

## Report of Analysis

|                    |                       |                  |                 |                |
|--------------------|-----------------------|------------------|-----------------|----------------|
| Client:            | Tetra Tech NUS, Inc.  |                  | Date Collected: | 10/30/24       |
| Project:           | CTO WE13              |                  | Date Received:  | 10/31/24       |
| Client Sample ID:  | RW5-SP201-90-20241030 |                  | SDG No.:        | P4652          |
| Lab Sample ID:     | P4652-05              |                  | Matrix:         | Water          |
| Analytical Method: | SW8270SIM             |                  | % Solid:        | 0              |
| Sample Wt/Vol:     | 960                   | Units: mL        | Final Vol:      | 1000 uL        |
| Soil Aliquot Vol:  |                       | uL               | Test:           | SVOC-SIMGroup1 |
| Extraction Type :  |                       | Decanted : N     | Level :         | LOW            |
| Injection Volume : |                       | GPC Factor : 1.0 | GPC Cleanup :   | N PH :         |
| Prep Method :      | SW3510C               |                  |                 |                |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BN034843.D        | 1         | 11/01/24 09:35 | 11/04/24 17:19 | PB164594      |

| CAS Number                | Parameter               | Conc. | Qualifier | MDL      | LOD  | LOQ / CRQL | Units    |
|---------------------------|-------------------------|-------|-----------|----------|------|------------|----------|
| <b>TARGETS</b>            |                         |       |           |          |      |            |          |
| 123-91-1                  | 1,4-Dioxane             | 0.21  | U         | 0.070    | 0.21 | 0.21       | ug/L     |
| <b>SURROGATES</b>         |                         |       |           |          |      |            |          |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.34  |           | 30 - 150 |      | 86%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.42  |           | 30 - 150 |      | 104%       | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.39  |           | 55 - 111 |      | 97%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.39  |           | 53 - 106 |      | 97%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.53  | *         | 58 - 132 |      | 133%       | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 5980  | 7.589     |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 18200 | 10.361    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 9000  | 14.222    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 17900 | 16.97     |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 9270  | 21.16     |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 6570  | 23.341    |          |      |            |          |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products