J             | ng/L |   | 120  | 50 - 150    |
| Perfluorodecanoic acid (PFDA)                            | 2.00        | 2.13       | J             | ng/L |   | 106  | 50 - 150    |

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# QC Sample Results

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

Job ID: 380-57979-1

## Method: 537.1 - Perfluorinated Alkyl Acids (LC/MS) (Continued)

**Lab Sample ID: MRL 380-51108/24-A**  
**Matrix: Water**  
**Analysis Batch: 51440**

**Client Sample ID: Lab Control Sample**  
**Prep Type: Total/NA**  
**Prep Batch: 51108**

| Analyte  | Spike Added | MRL Result | MRL Qualifier | Unit | D | %Rec | %Rec Limits |
|--|-------------|------------|---------------|------|---|------|-------------|
| Perfluorohexanesulfonic acid (PFHxS)                               | 1.83        | 1.77       | J             | ng/L |   | 97   | 50 - 150    |
| Perfluorobutanesulfonic acid (PFBS)                                | 1.77        | 2.00       | J             | ng/L |   | 113  | 50 - 150    |
| Perfluoroheptanoic acid (PFHpA)                                    | 2.00        | 1.98       | J             | ng/L |   | 99   | 50 - 150    |
| Perfluorononanoic acid (PFNA)                                      | 2.00        | 2.57       | J             | ng/L |   | 128  | 50 - 150    |
| Perfluorotetradecanoic acid (PFTA)                                 | 2.00        | 2.01       | J             | ng/L |   | 100  | 50 - 150    |
| Perfluorotridecanoic acid (PFTTrDA)                                | 2.00        | 2.15       | J             | ng/L |   | 107  | 50 - 150    |
| 9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid(9Cl-PF3ONS)     | 1.87        | 2.01       | J             | ng/L |   | 107  | 50 - 150    |
| 11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS) | 1.89        | 1.77       | J             | ng/L |   | 94   | 50 - 150    |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA)                        | 1.89        | 2.07       | J             | ng/L |   | 109  | 50 - 150    |

| Surrogate   | MRL %Recovery | MRL Qualifier | MRL Limits |
|-------------|---------------|---------------|------------|
| d5-NEtFOSAA | 97            |               | 70 - 130   |
| 13C2 PFHxA  | 122           |               | 70 - 130   |
| 13C2 PFDA   | 110           |               | 70 - 130   |
| 13C3-GenX   | 113           |               | 70 - 130   |

**Lab Sample ID: 380-58087-B-1-C LMS**  
**Matrix: Water**  
**Analysis Batch: 51333**

**Client Sample ID: Matrix Spike**  
**Prep Type: Total/NA**  
**Prep Batch: 51108**

| Analyte  | Sample Result | Sample Qualifier | Spike Added | LMS Result | LMS Qualifier | Unit | D | %Rec | %Rec Limits |
|--|---------------|------------------|-------------|------------|---------------|------|---|------|-------------|
| Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)      | <2.0          |                  | 2.02        | 2.29       |               | ng/L |   | 113  | 50 - 150    |
| Perfluorooctanesulfonic acid (PFOS)                      | <2.0          |                  | 1.87        | 2.06       |               | ng/L |   | 110  | 50 - 150    |
| Perfluoroundecanoic acid (PFUnA)                         | <2.0          |                  | 2.02        | 2.31       |               | ng/L |   | 115  | 50 - 150    |
| N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA) | <2.0          |                  | 2.02        | 2.30       |               | ng/L |   | 114  | 50 - 150    |
| N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)  | <2.0          |                  | 2.02        | 2.22       |               | ng/L |   | 110  | 50 - 150    |
| Perfluorohexanoic acid (PFHxA)                           | 12            |                  | 2.02        | 15.7       | 4             | ng/L |   | 203  | 50 - 150    |
| Perfluorododecanoic acid (PFDoA)                         | <2.0          |                  | 2.02        | 2.46       |               | ng/L |   | 122  | 50 - 150    |
| Perfluorooctanoic acid (PFOA)                            | 3.2           |                  | 2.02        | 5.73       |               | ng/L |   | 127  | 50 - 150    |
| Perfluorodecanoic acid (PFDA)                            | <2.0          |                  | 2.02        | 2.35       |               | ng/L |   | 116  | 50 - 150    |
| Perfluorohexanesulfonic acid (PFHxS)                     | <2.0          |                  | 1.84        | 2.36       |               | ng/L |   | 104  | 50 - 150    |
| Perfluorobutanesulfonic acid (PFBS)                      | 8.2           |                  | 1.78        | 10.5       | 4             | ng/L |   | 128  | 50 - 150    |
| Perfluoroheptanoic acid (PFHpA)                          | 3.4           |                  | 2.02        | 6.00       |               | ng/L |   | 127  | 50 - 150    |
| Perfluorononanoic acid (PFNA)                            | <2.0          |                  | 2.02        | 2.98       |               | ng/L |   | 148  | 50 - 150    |
| Perfluorotetradecanoic acid (PFTA)                       | <2.0          |                  | 2.02        | 2.38       |               | ng/L |   | 118  | 50 - 150    |

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# QC Sample Results

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

Job ID: 380-57979-1

## Method: 537.1 - Perfluorinated Alkyl Acids (LC/MS) (Continued)

**Lab Sample ID: 380-58087-B-1-C LMS**  
**Matrix: Water**  
**Analysis Batch: 51333**

**Client Sample ID: Matrix Spike**  
**Prep Type: Total/NA**  
**Prep Batch: 51108**

| Analyte   | Sample Result    | Sample Qualifier | Spike Added   | LMS Result | LMS Qualifier | Unit | D | %Rec | %Rec Limits |
|---|------------------|------------------|---------------|------------|---------------|------|---|------|-------------|
| Perfluorotridecanoic acid (PFTTrDA)                               | <2.0             |                  | 2.02          | 2.64       |               | ng/L |   | 131  | 50 - 150    |
| 9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid(9Cl-PF3ONS)    | <2.0             |                  | 1.89          | 2.15       |               | ng/L |   | 114  | 50 - 150    |
| 11-Chloroeicosafuoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS) | <2.0             |                  | 1.91          | 1.94       | J             | ng/L |   | 102  | 50 - 150    |
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA)                       | <2.0             |                  | 1.91          | 2.39       |               | ng/L |   | 126  | 50 - 150    |
| <b>LMS LMS</b>  |                  |                  |               |            |               |      |   |      |             |
| <b>Surrogate</b>  | <b>%Recovery</b> | <b>Qualifier</b> | <b>Limits</b> |            |               |      |   |      |             |
| d5-NEtFOSAA   | 117              |                  | 70 - 130      |            |               |      |   |      |             |
| 13C2 PFHxA  | 129              |                  | 70 - 130      |            |               |      |   |      |             |
| 13C2 PFDA   | 119              |                  | 70 - 130      |            |               |      |   |      |             |
| 13C3-GenX   | 119              |                  | 70 - 130      |            |               |      |   |      |             |

**Lab Sample ID: 380-58087-C-1-C LMSD**  
**Matrix: Water**  
**Analysis Batch: 51333**

**Client Sample ID: Matrix Spike Duplicate**  
**Prep Type: Total/NA**  
**Prep Batch: 51108**

| Analyte   | Sample Result | Sample Qualifier | Spike Added | LMSD Result | LMSD Qualifier | Unit | D | %Rec | %Rec Limits | RPD | RPD Limit |
|---|---------------|------------------|-------------|-------------|----------------|------|---|------|-------------|-----|-----------|
| Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)               | <2.0          |                  | 2.01        | 1.84        | J              | ng/L |   | 92   | 50 - 150    | 22  | 50        |
| Perfluorooctanesulfonic acid (PFOS)                               | <2.0          |                  | 1.86        | 1.97        | J              | ng/L |   | 106  | 50 - 150    | 5   | 50        |
| Perfluoroundecanoic acid (PFUnA)                                  | <2.0          |                  | 2.01        | 2.07        |                | ng/L |   | 103  | 50 - 150    | 11  | 50        |
| N-methylperfluorooctanesulfonamidoacetic acid (NMeFOSAA)          | <2.0          |                  | 2.01        | 2.10        |                | ng/L |   | 105  | 50 - 150    | 9   | 50        |
| N-ethylperfluorooctanesulfonamidoacetic acid (NEtFOSAA)           | <2.0          |                  | 2.01        | 2.12        |                | ng/L |   | 106  | 50 - 150    | 4   | 50        |
| Perfluorohexanoic acid (PFHxA)                                    | 12            |                  | 2.01        | 14.0        | 4              | ng/L |   | 121  | 50 - 150    | 11  | 50        |
| Perfluorododecanoic acid (PFDoA)                                  | <2.0          |                  | 2.01        | 2.20        |                | ng/L |   | 110  | 50 - 150    | 11  | 50        |
| Perfluorooctanoic acid (PFOA)                                     | 3.2           |                  | 2.01        | 5.22        |                | ng/L |   | 102  | 50 - 150    | 9   | 50        |
| Perfluorodecanoic acid (PFDA)                                     | <2.0          |                  | 2.01        | 2.15        |                | ng/L |   | 107  | 50 - 150    | 9   | 50        |
| Perfluorohexanesulfonic acid (PFHxS)                              | <2.0          |                  | 1.83        | 2.28        |                | ng/L |   | 100  | 50 - 150    | 3   | 50        |
| Perfluorobutanesulfonic acid (PFBS)                               | 8.2           |                  | 1.78        | 10.2        | 4              | ng/L |   | 113  | 50 - 150    | 3   | 50        |
| Perfluoroheptanoic acid (PFHpA)                                   | 3.4           |                  | 2.01        | 5.45        |                | ng/L |   | 100  | 50 - 150    | 10  | 50        |
| Perfluorononanoic acid (PFNA)                                     | <2.0          |                  | 2.01        | 2.60        |                | ng/L |   | 129  | 50 - 150    | 14  | 50        |
| Perfluorotetradecanoic acid (PFTA)                                | <2.0          |                  | 2.01        | 2.19        |                | ng/L |   | 109  | 50 - 150    | 8   | 50        |
| Perfluorotridecanoic acid (PFTTrDA)                               | <2.0          |                  | 2.01        | 2.46        |                | ng/L |   | 122  | 50 - 150    | 7   | 50        |
| 9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid(9Cl-PF3ONS)    | <2.0          |                  | 1.88        | 2.15        |                | ng/L |   | 114  | 50 - 150    | 0   | 50        |
| 11-Chloroeicosafuoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS) | <2.0          |                  | 1.90        | 1.90        | J              | ng/L |   | 100  | 50 - 150    | 2   | 50        |

Eurofins Eaton Analytical Pomona



# QC Sample Results

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

Job ID: 380-57979-1

## Method: 537.1 - Perfluorinated Alkyl Acids (LC/MS) (Continued)

**Lab Sample ID: 380-58087-C-1-C LMSD**

**Client Sample ID: Matrix Spike Duplicate**

**Matrix: Water**

**Prep Type: Total/NA**

**Analysis Batch: 51333**

**Prep Batch: 51108**

| Analyte                                     | Sample Result         | Sample Qualifier      | Spike Added   | LMSD Result | LMSD Qualifier | Unit | D | %Rec | %Rec Limits | RPD | RPD Limit |
|---|-----------------------|-----------------------|---------------|-------------|----------------|------|---|------|-------------|-----|-----------|
| 4,8-Dioxa-3H-perfluorononanoic acid (ADONA) | <2.0                  |                       | 1.90          | 2.17        |                | ng/L |   | 114  | 50 - 150    | 10  | 50        |
| <b>Surrogate</b>                            | <b>LMSD %Recovery</b> | <b>LMSD Qualifier</b> | <b>Limits</b> |             |                |      |   |      |             |     |           |
| d5-NEtFOSAA                                 | 114                   |                       | 70 - 130      |             |                |      |   |      |             |     |           |
| 13C2 PFHxA                                  | 117                   |                       | 70 - 130      |             |                |      |   |      |             |     |           |
| 13C2 PFDA                                   | 108                   |                       | 70 - 130      |             |                |      |   |      |             |     |           |
| 13C3-GenX                                   | 107                   |                       | 70 - 130      |             |                |      |   |      |             |     |           |

# QC Association Summary

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

Job ID: 380-57979-1

## GC/MS Semi VOA

### Prep Batch: 51109

| Lab Sample ID       | Client Sample ID         | Prep Type | Matrix         | Method | Prep Batch |
|---------------------|--------------------------|-----------|----------------|--------|------------|
| 380-57979-1         | HALAWA WELLS UNITS 1 & 2 | Total/NA  | Drinking Water | 525.2  |            |
| MB 380-51109/1-A    | Method Blank             | Total/NA  | Water          | 525.2  |            |
| LCS 380-51109/3-A   | Lab Control Sample       | Total/NA  | Water          | 525.2  |            |
| LCSD 380-51109/4-A  | Lab Control Sample Dup   | Total/NA  | Water          | 525.2  |            |
| MRL 380-51109/2-A   | Lab Control Sample       | Total/NA  | Water          | 525.2  |            |
| 380-57897-P-1-A MS  | Matrix Spike             | Total/NA  | Water          | 525.2  |            |
| 380-57909-AR-1-A DU | Duplicate                | Total/NA  | Water          | 525.2  |            |

### Analysis Batch: 51305

| Lab Sample ID       | Client Sample ID         | Prep Type | Matrix         | Method | Prep Batch |
|---------------------|--------------------------|-----------|----------------|--------|------------|
| 380-57979-1         | HALAWA WELLS UNITS 1 & 2 | Total/NA  | Drinking Water | 525.2  | 51109      |
| MB 380-51109/1-A    | Method Blank             | Total/NA  | Water          | 525.2  | 51109      |
| LCS 380-51109/3-A   | Lab Control Sample       | Total/NA  | Water          | 525.2  | 51109      |
| LCSD 380-51109/4-A  | Lab Control Sample Dup   | Total/NA  | Water          | 525.2  | 51109      |
| MRL 380-51109/2-A   | Lab Control Sample       | Total/NA  | Water          | 525.2  | 51109      |
| 380-57897-P-1-A MS  | Matrix Spike             | Total/NA  | Water          | 525.2  | 51109      |
| 380-57909-AR-1-A DU | Duplicate                | Total/NA  | Water          | 525.2  | 51109      |

## LCMS

### Prep Batch: 51108

| Lab Sample ID        | Client Sample ID             | Prep Type | Matrix         | Method   | Prep Batch |
|----------------------|------------------------------|-----------|----------------|----------|------------|
| 380-57979-1          | HALAWA WELLS UNITS 1 & 2     | Total/NA  | Drinking Water | 537.1 DW |            |
| 380-57979-3          | FB: HALAWA WELLS UNITS 1 & 2 | Total/NA  | Water          | 537.1 DW |            |
| MBL 380-51108/23-A   | Method Blank                 | Total/NA  | Water          | 537.1 DW |            |
| LCS 380-51108/25-A   | Lab Control Sample           | Total/NA  | Water          | 537.1 DW |            |
| LCSD 380-51108/26-A  | Lab Control Sample Dup       | Total/NA  | Water          | 537.1 DW |            |
| MRL 380-51108/24-A   | Lab Control Sample           | Total/NA  | Water          | 537.1 DW |            |
| 380-58087-B-1-C LMS  | Matrix Spike                 | Total/NA  | Water          | 537.1 DW |            |
| 380-58087-C-1-C LMSD | Matrix Spike Duplicate       | Total/NA  | Water          | 537.1 DW |            |

### Analysis Batch: 51333

| Lab Sample ID        | Client Sample ID             | Prep Type | Matrix | Method | Prep Batch |
|----------------------|------------------------------|-----------|--------|--------|------------|
| 380-57979-3          | FB: HALAWA WELLS UNITS 1 & 2 | Total/NA  | Water  | 537.1  | 51108      |
| MBL 380-51108/23-A   | Method Blank                 | Total/NA  | Water  | 537.1  | 51108      |
| LCS 380-51108/25-A   | Lab Control Sample           | Total/NA  | Water  | 537.1  | 51108      |
| LCSD 380-51108/26-A  | Lab Control Sample Dup       | Total/NA  | Water  | 537.1  | 51108      |
| 380-58087-B-1-C LMS  | Matrix Spike                 | Total/NA  | Water  | 537.1  | 51108      |
| 380-58087-C-1-C LMSD | Matrix Spike Duplicate       | Total/NA  | Water  | 537.1  | 51108      |

### Analysis Batch: 51440

| Lab Sample ID      | Client Sample ID         | Prep Type | Matrix         | Method | Prep Batch |
|--------------------|--------------------------|-----------|----------------|--------|------------|
| 380-57979-1        | HALAWA WELLS UNITS 1 & 2 | Total/NA  | Drinking Water | 537.1  | 51108      |
| MRL 380-51108/24-A | Lab Control Sample       | Total/NA  | Water          | 537.1  | 51108      |

### Prep Batch: 53035

| Lab Sample ID      | Client Sample ID             | Prep Type | Matrix         | Method | Prep Batch |
|--------------------|------------------------------|-----------|----------------|--------|------------|
| 380-57979-1        | HALAWA WELLS UNITS 1 & 2     | Total/NA  | Drinking Water | 533    |            |
| 380-57979-3        | FB: HALAWA WELLS UNITS 1 & 2 | Total/NA  | Water          | 533    |            |
| MBL 380-53035/21-A | Method Blank                 | Total/NA  | Water          | 533    |            |
| LCS 380-53035/23-A | Lab Control Sample           | Total/NA  | Water          | 533    |            |

Eurofins Eaton Analytical Pomona

# QC Association Summary

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

Job ID: 380-57979-1

## LCMS (Continued)

### Prep Batch: 53035 (Continued)

| Lab Sample ID        | Client Sample ID       | Prep Type | Matrix | Method | Prep Batch |
|----------------------|------------------------|-----------|--------|--------|------------|
| LCSD 380-53035/24-A  | Lab Control Sample Dup | Total/NA  | Water  | 533    |            |
| MRL 380-53035/22-A   | Lab Control Sample     | Total/NA  | Water  | 533    |            |
| 380-57432-O-1-D LMS  | Matrix Spike           | Total/NA  | Water  | 533    |            |
| 380-57432-P-1-C LMSD | Matrix Spike Duplicate | Total/NA  | Water  | 533    |            |

### Analysis Batch: 53385

| Lab Sample ID        | Client Sample ID             | Prep Type | Matrix         | Method | Prep Batch |
|----------------------|------------------------------|-----------|----------------|--------|------------|
| 380-57979-1          | HALAWA WELLS UNITS 1 & 2     | Total/NA  | Drinking Water | 533    | 53035      |
| 380-57979-3          | FB: HALAWA WELLS UNITS 1 & 2 | Total/NA  | Water          | 533    | 53035      |
| MBL 380-53035/21-A   | Method Blank                 | Total/NA  | Water          | 533    | 53035      |
| LCS 380-53035/23-A   | Lab Control Sample           | Total/NA  | Water          | 533    | 53035      |
| LCSD 380-53035/24-A  | Lab Control Sample Dup       | Total/NA  | Water          | 533    | 53035      |
| MRL 380-53035/22-A   | Lab Control Sample           | Total/NA  | Water          | 533    | 53035      |
| 380-57432-O-1-D LMS  | Matrix Spike                 | Total/NA  | Water          | 533    | 53035      |
| 380-57432-P-1-C LMSD | Matrix Spike Duplicate       | Total/NA  | Water          | 533    | 53035      |

# Lab Chronicle

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

Job ID: 380-57979-1

## Client Sample ID: HALAWA WELLS UNITS 1 & 2

## Lab Sample ID: 380-57979-1

Date Collected: 08/02/23 10:00

Matrix: Drinking Water

Date Received: 08/08/23 10:00

| Prep Type | Batch Type | Batch Method | Run | Dilution Factor | Batch Number | Analyst | Lab    | Prepared or Analyzed |
|-----------|------------|--------------|-----|-----------------|--------------|---------|--------|----------------------|
| Total/NA  | Prep       | 525.2        |     |                 | 51109        | OTM3    | EA POM | 08/10/23 08:15       |
| Total/NA  | Analysis   | 525.2        |     | 1               | 51305        | Q8LA    | EA POM | 08/11/23 11:56       |
| Total/NA  | Prep       | 533          |     |                 | 53035        | UMV1    | EA POM | 08/24/23 16:20       |
| Total/NA  | Analysis   | 533          |     | 1               | 53385        | UKYM    | EA POM | 08/27/23 06:48       |
| Total/NA  | Prep       | 537.1 DW     |     |                 | 51108        | US1B    | EA POM | 08/10/23 05:33       |
| Total/NA  | Analysis   | 537.1        |     | 1               | 51440        | Y7BM    | EA POM | 08/12/23 11:03       |

## Client Sample ID: FB: HALAWA WELLS UNITS 1 & 2

## Lab Sample ID: 380-57979-3

Date Collected: 08/02/23 10:00

Matrix: Water

Date Received: 08/08/23 10:00

| Prep Type | Batch Type | Batch Method | Run | Dilution Factor | Batch Number | Analyst | Lab    | Prepared or Analyzed |
|-----------|------------|--------------|-----|-----------------|--------------|---------|--------|----------------------|
| Total/NA  | Prep       | 533          |     |                 | 53035        | UMV1    | EA POM | 08/24/23 16:20       |
| Total/NA  | Analysis   | 533          |     | 1               | 53385        | UKYM    | EA POM | 08/27/23 07:07       |
| Total/NA  | Prep       | 537.1 DW     |     |                 | 51108        | US1B    | EA POM | 08/10/23 05:33       |
| Total/NA  | Analysis   | 537.1        |     | 1               | 51333        | Y7BM    | EA POM | 08/11/23 12:14       |

**Laboratory References:**

EA POM = Eurofins Eaton Analytical Pomona, 941 Corporate Center Drive, Pomona, CA 91768-2642, TEL (626)386-1100

# Accreditation/Certification Summary

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

Job ID: 380-57979-1

## Laboratory: Eurofins Eaton Analytical Pomona

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

| Authority | Program | Identification Number | Expiration Date |
|-----------|---------|-----------------------|-----------------|
| Hawaii    | State   | CA00006               | 01-31-24        |

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 | 1-Methylnaphthalene              |
| 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 | 2-Methylnaphthalene              |
| 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 | 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 | 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 | 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           |
| 525.2           | 525.2       | Drinking Water | Isophorone                       |

# Accreditation/Certification Summary

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

Job ID: 380-57979-1

## Laboratory: Eurofins Eaton Analytical Pomona (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        | 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  |
| 533   | 533         | Drinking Water        | 11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS) |
| 533   | 533         | Drinking Water        | 1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)                |
| 533   | 533         | Drinking Water        | 1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)                |
| 533   | 533         | Drinking Water        | 1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)                |
| 533   | 533         | Drinking Water        | 4,8-Dioxa-3H-perfluorononanoic acid (ADONA)                        |
| 533   | 533         | Drinking Water        | 9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid(9Cl-PF3ONS)     |
| 533   | 533         | Drinking Water        | Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)                |
| 533   | 533         | Drinking Water        | Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)                         |
| 533   | 533         | Drinking Water        | Perfluoro (2-ethoxyethane) sulfonic acid (PFEEESA)                 |
| 533   | 533         | Drinking Water        | Perfluoro-3-methoxypropanoic acid (PFMPA)                          |
| 533   | 533         | Drinking Water        | Perfluoro-4-methoxybutanoic acid (PFMBA)                           |
| 533   | 533         | Drinking Water        | Perfluorobutanoic acid (PFBA)                                      |
| 533   | 533         | Drinking Water        | Perfluoroheptanesulfonic acid (PFHpS)                              |
| 533   | 533         | Drinking Water        | Perfluoropentanesulfonic acid (PFPeS)                              |
| 533   | 533         | Drinking Water        | Perfluoropentanoic acid (PFPeA)                                    |
| 533   | 533         | Water                 | 11-Chloroeicosafluoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS) |
| 533   | 533         | Water                 | 1H,1H,2H,2H-Perfluorodecane sulfonic acid (8:2 FTS)                |
| 533   | 533         | Water                 | 1H,1H,2H,2H-Perfluorohexane sulfonic acid (4:2 FTS)                |
| 533   | 533         | Water                 | 1H,1H,2H,2H-Perfluorooctane sulfonic acid (6:2 FTS)                |
| 533   | 533         | Water                 | 4,8-Dioxa-3H-perfluorononanoic acid (ADONA)                        |
| 533   | 533         | Water                 | 9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid(9Cl-PF3ONS)     |

# Accreditation/Certification Summary

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

Job ID: 380-57979-1

## Laboratory: Eurofins Eaton Analytical Pomona (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. |                    |                       |   |
| <u>Analysis Method</u>  | <u>Prep Method</u> | <u>Matrix</u>         | <u>Analyte</u>  |
| 533   | 533                | Water                 | Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)               |
| 533   | 533                | Water                 | Nonafluoro-3,6-dioxaheptanoic acid (NFDHA)                        |
| 533   | 533                | Water                 | Perfluoro (2-ethoxyethane) sulfonic acid (PFEEESA)                |
| 533   | 533                | Water                 | Perfluoro-3-methoxypropanoic acid (PFMPA)                         |
| 533   | 533                | Water                 | Perfluoro-4-methoxybutanoic acid (PFMBA)                          |
| 533   | 533                | Water                 | Perfluorobutanoic acid (PFBA)                                     |
| 533   | 533                | Water                 | Perfluoroheptanesulfonic acid (PFHpS)                             |
| 533   | 533                | Water                 | Perfluoropentanesulfonic acid (PFPeS)                             |
| 533   | 533                | Water                 | Perfluoropentanoic acid (PFPeA)                                   |
| 537.1   | 537.1 DW           | Drinking Water        | 11-Chloroeicosafuoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS) |
| 537.1   | 537.1 DW           | Drinking Water        | 4,8-Dioxa-3H-perfluorononanoic acid (ADONA)                       |
| 537.1   | 537.1 DW           | Drinking Water        | 9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid(9Cl-PF3ONS)    |
| 537.1   | 537.1 DW           | Drinking Water        | Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)               |
| 537.1   | 537.1 DW           | Water                 | 11-Chloroeicosafuoro-3-oxaundecane-1-sulfonic acid (11Cl-PF3OUdS) |
| 537.1   | 537.1 DW           | Water                 | 4,8-Dioxa-3H-perfluorononanoic acid (ADONA)                       |
| 537.1   | 537.1 DW           | Water                 | 9-Chlorohexadecafluoro-3-oxanonane-1-sulfonic acid(9Cl-PF3ONS)    |
| 537.1   | 537.1 DW           | Water                 | Hexafluoropropylene Oxide Dimer Acid (HFPO-DA/GenX)               |

# Method Summary

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

Job ID: 380-57979-1

| Method   | Method Description  | Protocol | Laboratory |
|----------|---|----------|------------|
| 525.2    | Semivolatile Organic Compounds (GC/MS)                                | EPA      | EA POM     |
| 533      | Perfluorinated and Polyfluorinated Alkyl Substances in Drinking Water | EPA      | EA POM     |
| 537.1    | Perfluorinated Alkyl Acids (LC/MS)                                    | EPA      | EA POM     |
| 525.2    | Extraction of Semivolatile Compounds                                  | EPA      | EA POM     |
| 533      | Extraction of Perfluorinated and Polyfluorinated Alkyl Acids          | EPA      | EA POM     |
| 537.1 DW | Extraction of Perfluorinated Alkyl Acids                              | EPA      | EA POM     |

**Protocol References:**

EPA = US Environmental Protection Agency

**Laboratory References:**

EA POM = Eurofins Eaton Analytical Pomona, 941 Corporate Center Drive, Pomona, CA 91768-2642, TEL (626)386-1100





# Sample Summary

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

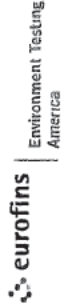
Job ID: 380-57979-1

| Lab Sample ID | Client Sample ID             | Matrix         | Collected      | Received       | PWSID Number |
|---------------|------------------------------|----------------|----------------|----------------|--------------|
| 380-57979-1   | HALAWA WELLS UNITS 1 & 2     | Drinking Water | 08/02/23 10:00 | 08/08/23 10:00 | HI0000331    |
| 380-57979-3   | FB: HALAWA WELLS UNITS 1 & 2 | Water          | 08/02/23 10:00 | 08/08/23 10:00 | HI0000331    |

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

### Chain of Custody Record



|   |  |  |  |   |  |   |  |  |  |
|---|--|--|--|---|--|---|--|--|--|
| <b>Client Information</b><br>Client Contact: <b>Dr. Ron Fenstermacher</b><br>Phone: <b>806-748-5640</b><br>City & County of Honolulu  |  | Lab PM: <b>Arada, Rachelle</b><br>E-Mail: <b>Rachelle.Arada@et.euronisus.com</b>   |  | Carrier Tracking Note:<br>State of Origin:  |  | COC No.: <b>380-27941-2757.2</b><br>Page: <b>Page 2 of 2</b><br>Job #:  |  |  |  |
| Address: <b>630 South Beretania Street, Chemistry Lab Honolulu HI, 96843</b><br>State Zip: <b>HI, 96843</b><br>Phone: <b>808-748-5091 (tel)</b><br>Email: <b>fenstermacher@hbws.org</b><br>Project Name: <b>RED-HILL/HBWS sites Event Desc: RUSH Weekly Red Hill Site</b>   |  | Due Date Requested:<br>TAT Requested (days):<br>Compliance Project: <b>Δ No</b><br>PO #: <b>C20525101 exp 05312023</b><br>W/O #: |  | <b>Analysis Requested</b><br>SUBCONTRACT - 825 PAH Physis LL (EAL) + TCS<br>SUBCONTRACT - 8015 Gas (Purgeable) LL (EAL)<br>SUBCONTRACT - 8915 Diesel LL (EAL) and Motor Oil<br>SUBCONTRACT - (MOD) 525plus PLUS TCS<br>SUBCONTRACT - 8015 Gas (Purgeable) LL (EAL)<br>SUBCONTRACT - 8015 Gas (Purgeable) LL (EAL) |  | Preservation Codes:<br>A - HCL<br>B - NaOH<br>O - AsH2O2<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>Other:<br>M - Hexane<br>N - None<br>O - AsH2O2<br>P - Ni2OAS<br>Q - Na2SO3<br>R - Na2SO4<br>S - H2SO4<br>T - TSP Dodecahydrate<br>U - Acetone<br>V - MCAA<br>W - pH 4-5<br>Y - Trisma<br>Z - other (specify) |  |  |  |
| Sample Identification<br><b>MOANALUA WELLS</b><br><b>AIE-A-GULCH WELLS-PUMP2</b><br><b>AIE-A WELLS-PUMPS 1&amp;2 (260)</b><br><b>HALAWA WELLS UNITS 1&amp;2</b><br><b>FB MOANALUA WELLS</b><br><b>FB-AIE-A-GULCH WELLS-PUMP2</b><br><b>FB-AIE-A WELLS-PUMPS 1&amp;2 (260)</b><br><b>FB HALAWA WELLS UNITS 1&amp;2</b> |  | Sample Date<br><b>5/2/2023</b><br><b>5/2/2023</b><br><b>5/2/2023</b><br><b>5/2/2023</b><br><b>5/2/2023</b><br><b>5/2/2023</b>    |  | Sample Time<br><b>1100</b><br><b>1130</b><br><b>1000</b><br><br><br><br>  |  | Sample Type (C=Comp, G=grab)<br><b>G</b><br><b>G</b><br><b>G</b><br><br><br><br>  |  | Matrix (W=water, S=solid, O=wastoid, BT=tissue, A=air)<br><b>Water</b><br><b>Water</b><br><b>Water</b><br><b>Water</b><br><b>Water</b><br><b>Water</b><br><b>Water</b> |  |
| Field Filtered Sample (Yes or No)<br><input checked="" type="checkbox"/>  |  | Perform MS/MSD (Yes or No)<br><input checked="" type="checkbox"/>  |  | RA Y N<br><b>2 4 2 2</b><br><b>2 4 2 2</b><br><b>2 4 2 2</b><br><br><b>2</b><br><b>2</b><br><b>2</b>  |  | Total Number of Containers<br><input checked="" type="checkbox"/>   |  |  |  |
| Special Instructions/Note:<br><b>Pump 2 10/2</b><br><b>Pump 2 8/4/23</b><br><b>Pump 1</b>   |  | Special Instructions/Note:<br><br><br><br>   |  | Special Instructions/Note:<br><br><br><br>  |  | Special Instructions/Note:<br><br><br><br>  |  |  |  |
| Possible Hazard Identification<br><input type="checkbox"/> Non-Hazard <input type="checkbox"/> Flammable <input type="checkbox"/> Skin Irritant<br>Deliverable Requested: I, II, III, IV, Other (specify)   |  | Empty Kit Relinquished by:<br>Relinquished by: [Redacted]<br>Relinquished by:<br>Relinquished by:                                |  | Sample Disposal (A fee may be assessed if samples are retained longer than 1 month)<br><input type="checkbox"/> Return To Client <input type="checkbox"/> Disposal By Lab <input type="checkbox"/> Archive For _____ Months   |  | Method of Shipment: <b>FED EX 7729 7646 7156</b><br>Date/Time: <b>08/08/2023 10:00</b><br>Company: <b>FEDEX</b>   |  |  |  |
| Date: <b>8/2/23</b><br>Date/Time: <b>1245</b><br>Date/Time:<br>Date/Time:   |  | Date: <b>8/2/23</b><br>Date/Time: <b>1245</b><br>Date/Time:<br>Date/Time:  |  | Date: <b>8/2/23</b><br>Date/Time: <b>1245</b><br>Date/Time:<br>Date/Time:   |  | Date: <b>8/2/23</b><br>Date/Time: <b>1245</b><br>Date/Time:<br>Date/Time:   |  |  |  |
| Custody Seal Intact<br><input type="checkbox"/> Yes <input type="checkbox"/> No   |  | Custody Seal No.:  |  | Cooler Temperature(s) °C and Other Remarks:<br><b>(-52A) 5.1°-0.2°-4.9°</b><br><b>GEL-FROZEN</b>  |  | Ver: 01/16/2019   |  |  |  |

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

**Chain of Custody Record**

**eurofins** Environment Testing  
 America

| <b>Client Information</b><br>Client Contact: Dr. Ron Fenstermacher<br>City & County of Honolulu<br>Address: 630 South Beretania Street, Chemistry Lab<br>Honolulu HI, 96843<br>Phone: 808-748-5840<br>Email: rfenstermacher@hbws.org   |             | Lab PIN: Arada, Rachelle<br>E-Mail: Rachelle.Arada@et.eurofins.com<br>State of Origin:  |  | Carrier Tracking No(s):<br>Page 2 of 2<br>Job #:  |  | ICC No: 380-27941-2757.2  |    |  |  |
|--|-------------|---|--|---|--|---|----|--|--|
| Due Date Requested:<br>TAT Requested (days):<br>Compliance Project: $\Delta$ No<br>PO #: C20525101 exp 05312023<br>WO #:<br>Project #: 38001111<br>SSO#:<br>Project Name: RED-HILL/HBWS sites Event Desc: RUSH Weekly Red Hill<br>Site:  |             | <b>Analysis Requested</b><br>Perform MS/MSD (Yes or No) <input checked="" type="checkbox"/><br>SUBCONTRACT - 625 PAH Physis LL (EAL) + MCS <input type="checkbox"/><br>SUBCONTRACT - 615 Diesel LL (EAL) and Motor Oil <input type="checkbox"/><br>SUBCONTRACT - (MOD) 525plus PLUS TCS <input type="checkbox"/><br>SUBCONTRACT - 615 Gas (Purgeable) LL (EAL) <input type="checkbox"/><br>SUBCONTRACT - 637.1 Full List <input type="checkbox"/><br>SUBCONTRACT - 633 - All Analytes <input type="checkbox"/><br>533 - All Analytes <input type="checkbox"/> |  | Preservation Codes:<br>A - HCL<br>B - NaOH<br>C - AsNaO2<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>Other:<br>M - Hexane<br>N - None<br>O - AsNaO2<br>P - Na2OAS<br>Q - Na2SO3<br>R - Na2SO4<br>S - H2SO4<br>T - TSP Dodecahydrate<br>U - Acetone<br>V - MCAA<br>W - PH 4.5<br>Y - Triana<br>Z - other (specify) |  | Total Number of Containers: <input checked="" type="checkbox"/> |    | Special Instructions/Note:<br>Pump 2 at 8/4/23<br>Pump 2<br>Pump 1 |  |
| Sample Identification  | Sample Date | Sample Time   | Sample Type (C=comp, G=grab)   | Preservation Code:  | Matrix (W=water, S=solid, O=oil, G=Gas, A=Air) | RA  | RY | YN   |  |
| MOANALUA WELLS   |             |   |  |   | Water  |   |    |  |  |
| AHEA-GULCH WELLS-PUMP2   | 8/1/2023    | 11:00   | G  |   | Water  | 3   | 3  |  |  |
| AHEA WELLS PUMPS 1&2 (260)   | 8/1/2023    | 11:30   | G  |   | Water  | 3   | 3  |  |  |
| HALAWA WELLS UNITS 1&2   | 8/2/2023    | 10:00   | G  |   | Water  | 3   | 3  |  |  |
| FB MOANALUA WELLS  |             |   |  |   | Water  |   |    |  |  |
| FB-AHEA GULCH WELLS PUMP2  | 8/2/2023    |   |  |   | Water  |   |    |  |  |
| FB-AHEA WELLS PUMPS 1&2 (260)  | 8/2/2023    |   |  |   | Water  |   |    |  |  |
| FB HALAWA WELLS UNITS 1&2  | 8/2/2023    |   |  |   | Water  |   |    |  |  |
| <b>Possible Hazard Identification</b><br><input type="checkbox"/> Non-Hazard <input type="checkbox"/> Flammable <input type="checkbox"/> Skin Irritant <input type="checkbox"/> Poison B <input type="checkbox"/> Unknown <input type="checkbox"/> Radiological<br><b>Deliverable Requested:</b> I, II, III, IV, Other (specify) |             |   |  |   |  |   |    |  |  |
| <b>Empty Kit Relinquished by:</b><br>Relinquished by: <i>Bryan Tsubota</i><br>Date/Time: 8/2/2023 12:45<br>Company: HBWS   |             |   | <b>Sample Disposal (A fee may be assessed if samples are retained longer than 1 month)</b><br><input type="checkbox"/> Return To Client <input type="checkbox"/> Disposal By Lab <input type="checkbox"/> Archive For _____ Months<br><b>Special Instructions/QC Requirements:</b> |   |  |   |    |  |  |
| <b>Relinquished by:</b><br>Relinquished by: <i>Bryan Tsubota</i><br>Date/Time: 8/2/2023 10:00<br>Company: Eurofins   |             |   | <b>Method of Shipment:</b> FED EX 7729 7646 7156<br>Date/Time: 08/08/2023 10:00<br>Company: Eurofins   |   |  |   |    |  |  |
| <b>Relinquished by:</b><br>Relinquished by: _____<br>Date/Time: _____<br>Company: _____  |             |   | <b>Method of Shipment:</b> _____<br>Date/Time: _____<br>Company: _____   |   |  |   |    |  |  |
| <b>Custody Seals Intact:</b><br>$\Delta$ Yes $\Delta$ No   |             |   | <b>Cooler Temperature(s) °C and Other Remarks:</b><br>752A 5.1 °C 2:49<br>GEL-FROZEN   |   |  |   |    |  |  |

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

**Chain of Custody Record**



|   |  |  |  |   |  |  |  |  |  |
|---|--|--|--|---|--|--|--|--|--|
| <b>Client Information</b><br>Client Contact: <b>Dr. Ron Fenstermacher</b><br>Phone: <b>806-748-5640</b><br>City & County of Honolulu  |  | Lab PM: <b>Arada, Rachelle</b><br>E-Mail: <b>Rachelle.Arada@et.euronisus.com</b><br>PWSID:                                       |  | Carrier Tracking Note:<br>State of Origin:  |  | COC No.: <b>380-27941-2757.2</b><br>Page: <b>Page 2 of 2</b><br>Job #:   |  |  |  |
| Address: <b>630 South Beretania Street, Chemistry Lab Honolulu HI, 96843</b><br>State Zip: <b>HI, 96843</b><br>Phone: <b>808-748-5091 (tel)</b><br>Email: <b>rfenstermacher@hbws.org</b><br>Project Name: <b>RED-HILL/HBWS sites Event Desc: RUSH Weekly Red Hill Site</b>  |  | Due Date Requested:<br>TAT Requested (days):<br>Compliance Project: <b>Δ No</b><br>PO #: <b>C20525101 exp 05312023</b><br>W/O #: |  | <b>Analysis Requested</b><br>SUBCONTRACT - 825 PAH Physis LL (EAL) + TCS<br>SUBCONTRACT - 8015 Gas (Purgeable) LL (EAL)<br>SUBCONTRACT - 8915 Diesel LL (EAL) and Motor Oil<br>SUBCONTRACT - (MOD) 525plus PLUS TCS<br>SUBCONTRACT - 8015 Gas (Purgeable) LL (EAL)<br>SUBCONTRACT - 8015 Gas (Purgeable) LL (EAL) |  | Preservation Codes:<br>A - HCL<br>B - NaOH<br>O - Ash/O2<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>Other:<br>M - Hexane<br>N - None<br>O - Ash/O2<br>P - Ni2OAS<br>Q - Na2SO3<br>R - Na2S2O3<br>S - H2SO4<br>T - TSP Dodecahydrate<br>U - Acetone<br>V - MCAA<br>W - pH 4-5<br>Y - Trisma<br>Z - other (specify) |  |  |  |
| Sample Identification<br><b>MOANALUA WELLS</b><br><b>AIE-A GULCH WELLS PUMP2</b><br><b>AIE-A WELLS PUMPS 1&amp;2 (260)</b><br><b>HALAWA WELLS UNITS 1&amp;2</b><br><b>FB MOANALUA WELLS</b><br><b>FB AIE-A GULCH WELLS PUMP2</b><br><b>FB AIE-A WELLS PUMPS 1&amp;2 (260)</b><br><b>FB HALAWA WELLS UNITS 1&amp;2</b> |  | Sample Date<br><b>5/2/2023</b><br><b>5/2/2023</b><br><b>5/2/2023</b><br><b>5/2/2023</b><br><b>5/2/2023</b><br><b>5/2/2023</b>    |  | Sample Time<br><b>1100</b><br><b>1130</b><br><b>1000</b><br><br><br><br>  |  | Sample Type (C=Comp, G=grab)<br><b>G</b><br><b>G</b><br><b>G</b><br><br><br><br>   |  | Matrix (W=water, S=solid, O=wastoid, BT=tissue, A=air)<br><b>Water</b><br><b>Water</b><br><b>Water</b><br><b>Water</b><br><b>Water</b><br><b>Water</b><br><b>Water</b> |  |
| Field Filtered Sample (Yes or No)<br><input checked="" type="checkbox"/>  |  | Perform MS/MSD (Yes or No)<br><input checked="" type="checkbox"/>  |  | RA Y N<br><b>2 4 2 2</b><br><b>2 4 2 2</b><br><b>2 4 2 2</b><br><br><b>2</b><br><b>2</b><br><b>2</b>  |  | Total Number of Containers<br><input checked="" type="checkbox"/>  |  |  |  |
| Special Instructions/Note:<br><b>Pump 2 100</b><br><b>Pump 2 8/4/23</b><br><b>Pump 1</b>  |  | Special Instructions/Note:<br><br><br>   |  | Special Instructions/Note:<br><br><br>  |  | Special Instructions/Note:<br><br><br>   |  |  |  |
| Possible Hazard Identification<br><input type="checkbox"/> Non-Hazard <input type="checkbox"/> Flammable <input type="checkbox"/> Skin Irritant <input type="checkbox"/> Poison B <input type="checkbox"/> Unknown <input type="checkbox"/> Radiological  |  | Deliverable Requested: I, II, III, IV, Other (specify)   |  | Sample Disposal (A fee may be assessed if samples are retained longer than 1 month)<br><input type="checkbox"/> Return To Client <input type="checkbox"/> Disposal By Lab <input type="checkbox"/> Archive For _____ Months   |  | Method of Shipment: <b>FED EX 7729 7646 7156</b><br>Date/Time: _____<br>Company: _____   |  |  |  |
| Empty Kit Relinquished by: _____<br>Date/Time: <b>8/2/23</b>  |  | Relinquished by: _____<br>Date/Time: <b>1245</b>   |  | Relinquished by: _____<br>Date/Time: _____  |  | Relinquished by: _____<br>Date/Time: _____   |  |  |  |
| Custody Seals Intact<br><input type="checkbox"/> Yes <input type="checkbox"/> No  |  | Custody Seal No.: _____  |  | Cooler Temperature(s) °C and Other Remarks:<br><b>(752A) 5.1°-0.2°-4.9°</b><br><b>GEL FROZEN</b>  |  | Ver: 01/16/2019  |  |  |  |

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

**Chain of Custody Record**



Environment Testing  
 America

|  |  |   |  |  |  |   |  |
|--|--|---|--|--|--|---|--|
| <b>Client Information</b>  |  | Lab PIN:<br>Arada, Rachelle               |  | Carrier Tracking No(s)                           |  | ICC No:<br>380-27941-2757.2   |  |
| Client Contact:<br>Dr. Ron Fenstermacher   |  | E-Mail:<br>Rachelle.Arada@et.eurofins.com |  | State of Origin:                                 |  | Page<br>Page 2 of 2   |  |
| Company:<br>City & County of Honolulu  |  | PWSID:                                    |  | Job #:   |  | Preservation Codes:<br>M - Hexane<br>N - None<br>O - AsNaO2<br>P - Na2OAS<br>Q - Na2SO3<br>R - NaHSO4<br>S - H2SO4<br>T - TSP Dodecahydrate<br>U - Acetone<br>V - MCAA<br>W - PH 4.5<br>Y - Trizma<br>Z - other (specify)<br>Other: |  |
| Address:<br>630 South Beretania Street, Chemistry Lab<br>Honolulu<br>State, Zip:<br>HI, 96843  |  | Due Date Requested:                       |  | Analysis Requested                               |  | Total Number of Containers  |  |
| City   |  | TAT Requested (days):                     |  | 533 - All Analytes                               |  | Special Instructions/Note:  |  |
| State, Zip:  |  | Compliance Project: Δ No                  |  | 537.1_QW_PREC - 537.1 Full List                  |  | Pump 2 at 8/4/23  |  |
| HI, 96843  |  | PO #                                      |  | SUBCONTRACT - 8015 Gas (Purgeable) LL (EAL)      |  | Pump 2  |  |
| Phone:   |  | C20525101 exp 05312023                    |  | 525.2_PREC - (MOD) 525plus PLUS TCS              |  | Pump 1  |  |
| Email:   |  | WO #                                      |  | SUBCONTRACT - 8015 Diesel LL (EAL) and Motor Oil |  |   |  |
| fenstermacher@hbws.org   |  | Project #                                 |  | SUBCONTRACT - 8015 Gas (Purgeable) LL (EAL)      |  |   |  |
| RED-HILL/HBWS sites Event Desc: RUSH Weekly Red Hill   |  | 38001111                                  |  | SUBCONTRACT - 8015 PAH Physic LL (EAL) + TCS     |  |   |  |
| Site:  |  | SSOW#                                     |  | Perform MS/MSD (Yes or No)                       |  |   |  |
| Sample Identification  |  | Sample Date                               |  | Plaid Filtered Sample (Yes or No)                |  |   |  |
| MOANALUA WELLS   |  | 8/1/2023                                  |  | X  |  |   |  |
| AHEA-GULCH WELLS PUMP2   |  | 8/1/2023 11:00                            |  | RA Y N   |  |   |  |
| AHEA WELLS PUMPS 1&2 (260)   |  | 8/1/2023 11:30                            |  | 3 3  |  |   |  |
| HALAWA WELLS UNITS 1&2   |  | 8/2/2023 10:00                            |  | 3 3  |  |   |  |
| FB MOANALUA WELLS  |  |   |  |  |  |   |  |
| FB-AHEA GULCH WELLS PUMP2  |  | 8/2/2023                                  |  | 1 1  |  |   |  |
| FB-AHEA WELLS PUMPS 1&2 (260)  |  | 8/2/2023                                  |  | 1 1  |  |   |  |
| FB HALAWA WELLS UNITS 1&2  |  | 8/2/2023                                  |  | 1 1  |  |   |  |
| Possible Hazard Identification   |  | Sample Time                               |  | Matrix   |  |   |  |
| <input type="checkbox"/> Non-Hazard <input type="checkbox"/> Flammable <input type="checkbox"/> Skin Irritant <input type="checkbox"/> Poison B <input type="checkbox"/> Unknown <input type="checkbox"/> Radiological |  | Preservation Code:                        |  | Water  |  |   |  |
| Deliverable Requested: I, II, III, IV, Other (specify)   |  | Sample Type (C=comp, G=grab)              |  | Water  |  |   |  |
| Empty Kit Relinquished by:   |  | Date                                      |  | Date   |  | Date  |  |
| Relinquished by: <i>Byron Inub</i>   |  | 8/2/2023 12:45                            |  | Company HBWS                                     |  | Company   |  |
| Relinquished by:   |  | Date/Time:                                |  | Company  |  | Company   |  |
| Relinquished by:   |  | Date/Time:                                |  | Company  |  | Company   |  |
| Custody Seals Intact:  |  | Custody Seal No.:                         |  | Cooler Temperature(s) °C and Other Remarks:      |  | Cooler Temperature(s) °C and Other Remarks:   |  |
| Δ Yes Δ No   |  |   |  | 752A 5.1'-0.2'-49°                               |  | GEL-FASTEN  |  |

Sample Disposal (A fee may be assessed if samples are retained longer than 1 month)  
 Return To Client  Disposal By Lab  Archive For \_\_\_\_\_ Months

Method of Shipment: *Fed Ex 7729 7646 7156*  
 Date/Time: *08/08/2023 10:00* Company  
 Received by: *GA GREINER* Company  
 Date/Time: \_\_\_\_\_ Company  
 Date/Time: \_\_\_\_\_ Company



# Login Sample Receipt Checklist

Client: City & County of Honolulu

Job Number: 380-57979-1

**Login Number: 57979**  
**List Number: 1**  
**Creator: Elyas, Matthew**

**List Source: Eurofins Eaton Analytical Pomona**

| 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.          | True   |         |
| 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   |         |

