

## LAB CHRONICLE

|                 |                      |                   |                               |
|-----------------|----------------------|-------------------|-------------------------------|
| <b>OrderID:</b> | Q2050                | <b>OrderDate:</b> | 5/15/2025 9:46:00 AM          |
| <b>Client:</b>  | Tetra Tech NUS, Inc. | <b>Project:</b>   | NWIRP Bethpage 112G08005-WE13 |
| <b>Contact:</b> | Ernie Wu             | <b>Location:</b>  | L31,VOA Ref. #3 Water         |

| LabID           | ClientID                            | Matrix       | Test         | Method   | Sample Date     | Prep Date | Anal Date | Received        |
|-----------------|-------------------------------------|--------------|--------------|----------|-----------------|-----------|-----------|-----------------|
| <b>Q2050-01</b> | <b>BP-VPB-182-TB-2025<br/>0512</b>  | <b>Water</b> |              |          | <b>05/12/25</b> |           |           | <b>05/14/25</b> |
|                 |                                     |              | VOCMS Group1 | 8260-Low |                 |           | 05/15/25  |                 |
| <b>Q2050-02</b> | <b>BP-VPB-182-GW-60-6<br/>2</b>     | <b>Water</b> |              |          | <b>05/12/25</b> |           |           | <b>05/14/25</b> |
|                 |                                     |              | VOCMS Group1 | 8260-Low |                 |           | 05/15/25  |                 |
| <b>Q2050-03</b> | <b>BP-VPB-182-GW-100-<br/>102</b>   | <b>Water</b> |              |          | <b>05/12/25</b> |           |           | <b>05/14/25</b> |
|                 |                                     |              | VOCMS Group1 | 8260-Low |                 |           | 05/15/25  |                 |
| <b>Q2050-04</b> | <b>BP-VPB-182-GW-150-<br/>152</b>   | <b>Water</b> |              |          | <b>05/13/25</b> |           |           | <b>05/14/25</b> |
|                 |                                     |              | VOCMS Group1 | 8260-Low |                 |           | 05/15/25  |                 |
| <b>Q2050-05</b> | <b>BP-VPB-182-GW-205-<br/>207</b>   | <b>Water</b> |              |          | <b>05/14/25</b> |           |           | <b>05/14/25</b> |
|                 |                                     |              | VOCMS Group1 | 8260-Low |                 |           | 05/15/25  |                 |
| <b>Q2050-06</b> | <b>BP-VPB-182-GW-220-<br/>222</b>   | <b>Water</b> |              |          | <b>05/14/25</b> |           |           | <b>05/14/25</b> |
|                 |                                     |              | VOCMS Group1 | 8260-Low |                 |           | 05/15/25  |                 |
| <b>Q2050-07</b> | <b>BP-VPB-182-DUP-202<br/>50514</b> | <b>Water</b> |              |          | <b>05/14/25</b> |           |           | <b>05/14/25</b> |
|                 |                                     |              | VOCMS Group1 | 8260-Low |                 |           | 05/15/25  |                 |
| <b>Q2050-08</b> | <b>BP-VPB-182-EB-2025<br/>0514</b>  | <b>Water</b> |              |          | <b>05/14/25</b> |           |           | <b>05/14/25</b> |
|                 |                                     |              | VOCMS Group1 | 8260-Low |                 |           | 05/15/25  |                 |
| <b>Q2050-09</b> | <b>BP-VPB-182-GW-240-<br/>242</b>   | <b>Water</b> |              |          | <b>05/14/25</b> |           |           | <b>05/14/25</b> |
|                 |                                     |              | VOCMS Group1 | 8260-Low |                 |           | 05/15/25  |                 |

**Hit Summary Sheet**  
**SW-846**

**SDG No.:** Q2050  
**Client:** Tetra Tech NUS, Inc.

| Sample ID                     | Client ID                                               | Matrix                      | Parameter | Concentration | C | MDL  | LOD  | RDL  | Units |
|-------------------------------|---------------------------------------------------------|-----------------------------|-----------|---------------|---|------|------|------|-------|
| <b>Client ID:</b><br>Q2050-01 | <b>BP-VPB-182-TB-20250512</b><br>BP-VPB-182-TB-20 Water | Acetone                     |           | 2.50          | J | 1.50 | 3.80 | 5.00 | ug/L  |
|                               |                                                         | <b>Total Voc :</b>          |           | 2.50          |   |      |      |      |       |
|                               |                                                         | <b>Total Concentration:</b> |           | 2.50          |   |      |      |      |       |
| <b>Client ID:</b><br>Q2050-02 | <b>BP-VPB-182-GW-60-62</b><br>BP-VPB-182-GW-6 Water     | Acetone                     |           | 7.80          |   | 1.50 | 3.80 | 5.00 | ug/L  |
| Q2050-02                      | BP-VPB-182-GW-6 Water                                   | Chloroform                  |           | 1.10          |   | 0.25 | 0.50 | 1.00 | ug/L  |
| Q2050-02                      | BP-VPB-182-GW-6 Water                                   | Tetrachloroethene           |           | 1.30          |   | 0.23 | 0.50 | 1.00 | ug/L  |
|                               |                                                         | <b>Total Voc :</b>          |           | 10.2          |   |      |      |      |       |
|                               |                                                         | <b>Total Concentration:</b> |           | 10.2          |   |      |      |      |       |
| <b>Client ID:</b><br>Q2050-03 | <b>BP-VPB-182-GW-100-102</b><br>BP-VPB-182-GW-1 Water   | Acetone                     |           | 3.80          | J | 1.50 | 3.80 | 5.00 | ug/L  |
| Q2050-03                      | BP-VPB-182-GW-1 Water                                   | Chloroform                  |           | 0.93          | J | 0.25 | 0.50 | 1.00 | ug/L  |
| Q2050-03                      | BP-VPB-182-GW-1 Water                                   | Tetrachloroethene           |           | 1.10          |   | 0.23 | 0.50 | 1.00 | ug/L  |
|                               |                                                         | <b>Total Voc :</b>          |           | 5.83          |   |      |      |      |       |
|                               |                                                         | <b>Total Concentration:</b> |           | 5.83          |   |      |      |      |       |
| <b>Client ID:</b><br>Q2050-04 | <b>BP-VPB-182-GW-150-152</b><br>BP-VPB-182-GW-1 Water   | Acetone                     |           | 4.00          | J | 1.50 | 3.80 | 5.00 | ug/L  |
|                               |                                                         | <b>Total Voc :</b>          |           | 4.00          |   |      |      |      |       |
|                               |                                                         | <b>Total Concentration:</b> |           | 4.00          |   |      |      |      |       |
| <b>Client ID:</b><br>Q2050-05 | <b>BP-VPB-182-GW-205-207</b><br>BP-VPB-182-GW-2 Water   | Acetone                     |           | 5.20          |   | 1.50 | 3.80 | 5.00 | ug/L  |
|                               |                                                         | <b>Total Voc :</b>          |           | 5.20          |   |      |      |      |       |
|                               |                                                         | <b>Total Concentration:</b> |           | 5.20          |   |      |      |      |       |
| <b>Client ID:</b><br>Q2050-06 | <b>BP-VPB-182-GW-220-222</b><br>BP-VPB-182-GW-2 Water   | Acetone                     |           | 3.20          | J | 1.50 | 3.80 | 5.00 | ug/L  |
|                               |                                                         | <b>Total Voc :</b>          |           | 3.20          |   |      |      |      |       |
|                               |                                                         | <b>Total Concentration:</b> |           | 3.20          |   |      |      |      |       |
| <b>Client ID:</b><br>Q2050-07 | <b>BP-VPB-182-DUP-20250514</b><br>BP-VPB-182-DUP- Water | Acetone                     |           | 2.80          | J | 1.50 | 3.80 | 5.00 | ug/L  |
|                               |                                                         | <b>Total Voc :</b>          |           | 2.80          |   |      |      |      |       |
|                               |                                                         | <b>Total Concentration:</b> |           | 2.80          |   |      |      |      |       |
| <b>Client ID:</b><br>Q2050-08 | <b>BP-VPB-182-EB-20250514</b><br>BP-VPB-182-EB-20 Water | Acetone                     |           | 4.30          | J | 1.50 | 3.80 | 5.00 | ug/L  |
|                               |                                                         | <b>Total Voc :</b>          |           | 4.30          |   |      |      |      |       |
|                               |                                                         | <b>Total Concentration:</b> |           | 4.30          |   |      |      |      |       |
| <b>Client ID:</b><br>Q2050-09 | <b>BP-VPB-182-GW-240-242</b><br>BP-VPB-182-GW-2 Water   | Acetone                     |           | 2.90          | J | 1.50 | 3.80 | 5.00 | ug/L  |
|                               |                                                         | <b>Total Voc :</b>          |           | 2.90          |   |      |      |      |       |
|                               |                                                         | <b>Total Concentration:</b> |           | 2.90          |   |      |      |      |       |



# SAMPLE DATA

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/12/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-TB-20250512        |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-01                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046209.D        | 1         |           | 05/15/25 13:24 | VX051525      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOD  | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |      |            |       |
| 74-87-3        | Chloromethane                  | 0.50  | U         | 0.32  | 0.50 | 1.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 0.75  | U         | 0.26  | 0.75 | 1.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 3.80  | U         | 1.40  | 3.80 | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 0.75  | U         | 0.47  | 0.75 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 0.50  | U         | 0.33  | 0.50 | 1.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 0.75  | U         | 0.23  | 0.75 | 1.00       | ug/L  |
| 67-64-1        | Acetone                        | 2.50  | J         | 1.50  | 3.80 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 0.75  | U         | 0.21  | 0.75 | 1.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 0.50  | U         | 0.28  | 0.50 | 1.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 2.50  | U         | 0.98  | 2.50 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 0.75  | U         | 0.19  | 0.75 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                     | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 71-43-2        | Benzene                        | 0.50  | U         | 0.15  | 0.50 | 1.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 0.75  | U         | 0.090 | 0.75 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 2.50  | U         | 0.68  | 2.50 | 5.00       | ug/L  |
| 108-88-3       | Toluene                        | 0.50  | U         | 0.14  | 0.50 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene          | 0.50  | U         | 0.17  | 0.50 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 79-00-5        | 1,1,2-Trichloroethane          | 0.50  | U         | 0.21  | 0.50 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                     | 2.50  | U         | 0.89  | 2.50 | 5.00       | ug/L  |

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/12/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-TB-20250512        |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-01                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046209.D        | 1         |           | 05/15/25 13:24 | VX051525      |

| CAS Number                            | Parameter                 | Conc.  | Qualifier | MDL      | LOD  | LOQ / CRQL | Units   |
|---------------------------------------|---------------------------|--------|-----------|----------|------|------------|---------|
| 124-48-1                              | Dibromochloromethane      | 0.50   | U         | 0.18     | 0.50 | 1.00       | ug/L    |
| 127-18-4                              | Tetrachloroethene         | 0.50   | U         | 0.23     | 0.50 | 1.00       | ug/L    |
| 108-90-7                              | Chlorobenzene             | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-41-4                              | Ethyl Benzene             | 0.50   | U         | 0.13     | 0.50 | 1.00       | ug/L    |
| 179601-23-1                           | m/p-Xylenes               | 1.00   | U         | 0.24     | 1.00 | 2.00       | ug/L    |
| 95-47-6                               | o-Xylene                  | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-42-5                              | Styrene                   | 0.50   | U         | 0.15     | 0.50 | 1.00       | ug/L    |
| 75-25-2                               | Bromoform                 | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 98-82-8                               | Isopropylbenzene          | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 79-34-5                               | 1,1,2,2-Tetrachloroethane | 0.50   | U         | 0.26     | 0.50 | 1.00       | ug/L    |
| 541-73-1                              | 1,3-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| 106-46-7                              | 1,4-Dichlorobenzene       | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 95-50-1                               | 1,2-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| <b>SURROGATES</b>                     |                           |        |           |          |      |            |         |
| 17060-07-0                            | 1,2-Dichloroethane-d4     | 55.1   |           | 81 - 118 |      | 110%       | SPK: 50 |
| 1868-53-7                             | Dibromofluoromethane      | 52.2   |           | 80 - 119 |      | 104%       | SPK: 50 |
| 2037-26-5                             | Toluene-d8                | 51.0   |           | 89 - 112 |      | 102%       | SPK: 50 |
| 460-00-4                              | 4-Bromofluorobenzene      | 50.7   |           | 85 - 114 |      | 101%       | SPK: 50 |
| <b>INTERNAL STANDARDS</b>             |                           |        |           |          |      |            |         |
| 363-72-4                              | Pentafluorobenzene        | 62700  | 5.544     |          |      |            |         |
| 540-36-3                              | 1,4-Difluorobenzene       | 125000 | 6.757     |          |      |            |         |
| 3114-55-4                             | Chlorobenzene-d5          | 120000 | 10.049    |          |      |            |         |
| 3855-82-1                             | 1,4-Dichlorobenzene-d4    | 51100  | 12.018    |          |      |            |         |
| <b>TENTATIVE IDENTIFIED COMPOUNDS</b> |                           |        |           |          |      |            |         |
| 75-43-4                               | Dichlorofluoromethane     | N.D    |           |          |      |            |         |

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/12/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-TB-20250512        |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-01                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046209.D        | 1         |           | 05/15/25 13:24 | VX051525      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOD | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|-----|------------|-------|
|------------|-----------|-------|-----------|-----|-----|------------|-------|

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

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/12/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-60-62           |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-02                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046208.D        | 1         |           | 05/15/25 13:01 | VX051525      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOD  | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |      |            |       |
| 74-87-3        | Chloromethane                  | 0.50  | U         | 0.32  | 0.50 | 1.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 0.75  | U         | 0.26  | 0.75 | 1.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 3.80  | U         | 1.40  | 3.80 | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 0.75  | U         | 0.47  | 0.75 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 0.50  | U         | 0.33  | 0.50 | 1.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 0.75  | U         | 0.23  | 0.75 | 1.00       | ug/L  |
| 67-64-1        | Acetone                        | 7.80  |           | 1.50  | 3.80 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 0.75  | U         | 0.21  | 0.75 | 1.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 0.50  | U         | 0.28  | 0.50 | 1.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 2.50  | U         | 0.98  | 2.50 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 0.75  | U         | 0.19  | 0.75 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                     | 1.10  |           | 0.25  | 0.50 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 71-43-2        | Benzene                        | 0.50  | U         | 0.15  | 0.50 | 1.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 0.75  | U         | 0.090 | 0.75 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 2.50  | U         | 0.68  | 2.50 | 5.00       | ug/L  |
| 108-88-3       | Toluene                        | 0.50  | U         | 0.14  | 0.50 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene          | 0.50  | U         | 0.17  | 0.50 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 79-00-5        | 1,1,2-Trichloroethane          | 0.50  | U         | 0.21  | 0.50 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                     | 2.50  | U         | 0.89  | 2.50 | 5.00       | ug/L  |

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/12/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-60-62           |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-02                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046208.D        | 1         |           | 05/15/25 13:01 | VX051525      |

| CAS Number                            | Parameter                 | Conc.  | Qualifier | MDL      | LOD  | LOQ / CRQL | Units   |
|---------------------------------------|---------------------------|--------|-----------|----------|------|------------|---------|
| 124-48-1                              | Dibromochloromethane      | 0.50   | U         | 0.18     | 0.50 | 1.00       | ug/L    |
| 127-18-4                              | Tetrachloroethene         | 1.30   |           | 0.23     | 0.50 | 1.00       | ug/L    |
| 108-90-7                              | Chlorobenzene             | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-41-4                              | Ethyl Benzene             | 0.50   | U         | 0.13     | 0.50 | 1.00       | ug/L    |
| 179601-23-1                           | m/p-Xylenes               | 1.00   | U         | 0.24     | 1.00 | 2.00       | ug/L    |
| 95-47-6                               | o-Xylene                  | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-42-5                              | Styrene                   | 0.50   | U         | 0.15     | 0.50 | 1.00       | ug/L    |
| 75-25-2                               | Bromoform                 | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 98-82-8                               | Isopropylbenzene          | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 79-34-5                               | 1,1,2,2-Tetrachloroethane | 0.50   | U         | 0.26     | 0.50 | 1.00       | ug/L    |
| 541-73-1                              | 1,3-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| 106-46-7                              | 1,4-Dichlorobenzene       | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 95-50-1                               | 1,2-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| <b>SURROGATES</b>                     |                           |        |           |          |      |            |         |
| 17060-07-0                            | 1,2-Dichloroethane-d4     | 54.7   |           | 81 - 118 |      | 109%       | SPK: 50 |
| 1868-53-7                             | Dibromofluoromethane      | 51.3   |           | 80 - 119 |      | 103%       | SPK: 50 |
| 2037-26-5                             | Toluene-d8                | 49.8   |           | 89 - 112 |      | 100%       | SPK: 50 |
| 460-00-4                              | 4-Bromofluorobenzene      | 50.0   |           | 85 - 114 |      | 100%       | SPK: 50 |
| <b>INTERNAL STANDARDS</b>             |                           |        |           |          |      |            |         |
| 363-72-4                              | Pentafluorobenzene        | 61900  | 5.55      |          |      |            |         |
| 540-36-3                              | 1,4-Difluorobenzene       | 123000 | 6.757     |          |      |            |         |
| 3114-55-4                             | Chlorobenzene-d5          | 115000 | 10.055    |          |      |            |         |
| 3855-82-1                             | 1,4-Dichlorobenzene-d4    | 50500  | 12.018    |          |      |            |         |
| <b>TENTATIVE IDENTIFIED COMPOUNDS</b> |                           |        |           |          |      |            |         |
| 75-43-4                               | Dichlorofluoromethane     | N.D    |           |          |      |            |         |



## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/12/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-60-62           |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-02                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046208.D        | 1         |           | 05/15/25 13:01 | VX051525      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOD | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|-----|------------|-------|
|------------|-----------|-------|-----------|-----|-----|------------|-------|

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

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/12/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-100-102         |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-03                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046211.D        | 1         |           | 05/15/25 14:11 | VX051525      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOD  | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |      |            |       |
| 74-87-3        | Chloromethane                  | 0.50  | U         | 0.32  | 0.50 | 1.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 0.75  | U         | 0.26  | 0.75 | 1.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 3.80  | U         | 1.40  | 3.80 | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 0.75  | U         | 0.47  | 0.75 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 0.50  | U         | 0.33  | 0.50 | 1.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 0.75  | U         | 0.23  | 0.75 | 1.00       | ug/L  |
| 67-64-1        | Acetone                        | 3.80  | J         | 1.50  | 3.80 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 0.75  | U         | 0.21  | 0.75 | 1.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 0.50  | U         | 0.28  | 0.50 | 1.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 2.50  | U         | 0.98  | 2.50 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 0.75  | U         | 0.19  | 0.75 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                     | 0.93  | J         | 0.25  | 0.50 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 71-43-2        | Benzene                        | 0.50  | U         | 0.15  | 0.50 | 1.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 0.75  | U         | 0.090 | 0.75 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 2.50  | U         | 0.68  | 2.50 | 5.00       | ug/L  |
| 108-88-3       | Toluene                        | 0.50  | U         | 0.14  | 0.50 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene          | 0.50  | U         | 0.17  | 0.50 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 79-00-5        | 1,1,2-Trichloroethane          | 0.50  | U         | 0.21  | 0.50 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                     | 2.50  | U         | 0.89  | 2.50 | 5.00       | ug/L  |

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/12/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-100-102         |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-03                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046211.D        | 1         |           | 05/15/25 14:11 | VX051525      |

| CAS Number                            | Parameter                 | Conc.  | Qualifier | MDL      | LOD  | LOQ / CRQL | Units   |
|---------------------------------------|---------------------------|--------|-----------|----------|------|------------|---------|
| 124-48-1                              | Dibromochloromethane      | 0.50   | U         | 0.18     | 0.50 | 1.00       | ug/L    |
| 127-18-4                              | Tetrachloroethene         | 1.10   |           | 0.23     | 0.50 | 1.00       | ug/L    |
| 108-90-7                              | Chlorobenzene             | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-41-4                              | Ethyl Benzene             | 0.50   | U         | 0.13     | 0.50 | 1.00       | ug/L    |
| 179601-23-1                           | m/p-Xylenes               | 1.00   | U         | 0.24     | 1.00 | 2.00       | ug/L    |
| 95-47-6                               | o-Xylene                  | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-42-5                              | Styrene                   | 0.50   | U         | 0.15     | 0.50 | 1.00       | ug/L    |
| 75-25-2                               | Bromoform                 | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 98-82-8                               | Isopropylbenzene          | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 79-34-5                               | 1,1,2,2-Tetrachloroethane | 0.50   | U         | 0.26     | 0.50 | 1.00       | ug/L    |
| 541-73-1                              | 1,3-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| 106-46-7                              | 1,4-Dichlorobenzene       | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 95-50-1                               | 1,2-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| <b>SURROGATES</b>                     |                           |        |           |          |      |            |         |
| 17060-07-0                            | 1,2-Dichloroethane-d4     | 54.9   |           | 81 - 118 |      | 110%       | SPK: 50 |
| 1868-53-7                             | Dibromofluoromethane      | 53.0   |           | 80 - 119 |      | 106%       | SPK: 50 |
| 2037-26-5                             | Toluene-d8                | 51.3   |           | 89 - 112 |      | 103%       | SPK: 50 |
| 460-00-4                              | 4-Bromofluorobenzene      | 49.6   |           | 85 - 114 |      | 99%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b>             |                           |        |           |          |      |            |         |
| 363-72-4                              | Pentafluorobenzene        | 61200  | 5.55      |          |      |            |         |
| 540-36-3                              | 1,4-Difluorobenzene       | 121000 | 6.757     |          |      |            |         |
| 3114-55-4                             | Chlorobenzene-d5          | 115000 | 10.055    |          |      |            |         |
| 3855-82-1                             | 1,4-Dichlorobenzene-d4    | 47600  | 12.018    |          |      |            |         |
| <b>TENTATIVE IDENTIFIED COMPOUNDS</b> |                           |        |           |          |      |            |         |
| 75-43-4                               | Dichlorofluoromethane     | N.D    |           |          |      |            |         |

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/12/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-100-102         |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-03                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046211.D        | 1         |           | 05/15/25 14:11 | VX051525      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOD | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|-----|------------|-------|
|------------|-----------|-------|-----------|-----|-----|------------|-------|

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

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/13/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-150-152         |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-04                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046212.D        | 1         |           | 05/15/25 14:34 | VX051525      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOD  | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |      |            |       |
| 74-87-3        | Chloromethane                  | 0.50  | U         | 0.32  | 0.50 | 1.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 0.75  | U         | 0.26  | 0.75 | 1.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 3.80  | U         | 1.40  | 3.80 | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 0.75  | U         | 0.47  | 0.75 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 0.50  | U         | 0.33  | 0.50 | 1.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 0.75  | U         | 0.23  | 0.75 | 1.00       | ug/L  |
| 67-64-1        | Acetone                        | 4.00  | J         | 1.50  | 3.80 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 0.75  | U         | 0.21  | 0.75 | 1.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 0.50  | U         | 0.28  | 0.50 | 1.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 2.50  | U         | 0.98  | 2.50 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 0.75  | U         | 0.19  | 0.75 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                     | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 71-43-2        | Benzene                        | 0.50  | U         | 0.15  | 0.50 | 1.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 0.75  | U         | 0.090 | 0.75 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 2.50  | U         | 0.68  | 2.50 | 5.00       | ug/L  |
| 108-88-3       | Toluene                        | 0.50  | U         | 0.14  | 0.50 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene          | 0.50  | U         | 0.17  | 0.50 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 79-00-5        | 1,1,2-Trichloroethane          | 0.50  | U         | 0.21  | 0.50 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                     | 2.50  | U         | 0.89  | 2.50 | 5.00       | ug/L  |

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/13/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-150-152         |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-04                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046212.D        | 1         |           | 05/15/25 14:34 | VX051525      |

| CAS Number                            | Parameter                 | Conc.  | Qualifier | MDL      | LOD  | LOQ / CRQL | Units   |
|---------------------------------------|---------------------------|--------|-----------|----------|------|------------|---------|
| 124-48-1                              | Dibromochloromethane      | 0.50   | U         | 0.18     | 0.50 | 1.00       | ug/L    |
| 127-18-4                              | Tetrachloroethene         | 0.50   | U         | 0.23     | 0.50 | 1.00       | ug/L    |
| 108-90-7                              | Chlorobenzene             | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-41-4                              | Ethyl Benzene             | 0.50   | U         | 0.13     | 0.50 | 1.00       | ug/L    |
| 179601-23-1                           | m/p-Xylenes               | 1.00   | U         | 0.24     | 1.00 | 2.00       | ug/L    |
| 95-47-6                               | o-Xylene                  | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-42-5                              | Styrene                   | 0.50   | U         | 0.15     | 0.50 | 1.00       | ug/L    |
| 75-25-2                               | Bromoform                 | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 98-82-8                               | Isopropylbenzene          | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 79-34-5                               | 1,1,2,2-Tetrachloroethane | 0.50   | U         | 0.26     | 0.50 | 1.00       | ug/L    |
| 541-73-1                              | 1,3-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| 106-46-7                              | 1,4-Dichlorobenzene       | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 95-50-1                               | 1,2-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| <b>SURROGATES</b>                     |                           |        |           |          |      |            |         |
| 17060-07-0                            | 1,2-Dichloroethane-d4     | 55.7   |           | 81 - 118 |      | 111%       | SPK: 50 |
| 1868-53-7                             | Dibromofluoromethane      | 52.4   |           | 80 - 119 |      | 105%       | SPK: 50 |
| 2037-26-5                             | Toluene-d8                | 50.5   |           | 89 - 112 |      | 101%       | SPK: 50 |
| 460-00-4                              | 4-Bromofluorobenzene      | 50.1   |           | 85 - 114 |      | 100%       | SPK: 50 |
| <b>INTERNAL STANDARDS</b>             |                           |        |           |          |      |            |         |
| 363-72-4                              | Pentafluorobenzene        | 63700  | 5.55      |          |      |            |         |
| 540-36-3                              | 1,4-Difluorobenzene       | 128000 | 6.757     |          |      |            |         |
| 3114-55-4                             | Chlorobenzene-d5          | 120000 | 10.049    |          |      |            |         |
| 3855-82-1                             | 1,4-Dichlorobenzene-d4    | 51500  | 12.018    |          |      |            |         |
| <b>TENTATIVE IDENTIFIED COMPOUNDS</b> |                           |        |           |          |      |            |         |
| 75-43-4                               | Dichlorofluoromethane     | N.D    |           |          |      |            |         |

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/13/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-150-152         |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-04                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046212.D        | 1         |           | 05/15/25 14:34 | VX051525      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOD | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|-----|------------|-------|
|------------|-----------|-------|-----------|-----|-----|------------|-------|

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

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/14/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-205-207         |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-05                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046213.D        | 1         |           | 05/15/25 14:58 | VX051525      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOD  | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |      |            |       |
| 74-87-3        | Chloromethane                  | 0.50  | U         | 0.32  | 0.50 | 1.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 0.75  | U         | 0.26  | 0.75 | 1.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 3.80  | U         | 1.40  | 3.80 | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 0.75  | U         | 0.47  | 0.75 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 0.50  | U         | 0.33  | 0.50 | 1.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 0.75  | U         | 0.23  | 0.75 | 1.00       | ug/L  |
| 67-64-1        | Acetone                        | 5.20  |           | 1.50  | 3.80 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 0.75  | U         | 0.21  | 0.75 | 1.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 0.50  | U         | 0.28  | 0.50 | 1.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 2.50  | U         | 0.98  | 2.50 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 0.75  | U         | 0.19  | 0.75 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                     | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 71-43-2        | Benzene                        | 0.50  | U         | 0.15  | 0.50 | 1.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 0.75  | U         | 0.090 | 0.75 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 2.50  | U         | 0.68  | 2.50 | 5.00       | ug/L  |
| 108-88-3       | Toluene                        | 0.50  | U         | 0.14  | 0.50 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene          | 0.50  | U         | 0.17  | 0.50 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 79-00-5        | 1,1,2-Trichloroethane          | 0.50  | U         | 0.21  | 0.50 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                     | 2.50  | U         | 0.89  | 2.50 | 5.00       | ug/L  |



## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/14/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-205-207         |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-05                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046213.D        | 1         |           | 05/15/25 14:58 | VX051525      |

| CAS Number                            | Parameter                 | Conc.  | Qualifier | MDL      | LOD  | LOQ / CRQL | Units   |
|---------------------------------------|---------------------------|--------|-----------|----------|------|------------|---------|
| 124-48-1                              | Dibromochloromethane      | 0.50   | U         | 0.18     | 0.50 | 1.00       | ug/L    |
| 127-18-4                              | Tetrachloroethene         | 0.50   | U         | 0.23     | 0.50 | 1.00       | ug/L    |
| 108-90-7                              | Chlorobenzene             | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-41-4                              | Ethyl Benzene             | 0.50   | U         | 0.13     | 0.50 | 1.00       | ug/L    |
| 179601-23-1                           | m/p-Xylenes               | 1.00   | U         | 0.24     | 1.00 | 2.00       | ug/L    |
| 95-47-6                               | o-Xylene                  | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-42-5                              | Styrene                   | 0.50   | U         | 0.15     | 0.50 | 1.00       | ug/L    |
| 75-25-2                               | Bromoform                 | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 98-82-8                               | Isopropylbenzene          | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 79-34-5                               | 1,1,2,2-Tetrachloroethane | 0.50   | U         | 0.26     | 0.50 | 1.00       | ug/L    |
| 541-73-1                              | 1,3-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| 106-46-7                              | 1,4-Dichlorobenzene       | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 95-50-1                               | 1,2-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| <b>SURROGATES</b>                     |                           |        |           |          |      |            |         |
| 17060-07-0                            | 1,2-Dichloroethane-d4     | 55.4   |           | 81 - 118 |      | 111%       | SPK: 50 |
| 1868-53-7                             | Dibromofluoromethane      | 52.0   |           | 80 - 119 |      | 104%       | SPK: 50 |
| 2037-26-5                             | Toluene-d8                | 51.5   |           | 89 - 112 |      | 103%       | SPK: 50 |
| 460-00-4                              | 4-Bromofluorobenzene      | 49.5   |           | 85 - 114 |      | 99%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b>             |                           |        |           |          |      |            |         |
| 363-72-4                              | Pentafluorobenzene        | 60500  | 5.55      |          |      |            |         |
| 540-36-3                              | 1,4-Difluorobenzene       | 121000 | 6.757     |          |      |            |         |
| 3114-55-4                             | Chlorobenzene-d5          | 113000 | 10.055    |          |      |            |         |
| 3855-82-1                             | 1,4-Dichlorobenzene-d4    | 47400  | 12.018    |          |      |            |         |
| <b>TENTATIVE IDENTIFIED COMPOUNDS</b> |                           |        |           |          |      |            |         |
| 75-43-4                               | Dichlorofluoromethane     | N.D    |           |          |      |            |         |

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/14/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-205-207         |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-05                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046213.D        | 1         |           | 05/15/25 14:58 | VX051525      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOD | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|-----|------------|-------|
|------------|-----------|-------|-----------|-----|-----|------------|-------|

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

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/14/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-220-222         |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-06                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046214.D        | 1         |           | 05/15/25 15:21 | VX051525      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOD  | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |      |            |       |
| 74-87-3        | Chloromethane                  | 0.50  | U         | 0.32  | 0.50 | 1.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 0.75  | U         | 0.26  | 0.75 | 1.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 3.80  | U         | 1.40  | 3.80 | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 0.75  | U         | 0.47  | 0.75 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 0.50  | U         | 0.33  | 0.50 | 1.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 0.75  | U         | 0.23  | 0.75 | 1.00       | ug/L  |
| 67-64-1        | Acetone                        | 3.20  | J         | 1.50  | 3.80 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 0.75  | U         | 0.21  | 0.75 | 1.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 0.50  | U         | 0.28  | 0.50 | 1.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 2.50  | U         | 0.98  | 2.50 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 0.75  | U         | 0.19  | 0.75 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                     | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 71-43-2        | Benzene                        | 0.50  | U         | 0.15  | 0.50 | 1.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 0.75  | U         | 0.090 | 0.75 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 2.50  | U         | 0.68  | 2.50 | 5.00       | ug/L  |
| 108-88-3       | Toluene                        | 0.50  | U         | 0.14  | 0.50 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene          | 0.50  | U         | 0.17  | 0.50 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 79-00-5        | 1,1,2-Trichloroethane          | 0.50  | U         | 0.21  | 0.50 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                     | 2.50  | U         | 0.89  | 2.50 | 5.00       | ug/L  |

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/14/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-220-222         |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-06                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046214.D        | 1         |           | 05/15/25 15:21 | VX051525      |

| CAS Number                            | Parameter                 | Conc.  | Qualifier | MDL      | LOD  | LOQ / CRQL | Units   |
|---------------------------------------|---------------------------|--------|-----------|----------|------|------------|---------|
| 124-48-1                              | Dibromochloromethane      | 0.50   | U         | 0.18     | 0.50 | 1.00       | ug/L    |
| 127-18-4                              | Tetrachloroethene         | 0.50   | U         | 0.23     | 0.50 | 1.00       | ug/L    |
| 108-90-7                              | Chlorobenzene             | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-41-4                              | Ethyl Benzene             | 0.50   | U         | 0.13     | 0.50 | 1.00       | ug/L    |
| 179601-23-1                           | m/p-Xylenes               | 1.00   | U         | 0.24     | 1.00 | 2.00       | ug/L    |
| 95-47-6                               | o-Xylene                  | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-42-5                              | Styrene                   | 0.50   | U         | 0.15     | 0.50 | 1.00       | ug/L    |
| 75-25-2                               | Bromoform                 | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 98-82-8                               | Isopropylbenzene          | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 79-34-5                               | 1,1,2,2-Tetrachloroethane | 0.50   | U         | 0.26     | 0.50 | 1.00       | ug/L    |
| 541-73-1                              | 1,3-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| 106-46-7                              | 1,4-Dichlorobenzene       | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 95-50-1                               | 1,2-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| <b>SURROGATES</b>                     |                           |        |           |          |      |            |         |
| 17060-07-0                            | 1,2-Dichloroethane-d4     | 54.6   |           | 81 - 118 |      | 109%       | SPK: 50 |
| 1868-53-7                             | Dibromofluoromethane      | 51.6   |           | 80 - 119 |      | 103%       | SPK: 50 |
| 2037-26-5                             | Toluene-d8                | 50.8   |           | 89 - 112 |      | 102%       | SPK: 50 |
| 460-00-4                              | 4-Bromofluorobenzene      | 48.4   |           | 85 - 114 |      | 97%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b>             |                           |        |           |          |      |            |         |
| 363-72-4                              | Pentafluorobenzene        | 60700  | 5.55      |          |      |            |         |
| 540-36-3                              | 1,4-Difluorobenzene       | 121000 | 6.763     |          |      |            |         |
| 3114-55-4                             | Chlorobenzene-d5          | 112000 | 10.055    |          |      |            |         |
| 3855-82-1                             | 1,4-Dichlorobenzene-d4    | 45000  | 12.018    |          |      |            |         |
| <b>TENTATIVE IDENTIFIED COMPOUNDS</b> |                           |        |           |          |      |            |         |
| 75-43-4                               | Dichlorofluoromethane     | N.D    |           |          |      |            |         |

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/14/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-220-222         |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-06                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046214.D        | 1         |           | 05/15/25 15:21 | VX051525      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOD | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|-----|------------|-------|
|------------|-----------|-------|-----------|-----|-----|------------|-------|

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

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/14/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-DUP-20250514       |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-07                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046215.D        | 1         |           | 05/15/25 15:44 | VX051525      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOD  | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |      |            |       |
| 74-87-3        | Chloromethane                  | 0.50  | U         | 0.32  | 0.50 | 1.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 0.75  | U         | 0.26  | 0.75 | 1.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 3.80  | U         | 1.40  | 3.80 | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 0.75  | U         | 0.47  | 0.75 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 0.50  | U         | 0.33  | 0.50 | 1.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 0.75  | U         | 0.23  | 0.75 | 1.00       | ug/L  |
| 67-64-1        | Acetone                        | 2.80  | J         | 1.50  | 3.80 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 0.75  | U         | 0.21  | 0.75 | 1.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 0.50  | U         | 0.28  | 0.50 | 1.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 2.50  | U         | 0.98  | 2.50 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 0.75  | U         | 0.19  | 0.75 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                     | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 71-43-2        | Benzene                        | 0.50  | U         | 0.15  | 0.50 | 1.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 0.75  | U         | 0.090 | 0.75 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 2.50  | U         | 0.68  | 2.50 | 5.00       | ug/L  |
| 108-88-3       | Toluene                        | 0.50  | U         | 0.14  | 0.50 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene          | 0.50  | U         | 0.17  | 0.50 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 79-00-5        | 1,1,2-Trichloroethane          | 0.50  | U         | 0.21  | 0.50 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                     | 2.50  | U         | 0.89  | 2.50 | 5.00       | ug/L  |

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/14/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-DUP-20250514       |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-07                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046215.D        | 1         |           | 05/15/25 15:44 | VX051525      |

| CAS Number                            | Parameter                 | Conc.  | Qualifier | MDL      | LOD  | LOQ / CRQL | Units   |
|---------------------------------------|---------------------------|--------|-----------|----------|------|------------|---------|
| 124-48-1                              | Dibromochloromethane      | 0.50   | U         | 0.18     | 0.50 | 1.00       | ug/L    |
| 127-18-4                              | Tetrachloroethene         | 0.50   | U         | 0.23     | 0.50 | 1.00       | ug/L    |
| 108-90-7                              | Chlorobenzene             | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-41-4                              | Ethyl Benzene             | 0.50   | U         | 0.13     | 0.50 | 1.00       | ug/L    |
| 179601-23-1                           | m/p-Xylenes               | 1.00   | U         | 0.24     | 1.00 | 2.00       | ug/L    |
| 95-47-6                               | o-Xylene                  | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-42-5                              | Styrene                   | 0.50   | U         | 0.15     | 0.50 | 1.00       | ug/L    |
| 75-25-2                               | Bromoform                 | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 98-82-8                               | Isopropylbenzene          | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 79-34-5                               | 1,1,2,2-Tetrachloroethane | 0.50   | U         | 0.26     | 0.50 | 1.00       | ug/L    |
| 541-73-1                              | 1,3-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| 106-46-7                              | 1,4-Dichlorobenzene       | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 95-50-1                               | 1,2-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| <b>SURROGATES</b>                     |                           |        |           |          |      |            |         |
| 17060-07-0                            | 1,2-Dichloroethane-d4     | 55.6   |           | 81 - 118 |      | 111%       | SPK: 50 |
| 1868-53-7                             | Dibromofluoromethane      | 52.0   |           | 80 - 119 |      | 104%       | SPK: 50 |
| 2037-26-5                             | Toluene-d8                | 50.6   |           | 89 - 112 |      | 101%       | SPK: 50 |
| 460-00-4                              | 4-Bromofluorobenzene      | 51.4   |           | 85 - 114 |      | 103%       | SPK: 50 |
| <b>INTERNAL STANDARDS</b>             |                           |        |           |          |      |            |         |
| 363-72-4                              | Pentafluorobenzene        | 61100  | 5.544     |          |      |            |         |
| 540-36-3                              | 1,4-Difluorobenzene       | 122000 | 6.757     |          |      |            |         |
| 3114-55-4                             | Chlorobenzene-d5          | 117000 | 10.049    |          |      |            |         |
| 3855-82-1                             | 1,4-Dichlorobenzene-d4    | 50000  | 12.018    |          |      |            |         |
| <b>TENTATIVE IDENTIFIED COMPOUNDS</b> |                           |        |           |          |      |            |         |
| 75-43-4                               | Dichlorofluoromethane     | N.D    |           |          |      |            |         |

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/14/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-DUP-20250514       |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-07                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046215.D        | 1         |           | 05/15/25 15:44 | VX051525      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOD | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|-----|------------|-------|
|------------|-----------|-------|-----------|-----|-----|------------|-------|

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



## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/14/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-EB-20250514        |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-08                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046210.D        | 1         |           | 05/15/25 13:47 | VX051525      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOD  | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |      |            |       |
| 74-87-3        | Chloromethane                  | 0.50  | U         | 0.32  | 0.50 | 1.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 0.75  | U         | 0.26  | 0.75 | 1.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 3.80  | U         | 1.40  | 3.80 | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 0.75  | U         | 0.47  | 0.75 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 0.50  | U         | 0.33  | 0.50 | 1.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 0.75  | U         | 0.23  | 0.75 | 1.00       | ug/L  |
| 67-64-1        | Acetone                        | 4.30  | J         | 1.50  | 3.80 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 0.75  | U         | 0.21  | 0.75 | 1.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 0.50  | U         | 0.28  | 0.50 | 1.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 2.50  | U         | 0.98  | 2.50 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 0.75  | U         | 0.19  | 0.75 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                     | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 71-43-2        | Benzene                        | 0.50  | U         | 0.15  | 0.50 | 1.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 0.75  | U         | 0.090 | 0.75 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 2.50  | U         | 0.68  | 2.50 | 5.00       | ug/L  |
| 108-88-3       | Toluene                        | 0.50  | U         | 0.14  | 0.50 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene          | 0.50  | U         | 0.17  | 0.50 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 79-00-5        | 1,1,2-Trichloroethane          | 0.50  | U         | 0.21  | 0.50 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                     | 2.50  | U         | 0.89  | 2.50 | 5.00       | ug/L  |

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/14/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-EB-20250514        |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-08                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046210.D        | 1         |           | 05/15/25 13:47 | VX051525      |

| CAS Number                            | Parameter                 | Conc.  | Qualifier | MDL      | LOD  | LOQ / CRQL | Units   |
|---------------------------------------|---------------------------|--------|-----------|----------|------|------------|---------|
| 124-48-1                              | Dibromochloromethane      | 0.50   | U         | 0.18     | 0.50 | 1.00       | ug/L    |
| 127-18-4                              | Tetrachloroethene         | 0.50   | U         | 0.23     | 0.50 | 1.00       | ug/L    |
| 108-90-7                              | Chlorobenzene             | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-41-4                              | Ethyl Benzene             | 0.50   | U         | 0.13     | 0.50 | 1.00       | ug/L    |
| 179601-23-1                           | m/p-Xylenes               | 1.00   | U         | 0.24     | 1.00 | 2.00       | ug/L    |
| 95-47-6                               | o-Xylene                  | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-42-5                              | Styrene                   | 0.50   | U         | 0.15     | 0.50 | 1.00       | ug/L    |
| 75-25-2                               | Bromoform                 | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 98-82-8                               | Isopropylbenzene          | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 79-34-5                               | 1,1,2,2-Tetrachloroethane | 0.50   | U         | 0.26     | 0.50 | 1.00       | ug/L    |
| 541-73-1                              | 1,3-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| 106-46-7                              | 1,4-Dichlorobenzene       | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 95-50-1                               | 1,2-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| <b>SURROGATES</b>                     |                           |        |           |          |      |            |         |
| 17060-07-0                            | 1,2-Dichloroethane-d4     | 54.1   |           | 81 - 118 |      | 108%       | SPK: 50 |
| 1868-53-7                             | Dibromofluoromethane      | 51.8   |           | 80 - 119 |      | 104%       | SPK: 50 |
| 2037-26-5                             | Toluene-d8                | 50.5   |           | 89 - 112 |      | 101%       | SPK: 50 |
| 460-00-4                              | 4-Bromofluorobenzene      | 51.9   |           | 85 - 114 |      | 104%       | SPK: 50 |
| <b>INTERNAL STANDARDS</b>             |                           |        |           |          |      |            |         |
| 363-72-4                              | Pentafluorobenzene        | 61600  | 5.544     |          |      |            |         |
| 540-36-3                              | 1,4-Difluorobenzene       | 122000 | 6.757     |          |      |            |         |
| 3114-55-4                             | Chlorobenzene-d5          | 116000 | 10.049    |          |      |            |         |
| 3855-82-1                             | 1,4-Dichlorobenzene-d4    | 50300  | 12.018    |          |      |            |         |
| <b>TENTATIVE IDENTIFIED COMPOUNDS</b> |                           |        |           |          |      |            |         |
| 75-43-4                               | Dichlorofluoromethane     | N.D    |           |          |      |            |         |

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/14/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-EB-20250514        |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-08                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046210.D        | 1         |           | 05/15/25 13:47 | VX051525      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOD | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|-----|------------|-------|
|------------|-----------|-------|-----------|-----|-----|------------|-------|

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

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/14/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-240-242         |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-09                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046216.D        | 1         |           | 05/15/25 16:08 | VX051525      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOD  | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |      |            |       |
| 74-87-3        | Chloromethane                  | 0.50  | U         | 0.32  | 0.50 | 1.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 0.75  | U         | 0.26  | 0.75 | 1.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 3.80  | U         | 1.40  | 3.80 | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 0.75  | U         | 0.47  | 0.75 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 0.50  | U         | 0.33  | 0.50 | 1.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 0.75  | U         | 0.23  | 0.75 | 1.00       | ug/L  |
| 67-64-1        | Acetone                        | 2.90  | J         | 1.50  | 3.80 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 0.75  | U         | 0.21  | 0.75 | 1.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 0.50  | U         | 0.28  | 0.50 | 1.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 2.50  | U         | 0.98  | 2.50 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 0.75  | U         | 0.19  | 0.75 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                     | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 71-43-2        | Benzene                        | 0.50  | U         | 0.15  | 0.50 | 1.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 0.75  | U         | 0.090 | 0.75 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 2.50  | U         | 0.68  | 2.50 | 5.00       | ug/L  |
| 108-88-3       | Toluene                        | 0.50  | U         | 0.14  | 0.50 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene          | 0.50  | U         | 0.17  | 0.50 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 79-00-5        | 1,1,2-Trichloroethane          | 0.50  | U         | 0.21  | 0.50 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                     | 2.50  | U         | 0.89  | 2.50 | 5.00       | ug/L  |

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/14/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-240-242         |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-09                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046216.D        | 1         |           | 05/15/25 16:08 | VX051525      |

| CAS Number                            | Parameter                 | Conc.  | Qualifier | MDL      | LOD  | LOQ / CRQL | Units   |
|---------------------------------------|---------------------------|--------|-----------|----------|------|------------|---------|
| 124-48-1                              | Dibromochloromethane      | 0.50   | U         | 0.18     | 0.50 | 1.00       | ug/L    |
| 127-18-4                              | Tetrachloroethene         | 0.50   | U         | 0.23     | 0.50 | 1.00       | ug/L    |
| 108-90-7                              | Chlorobenzene             | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-41-4                              | Ethyl Benzene             | 0.50   | U         | 0.13     | 0.50 | 1.00       | ug/L    |
| 179601-23-1                           | m/p-Xylenes               | 1.00   | U         | 0.24     | 1.00 | 2.00       | ug/L    |
| 95-47-6                               | o-Xylene                  | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-42-5                              | Styrene                   | 0.50   | U         | 0.15     | 0.50 | 1.00       | ug/L    |
| 75-25-2                               | Bromoform                 | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 98-82-8                               | Isopropylbenzene          | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 79-34-5                               | 1,1,2,2-Tetrachloroethane | 0.50   | U         | 0.26     | 0.50 | 1.00       | ug/L    |
| 541-73-1                              | 1,3-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| 106-46-7                              | 1,4-Dichlorobenzene       | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 95-50-1                               | 1,2-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| <b>SURROGATES</b>                     |                           |        |           |          |      |            |         |
| 17060-07-0                            | 1,2-Dichloroethane-d4     | 55.0   |           | 81 - 118 |      | 110%       | SPK: 50 |
| 1868-53-7                             | Dibromofluoromethane      | 52.2   |           | 80 - 119 |      | 104%       | SPK: 50 |
| 2037-26-5                             | Toluene-d8                | 50.6   |           | 89 - 112 |      | 101%       | SPK: 50 |
| 460-00-4                              | 4-Bromofluorobenzene      | 50.4   |           | 85 - 114 |      | 101%       | SPK: 50 |
| <b>INTERNAL STANDARDS</b>             |                           |        |           |          |      |            |         |
| 363-72-4                              | Pentafluorobenzene        | 62400  | 5.55      |          |      |            |         |
| 540-36-3                              | 1,4-Difluorobenzene       | 125000 | 6.757     |          |      |            |         |
| 3114-55-4                             | Chlorobenzene-d5          | 120000 | 10.055    |          |      |            |         |
| 3855-82-1                             | 1,4-Dichlorobenzene-d4    | 50200  | 12.018    |          |      |            |         |
| <b>TENTATIVE IDENTIFIED COMPOUNDS</b> |                           |        |           |          |      |            |         |
| 75-43-4                               | Dichlorofluoromethane     | N.D    |           |          |      |            |         |

## Report of Analysis

|                    |                               |           |                 |              |    |
|--------------------|-------------------------------|-----------|-----------------|--------------|----|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: | 05/14/25     |    |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  | 05/14/25     |    |
| Client Sample ID:  | BP-VPB-182-GW-240-242         |           | SDG No.:        | Q2050        |    |
| Lab Sample ID:     | Q2050-09                      |           | Matrix:         | Water        |    |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |    |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000         | uL |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |    |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |    |
| Prep Method :      |                               |           |                 |              |    |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046216.D        | 1         |           | 05/15/25 16:08 | VX051525      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOD | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|-----|------------|-------|
|------------|-----------|-------|-----------|-----|-----|------------|-------|

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



# QC SUMMARY

## Surrogate Summary

SDG No.: **Q2050**

Client: **Tetra Tech NUS, Inc.**

Analytical Method: **SW8260-Low**

| Lab Sample ID | Client ID               | Parameter             | Spike | Result | RecoveryQual | Limits |      |
|---------------|-------------------------|-----------------------|-------|--------|--------------|--------|------|
|               |                         |                       |       |        |              | Low    | High |
| Q2050-01      | BP-VPB-182-TB-20250512  | 1,2-Dichloroethane-d4 | 50    | 55.1   | 110          | 81     | 118  |
|               |                         | Dibromofluoromethane  | 50    | 52.3   | 104          | 80     | 119  |
|               |                         | Toluene-d8            | 50    | 51.0   | 102          | 89     | 112  |
|               |                         | 4-Bromofluorobenzene  | 50    | 50.7   | 101          | 85     | 114  |
| Q2050-02      | BP-VPB-182-GW-60-62     | 1,2-Dichloroethane-d4 | 50    | 54.7   | 109          | 81     | 118  |
|               |                         | Dibromofluoromethane  | 50    | 51.3   | 103          | 80     | 119  |
|               |                         | Toluene-d8            | 50    | 49.8   | 100          | 89     | 112  |
|               |                         | 4-Bromofluorobenzene  | 50    | 50.0   | 100          | 85     | 114  |
| Q2050-03      | BP-VPB-182-GW-100-102   | 1,2-Dichloroethane-d4 | 50    | 54.9   | 110          | 81     | 118  |
|               |                         | Dibromofluoromethane  | 50    | 53.0   | 106          | 80     | 119  |
|               |                         | Toluene-d8            | 50    | 51.3   | 103          | 89     | 112  |
|               |                         | 4-Bromofluorobenzene  | 50    | 49.6   | 99           | 85     | 114  |
| Q2050-04      | BP-VPB-182-GW-150-152   | 1,2-Dichloroethane-d4 | 50    | 55.6   | 111          | 81     | 118  |
|               |                         | Dibromofluoromethane  | 50    | 52.4   | 105          | 80     | 119  |
|               |                         | Toluene-d8            | 50    | 50.5   | 101          | 89     | 112  |
|               |                         | 4-Bromofluorobenzene  | 50    | 50.0   | 100          | 85     | 114  |
| Q2050-05      | BP-VPB-182-GW-205-207   | 1,2-Dichloroethane-d4 | 50    | 55.5   | 111          | 81     | 118  |
|               |                         | Dibromofluoromethane  | 50    | 52.0   | 104          | 80     | 119  |
|               |                         | Toluene-d8            | 50    | 51.5   | 103          | 89     | 112  |
|               |                         | 4-Bromofluorobenzene  | 50    | 49.5   | 99           | 85     | 114  |
| Q2050-06      | BP-VPB-182-GW-220-222   | 1,2-Dichloroethane-d4 | 50    | 54.6   | 109          | 81     | 118  |
|               |                         | Dibromofluoromethane  | 50    | 51.6   | 103          | 80     | 119  |
|               |                         | Toluene-d8            | 50    | 50.8   | 102          | 89     | 112  |
|               |                         | 4-Bromofluorobenzene  | 50    | 48.4   | 97           | 85     | 114  |
| Q2050-07      | BP-VPB-182-DUP-20250514 | 1,2-Dichloroethane-d4 | 50    | 55.6   | 111          | 81     | 118  |
|               |                         | Dibromofluoromethane  | 50    | 52.0   | 104          | 80     | 119  |
|               |                         | Toluene-d8            | 50    | 50.6   | 101          | 89     | 112  |
|               |                         | 4-Bromofluorobenzene  | 50    | 51.4   | 103          | 85     | 114  |
| Q2050-08      | BP-VPB-182-EB-20250514  | 1,2-Dichloroethane-d4 | 50    | 54.1   | 108          | 81     | 118  |
|               |                         | Dibromofluoromethane  | 50    | 51.8   | 104          | 80     | 119  |
|               |                         | Toluene-d8            | 50    | 50.5   | 101          | 89     | 112  |
|               |                         | 4-Bromofluorobenzene  | 50    | 51.9   | 104          | 85     | 114  |
| Q2050-09      | BP-VPB-182-GW-240-242   | 1,2-Dichloroethane-d4 | 50    | 55.0   | 110          | 81     | 118  |
|               |                         | Dibromofluoromethane  | 50    | 52.1   | 104          | 80     | 119  |
|               |                         | Toluene-d8            | 50    | 50.6   | 101          | 89     | 112  |
|               |                         | 4-Bromofluorobenzene  | 50    | 50.4   | 101          | 85     | 114  |
| VX0515WBL01   | VX0515WBL01             | 1,2-Dichloroethane-d4 | 50    | 54.2   | 108          | 81     | 118  |
|               |                         | Dibromofluoromethane  | 50    | 51.5   | 103          | 80     | 119  |
|               |                         | Toluene-d8            | 50    | 50.2   | 100          | 89     | 112  |
|               |                         | 4-Bromofluorobenzene  | 50    | 49.0   | 98           | 85     | 114  |
| VX0515WBS01   | VX0515WBS01             | 1,2-Dichloroethane-d4 | 50    | 50.7   | 101          | 81     | 118  |
|               |                         | Dibromofluoromethane  | 50    | 50.0   | 100          | 80     | 119  |
|               |                         | Toluene-d8            | 50    | 49.7   | 99           | 89     | 112  |
|               |                         | 4-Bromofluorobenzene  | 50    | 48.9   | 98           | 85     | 114  |
| VX0515WBSD0   | VX0515WBSD01            | 1,2-Dichloroethane-d4 | 50    | 49.4   | 99           | 81     | 118  |
|               |                         | Dibromofluoromethane  | 50    | 49.3   | 99           | 80     | 119  |
|               |                         | Toluene-d8            | 50    | 48.5   | 97           | 89     | 112  |
|               |                         | 4-Bromofluorobenzene  | 50    | 47.9   | 96           | 85     | 114  |



Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q2050

Client: Tetra Tech NUS, Inc.

Analytical Method: SW8260-Low

Datafile : VX046206.D

| Lab Sample ID | Parameter                      | Spike | Result | Unit | Rec | RPD | Qual | Low | Limits<br>High | RPD |
|---------------|--------------------------------|-------|--------|------|-----|-----|------|-----|----------------|-----|
| VX0515WBS01   | Chloromethane                  | 20    | 19.5   | ug/L | 98  |     |      | 50  | 139            |     |
|               | Vinyl chloride                 | 20    | 19.1   | ug/L | 96  |     |      | 58  | 137            |     |
|               | Bromomethane                   | 20    | 18.9   | ug/L | 95  |     |      | 53  | 141            |     |
|               | Chloroethane                   | 20    | 22.5   | ug/L | 113 |     |      | 60  | 138            |     |
|               | Trichlorofluoromethane         | 20    | 20.4   | ug/L | 102 |     |      | 65  | 141            |     |
|               | 1,1,2-Trichlorotrifluoroethane | 20    | 19.9   | ug/L | 100 |     |      | 70  | 136            |     |
|               | 1,1-Dichloroethene             | 20    | 19.0   | ug/L | 95  |     |      | 71  | 131            |     |
|               | Acetone                        | 100   | 100    | ug/L | 100 |     |      | 39  | 160            |     |
|               | Carbon disulfide               | 20    | 16.7   | ug/L | 84  |     |      | 64  | 133            |     |
|               | Methyl tert-butyl Ether        | 20    | 19.9   | ug/L | 100 |     |      | 71  | 124            |     |
|               | Methylene Chloride             | 20    | 19.0   | ug/L | 95  |     |      | 74  | 124            |     |
|               | trans-1,2-Dichloroethene       | 20    | 19.5   | ug/L | 98  |     |      | 75  | 124            |     |
|               | 1,1-Dichloroethane             | 20    | 20.2   | ug/L | 101 |     |      | 77  | 125            |     |
|               | 2-Butanone                     | 100   | 110    | ug/L | 110 |     |      | 56  | 143            |     |
|               | Carbon Tetrachloride           | 20    | 19.5   | ug/L | 98  |     |      | 72  | 136            |     |
|               | cis-1,2-Dichloroethene         | 20    | 20.2   | ug/L | 101 |     |      | 78  | 123            |     |
|               | Chloroform                     | 20    | 21.0   | ug/L | 105 |     |      | 79  | 124            |     |
|               | 1,1,1-Trichloroethane          | 20    | 20.2   | ug/L | 101 |     |      | 74  | 131            |     |
|               | Methylcyclohexane              | 20    | 18.3   | ug/L | 92  |     |      | 72  | 132            |     |
|               | Benzene                        | 20    | 20.1   | ug/L | 101 |     |      | 79  | 120            |     |
|               | 1,2-Dichloroethane             | 20    | 20.6   | ug/L | 103 |     |      | 73  | 128            |     |
|               | Trichloroethene                | 20    | 19.3   | ug/L | 97  |     |      | 79  | 123            |     |
|               | 1,2-Dichloropropane            | 20    | 20.9   | ug/L | 104 |     |      | 78  | 122            |     |
|               | Bromodichloromethane           | 20    | 20.6   | ug/L | 103 |     |      | 79  | 125            |     |
|               | 4-Methyl-2-Pentanone           | 100   | 110    | ug/L | 110 |     |      | 67  | 130            |     |
|               | Toluene                        | 20    | 20.0   | ug/L | 100 |     |      | 80  | 121            |     |
|               | t-1,3-Dichloropropene          | 20    | 18.5   | ug/L | 93  |     |      | 73  | 127            |     |
|               | cis-1,3-Dichloropropene        | 20    | 19.5   | ug/L | 98  |     |      | 75  | 124            |     |
|               | 1,1,2-Trichloroethane          | 20    | 20.7   | ug/L | 104 |     |      | 80  | 119            |     |
|               | 2-Hexanone                     | 100   | 110    | ug/L | 110 |     |      | 57  | 139            |     |
|               | Dibromochloromethane           | 20    | 20.7   | ug/L | 104 |     |      | 74  | 126            |     |
|               | Tetrachloroethene              | 20    | 19.5   | ug/L | 98  |     |      | 74  | 129            |     |
|               | Chlorobenzene                  | 20    | 19.5   | ug/L | 98  |     |      | 82  | 118            |     |
|               | Ethyl Benzene                  | 20    | 19.8   | ug/L | 99  |     |      | 79  | 121            |     |
|               | m/p-Xylenes                    | 40    | 39.4   | ug/L | 99  |     |      | 80  | 121            |     |
|               | o-Xylene                       | 20    | 20.3   | ug/L | 102 |     |      | 78  | 122            |     |
|               | Styrene                        | 20    | 20.5   | ug/L | 103 |     |      | 78  | 123            |     |
|               | Bromoform                      | 20    | 19.1   | ug/L | 96  |     |      | 66  | 130            |     |
|               | Isopropylbenzene               | 20    | 20.1   | ug/L | 101 |     |      | 72  | 131            |     |
|               | 1,1,2,2-Tetrachloroethane      | 20    | 19.8   | ug/L | 99  |     |      | 71  | 121            |     |
|               | 1,3-Dichlorobenzene            | 20    | 19.7   | ug/L | 99  |     |      | 80  | 119            |     |
|               | 1,4-Dichlorobenzene            | 20    | 19.3   | ug/L | 97  |     |      | 79  | 118            |     |
|               | 1,2-Dichlorobenzene            | 20    | 20.3   | ug/L | 102 |     |      | 80  | 119            |     |

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q2050

Client: Tetra Tech NUS, Inc.

Analytical Method: SW8260-Low

Datafile : VX046207.D

| Lab Sample ID | Parameter                      | Spike | Result | Unit | Rec | RPD | Qual | Low | Limits High | RPD |
|---------------|--------------------------------|-------|--------|------|-----|-----|------|-----|-------------|-----|
| VX0515WBSD01  | Chloromethane                  | 20    | 19.3   | ug/L | 97  | 1   |      | 50  | 139         | 20  |
|               | Vinyl chloride                 | 20    | 18.2   | ug/L | 91  | 5   |      | 58  | 137         | 20  |
|               | Bromomethane                   | 20    | 17.8   | ug/L | 89  | 7   |      | 53  | 141         | 20  |
|               | Chloroethane                   | 20    | 19.8   | ug/L | 99  | 13  |      | 60  | 138         | 20  |
|               | Trichlorofluoromethane         | 20    | 19.8   | ug/L | 99  | 3   |      | 65  | 141         | 20  |
|               | 1,1,2-Trichlorotrifluoroethane | 20    | 19.3   | ug/L | 97  | 3   |      | 70  | 136         | 20  |
|               | 1,1-Dichloroethene             | 20    | 17.9   | ug/L | 90  | 5   |      | 71  | 131         | 20  |
|               | Acetone                        | 100   | 100    | ug/L | 100 | 0   |      | 39  | 160         | 20  |
|               | Carbon disulfide               | 20    | 16.1   | ug/L | 81  | 4   |      | 64  | 133         | 20  |
|               | Methyl tert-butyl Ether        | 20    | 20.1   | ug/L | 101 | 1   |      | 71  | 124         | 20  |
|               | Methylene Chloride             | 20    | 18.7   | ug/L | 94  | 1   |      | 74  | 124         | 20  |
|               | trans-1,2-Dichloroethene       | 20    | 19.0   | ug/L | 95  | 3   |      | 75  | 124         | 20  |
|               | 1,1-Dichloroethane             | 20    | 20.0   | ug/L | 100 | 1   |      | 77  | 125         | 20  |
|               | 2-Butanone                     | 100   | 110    | ug/L | 110 | 0   |      | 56  | 143         | 20  |
|               | Carbon Tetrachloride           | 20    | 19.2   | ug/L | 96  | 2   |      | 72  | 136         | 20  |
|               | cis-1,2-Dichloroethene         | 20    | 19.3   | ug/L | 97  | 4   |      | 78  | 123         | 20  |
|               | Chloroform                     | 20    | 20.6   | ug/L | 103 | 2   |      | 79  | 124         | 20  |
|               | 1,1,1-Trichloroethane          | 20    | 19.8   | ug/L | 99  | 2   |      | 74  | 131         | 20  |
|               | Methylcyclohexane              | 20    | 17.9   | ug/L | 90  | 2   |      | 72  | 132         | 20  |
|               | Benzene                        | 20    | 19.7   | ug/L | 99  | 2   |      | 79  | 120         | 20  |
|               | 1,2-Dichloroethane             | 20    | 20.7   | ug/L | 104 | 1   |      | 73  | 128         | 20  |
|               | Trichloroethene                | 20    | 18.8   | ug/L | 94  | 3   |      | 79  | 123         | 20  |
|               | 1,2-Dichloropropane            | 20    | 21.3   | ug/L | 106 | 2   |      | 78  | 122         | 20  |
|               | Bromodichloromethane           | 20    | 20.3   | ug/L | 102 | 1   |      | 79  | 125         | 20  |
|               | 4-Methyl-2-Pentanone           | 100   | 110    | ug/L | 110 | 0   |      | 67  | 130         | 20  |
|               | Toluene                        | 20    | 20.1   | ug/L | 101 | 1   |      | 80  | 121         | 20  |
|               | t-1,3-Dichloropropene          | 20    | 18.8   | ug/L | 94  | 1   |      | 73  | 127         | 20  |
|               | cis-1,3-Dichloropropene        | 20    | 19.6   | ug/L | 98  | 0   |      | 75  | 124         | 20  |
|               | 1,1,2-Trichloroethane          | 20    | 21.5   | ug/L | 108 | 4   |      | 80  | 119         | 20  |
|               | 2-Hexanone                     | 100   | 110    | ug/L | 110 | 0   |      | 57  | 139         | 20  |
|               | Dibromochloromethane           | 20    | 20.7   | ug/L | 104 | 0   |      | 74  | 126         | 20  |
|               | Tetrachloroethene              | 20    | 19.3   | ug/L | 97  | 1   |      | 74  | 129         | 20  |
|               | Chlorobenzene                  | 20    | 19.5   | ug/L | 98  | 0   |      | 82  | 118         | 20  |
|               | Ethyl Benzene                  | 20    | 19.7   | ug/L | 99  | 0   |      | 79  | 121         | 20  |
|               | m/p-Xylenes                    | 40    | 39.6   | ug/L | 99  | 0   |      | 80  | 121         | 20  |
|               | o-Xylene                       | 20    | 20.0   | ug/L | 100 | 2   |      | 78  | 122         | 20  |
|               | Styrene                        | 20    | 20.5   | ug/L | 103 | 0   |      | 78  | 123         | 20  |
|               | Bromoform                      | 20    | 19.9   | ug/L | 100 | 4   |      | 66  | 130         | 20  |
|               | Isopropylbenzene               | 20    | 20.1   | ug/L | 101 | 0   |      | 72  | 131         | 20  |
|               | 1,1,2,2-Tetrachloroethane      | 20    | 20.6   | ug/L | 103 | 4   |      | 71  | 121         | 20  |
|               | 1,3-Dichlorobenzene            | 20    | 19.5   | ug/L | 98  | 1   |      | 80  | 119         | 20  |
|               | 1,4-Dichlorobenzene            | 20    | 18.8   | ug/L | 94  | 3   |      | 79  | 118         | 20  |
|               | 1,2-Dichlorobenzene            | 20    | 20.5   | ug/L | 103 | 1   |      | 80  | 119         | 20  |

VOLATILE METHOD BLANK SUMMARY

EPA SAMPLE NO.

VX0515WBL01

Lab Name: CHEMTECH

Contract: TETR06

Lab Code: CHEM Case No.: Q2050

SAS No.: Q2050 SDG NO.: Q2050

Lab File ID: VX046202.D

Lab Sample ID: VX0515WBL01

Date Analyzed: 05/15/2025

Time Analyzed: 10:38

GC Column: DB-624UI ID: 0.18 (mm)

Heated Purge: (Y/N) N

Instrument ID: MSVOA\_X

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

| EPA<br>SAMPLE NO.       | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED |
|-------------------------|------------------|----------------|------------------|
| VX0515WBS01             | VX0515WBS01      | VX046206.D     | 05/15/2025       |
| VX0515WBSD01            | VX0515WBSD01     | VX046207.D     | 05/15/2025       |
| BP-VPB-182-GW-60-62     | Q2050-02         | VX046208.D     | 05/15/2025       |
| BP-VPB-182-TB-20250512  | Q2050-01         | VX046209.D     | 05/15/2025       |
| BP-VPB-182-EB-20250514  | Q2050-08         | VX046210.D     | 05/15/2025       |
| BP-VPB-182-GW-100-102   | Q2050-03         | VX046211.D     | 05/15/2025       |
| BP-VPB-182-GW-150-152   | Q2050-04         | VX046212.D     | 05/15/2025       |
| BP-VPB-182-GW-205-207   | Q2050-05         | VX046213.D     | 05/15/2025       |
| BP-VPB-182-GW-220-222   | Q2050-06         | VX046214.D     | 05/15/2025       |
| BP-VPB-182-DUP-20250514 | Q2050-07         | VX046215.D     | 05/15/2025       |
| BP-VPB-182-GW-240-242   | Q2050-09         | VX046216.D     | 05/15/2025       |

COMMENTS: \_\_\_\_\_

\_\_\_\_\_

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
BROMOFLUOROBENZENE (BFB)

Lab Name: CHEMTECH Contract: TETR06  
Lab Code: CHEM Case No.: Q2050 SAS No.: Q2050 SDG NO.: Q2050  
Lab File ID: VX046038.D BFB Injection Date: 05/05/2025  
Instrument ID: MSVOA\_X BFB Injection Time: 09:37  
GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: Y/N N

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 50  | 15.0 - 40.0% of mass 95            | 22.1                 |
| 75  | 30.0 - 60.0% of mass 95            | 56.2                 |
| 95  | Base Peak, 100% relative abundance | 100                  |
| 96  | 5.0 - 9.0% of mass 95              | 6.4                  |
| 173 | Less than 2.0% of mass 174         | 0.5 ( 0.7 ) 1        |
| 174 | 50.0 - 100.0% of mass 95           | 68.8                 |
| 175 | 5.0 - 9.0% of mass 174             | 5 ( 7.3 ) 1          |
| 176 | 95.0 - 101.0% of mass 174          | 66.7 ( 97 ) 1        |
| 177 | 5.0 - 9.0% of mass 176             | 4.6 ( 6.9 ) 2        |

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED | TIME<br>ANALYZED |
|-------------------|------------------|----------------|------------------|------------------|
| VSTDIC020         | VSTDIC020        | VX046041.D     | 05/05/2025       | 11:35            |
| VSTDIC050         | VSTDIC050        | VX046042.D     | 05/05/2025       | 11:58            |
| VSTDIC100         | VSTDIC100        | VX046043.D     | 05/05/2025       | 12:21            |
| VSTDIC150         | VSTDIC150        | VX046044.D     | 05/05/2025       | 12:45            |
| VSTDIC005         | VSTDIC005        | VX046046.D     | 05/05/2025       | 16:04            |
| VSTDIC001         | VSTDIC001        | VX046047.D     | 05/05/2025       | 16:27            |

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
BROMOFLUOROBENZENE (BFB)

Lab Name: CHEMTECH Contract: TETR06  
Lab Code: CHEM Case No.: Q2050 SAS No.: Q2050 SDG NO.: Q2050  
Lab File ID: VX046199.D BFB Injection Date: 05/15/2025  
Instrument ID: MSVOA\_X BFB Injection Time: 08:47  
GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: Y/N N

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 50  | 15.0 - 40.0% of mass 95            | 21.9                 |
| 75  | 30.0 - 60.0% of mass 95            | 58.6                 |
| 95  | Base Peak, 100% relative abundance | 100                  |
| 96  | 5.0 - 9.0% of mass 95              | 6.8                  |
| 173 | Less than 2.0% of mass 174         | 0.7 ( 1.1 ) 1        |
| 174 | 50.0 - 100.0% of mass 95           | 67.5                 |
| 175 | 5.0 - 9.0% of mass 174             | 5.3 ( 7.8 ) 1        |
| 176 | 95.0 - 101.0% of mass 174          | 65.6 ( 97.2 ) 1      |
| 177 | 5.0 - 9.0% of mass 176             | 4 ( 6.2 ) 2          |

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| EPA<br>SAMPLE NO.       | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED | TIME<br>ANALYZED |
|-------------------------|------------------|----------------|------------------|------------------|
| VSTDCCC050              | VSTDCCC050       | VX046200.D     | 05/15/2025       | 09:52            |
| VX0515WBL01             | VX0515WBL01      | VX046202.D     | 05/15/2025       | 10:38            |
| VX0515WBS01             | VX0515WBS01      | VX046206.D     | 05/15/2025       | 12:11            |
| VX0515WBSD01            | VX0515WBSD01     | VX046207.D     | 05/15/2025       | 12:38            |
| BP-VPB-182-GW-60-62     | Q2050-02         | VX046208.D     | 05/15/2025       | 13:01            |
| BP-VPB-182-TB-20250512  | Q2050-01         | VX046209.D     | 05/15/2025       | 13:24            |
| BP-VPB-182-EB-20250514  | Q2050-08         | VX046210.D     | 05/15/2025       | 13:47            |
| BP-VPB-182-GW-100-102   | Q2050-03         | VX046211.D     | 05/15/2025       | 14:11            |
| BP-VPB-182-GW-150-152   | Q2050-04         | VX046212.D     | 05/15/2025       | 14:34            |
| BP-VPB-182-GW-205-207   | Q2050-05         | VX046213.D     | 05/15/2025       | 14:58            |
| BP-VPB-182-GW-220-222   | Q2050-06         | VX046214.D     | 05/15/2025       | 15:21            |
| BP-VPB-182-DUP-20250514 | Q2050-07         | VX046215.D     | 05/15/2025       | 15:44            |
| BP-VPB-182-GW-240-242   | Q2050-09         | VX046216.D     | 05/15/2025       | 16:08            |
| VSTDCCC050EC            | VSTDCCC050       | VX046226.D     | 05/15/2025       | 20:01            |

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH Contract: TETR06  
Lab Code: CHEM Case No.: Q2050 SAS No.: Q2050 SDG NO.: Q2050  
Lab File ID: VX046200.D Date Analyzed: 05/15/2025  
Instrument ID: MSVOA\_X Time Analyzed: 09:52  
GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: (Y/N) N

|                         | IS1<br>AREA # | RT #  | IS2<br>AREA # | RT #  | IS3<br>AREA # | RT #   |
|-------------------------|---------------|-------|---------------|-------|---------------|--------|
| 12 HOUR STD             | 79845         | 5.54  | 140483        | 6.75  | 127630        | 10.05  |
| UPPER LIMIT             | 159690        | 6.044 | 280966        | 7.251 | 255260        | 10.549 |
| LOWER LIMIT             | 39922.5       | 5.044 | 70241.5       | 6.251 | 63815         | 9.549  |
| EPA SAMPLE NO.          |               |       |               |       |               |        |
| BP-VPB-182-TB-20250512  | 62722         | 5.54  | 124979        | 6.76  | 119897        | 10.05  |
| BP-VPB-182-GW-60-62     | 61866         | 5.55  | 123143        | 6.76  | 115076        | 10.06  |
| BP-VPB-182-GW-100-102   | 61221         | 5.55  | 120798        | 6.76  | 115485        | 10.06  |
| BP-VPB-182-GW-150-152   | 63694         | 5.55  | 128038        | 6.76  | 119578        | 10.05  |
| BP-VPB-182-GW-205-207   | 60533         | 5.55  | 121144        | 6.76  | 113426        | 10.06  |
| BP-VPB-182-GW-220-222   | 60707         | 5.55  | 121001        | 6.76  | 112446        | 10.06  |
| BP-VPB-182-DUP-20250514 | 61073         | 5.54  | 122275        | 6.76  | 116802        | 10.05  |
| BP-VPB-182-EB-20250514  | 61610         | 5.54  | 121862        | 6.76  | 116490        | 10.05  |
| BP-VPB-182-GW-240-242   | 62386         | 5.55  | 125278        | 6.76  | 120203        | 10.06  |
| VX0515WBL01             | 65765         | 5.54  | 130899        | 6.76  | 122283        | 10.05  |
| VX0515WBS01             | 89041         | 5.54  | 157824        | 6.76  | 139362        | 10.05  |
| VX0515WBSD01            | 86433         | 5.54  | 151750        | 6.76  | 132818        | 10.05  |

IS1 = Pentafluorobenzene  
IS2 = 1,4-Difluorobenzene  
IS3 = Chlorobenzene-d5

AREA UPPER LIMIT = +100% of internal standard area  
AREA LOWER LIMIT = -50% of internal standard area  
RT UPPER LIMIT = +0.50 minutes of internal standard RT  
RT LOWER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.  
\* Values outside of QC limits.

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH Contract: TETR06  
Lab Code: CHEM Case No.: Q2050 SAS No.: Q2050 SDG NO.: Q2050  
Lab File ID: VX046200.D Date Analyzed: 05/15/2025  
Instrument ID: MSVOA\_X Time Analyzed: 09:52  
GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: (Y/N) N

|                         | IS4<br>AREA # | RT #   |  |  |  |  |
|-------------------------|---------------|--------|--|--|--|--|
| 12 HOUR STD             | 62876         | 12.018 |  |  |  |  |
| UPPER LIMIT             | 125752        | 12.518 |  |  |  |  |
| LOWER LIMIT             | 31438         | 11.518 |  |  |  |  |
| EPA SAMPLE NO.          |               |        |  |  |  |  |
| BP-VPB-182-TB-20250512  | 51079         | 12.02  |  |  |  |  |
| BP-VPB-182-GW-60-62     | 50521         | 12.02  |  |  |  |  |
| BP-VPB-182-GW-100-102   | 47647         | 12.02  |  |  |  |  |
| BP-VPB-182-GW-150-152   | 51539         | 12.02  |  |  |  |  |
| BP-VPB-182-GW-205-207   | 47407         | 12.02  |  |  |  |  |
| BP-VPB-182-GW-220-222   | 44995         | 12.02  |  |  |  |  |
| BP-VPB-182-DUP-20250514 | 50006         | 12.02  |  |  |  |  |
| BP-VPB-182-EB-20250514  | 50288         | 12.02  |  |  |  |  |
| BP-VPB-182-GW-240-242   | 50190         | 12.02  |  |  |  |  |
| VX0515WBL01             | 51360         | 12.02  |  |  |  |  |
| VX0515WBS01             | 65777         | 12.02  |  |  |  |  |
| VX0515WBSD01            | 63406         | 12.02  |  |  |  |  |

IS4 = 1,4-Dichlorobenzene-d4

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT LOWER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.



# QC SAMPLE DATA



## Report of Analysis

|                    |                               |           |                 |              |
|--------------------|-------------------------------|-----------|-----------------|--------------|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: |              |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  |              |
| Client Sample ID:  | VX0515WBL01                   |           | SDG No.:        | Q2050        |
| Lab Sample ID:     | VX0515WBL01                   |           | Matrix:         | Water        |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |
| Prep Method :      |                               |           |                 |              |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046202.D        | 1         |           | 05/15/25 10:38 | VX051525      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOD  | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |      |            |       |
| 74-87-3        | Chloromethane                  | 0.50  | U         | 0.32  | 0.50 | 1.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 0.75  | U         | 0.26  | 0.75 | 1.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 3.80  | U         | 1.40  | 3.80 | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 0.75  | U         | 0.47  | 0.75 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 0.50  | U         | 0.33  | 0.50 | 1.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 0.75  | U         | 0.23  | 0.75 | 1.00       | ug/L  |
| 67-64-1        | Acetone                        | 3.80  | U         | 1.50  | 3.80 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 0.75  | U         | 0.21  | 0.75 | 1.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 0.50  | U         | 0.28  | 0.50 | 1.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 0.50  | U         | 0.23  | 0.50 | 1.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 2.50  | U         | 0.98  | 2.50 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 0.75  | U         | 0.19  | 0.75 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                     | 0.50  | U         | 0.25  | 0.50 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 71-43-2        | Benzene                        | 0.50  | U         | 0.15  | 0.50 | 1.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 0.75  | U         | 0.090 | 0.75 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 0.50  | U         | 0.20  | 0.50 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 0.50  | U         | 0.22  | 0.50 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 2.50  | U         | 0.68  | 2.50 | 5.00       | ug/L  |
| 108-88-3       | Toluene                        | 0.50  | U         | 0.14  | 0.50 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene          | 0.50  | U         | 0.17  | 0.50 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene        | 0.50  | U         | 0.16  | 0.50 | 1.00       | ug/L  |
| 79-00-5        | 1,1,2-Trichloroethane          | 0.50  | U         | 0.21  | 0.50 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                     | 2.50  | U         | 0.89  | 2.50 | 5.00       | ug/L  |

## Report of Analysis

|                    |                               |           |                 |              |
|--------------------|-------------------------------|-----------|-----------------|--------------|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: |              |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  |              |
| Client Sample ID:  | VX0515WBL01                   |           | SDG No.:        | Q2050        |
| Lab Sample ID:     | VX0515WBL01                   |           | Matrix:         | Water        |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |
| Prep Method :      |                               |           |                 |              |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046202.D        | 1         |           | 05/15/25 10:38 | VX051525      |

| CAS Number                | Parameter                 | Conc.  | Qualifier | MDL      | LOD  | LOQ / CRQL | Units   |
|---------------------------|---------------------------|--------|-----------|----------|------|------------|---------|
| 124-48-1                  | Dibromochloromethane      | 0.50   | U         | 0.18     | 0.50 | 1.00       | ug/L    |
| 127-18-4                  | Tetrachloroethene         | 0.50   | U         | 0.23     | 0.50 | 1.00       | ug/L    |
| 108-90-7                  | Chlorobenzene             | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-41-4                  | Ethyl Benzene             | 0.50   | U         | 0.13     | 0.50 | 1.00       | ug/L    |
| 179601-23-1               | m/p-Xylenes               | 1.00   | U         | 0.24     | 1.00 | 2.00       | ug/L    |
| 95-47-6                   | o-Xylene                  | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-42-5                  | Styrene                   | 0.50   | U         | 0.15     | 0.50 | 1.00       | ug/L    |
| 75-25-2                   | Bromoform                 | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 98-82-8                   | Isopropylbenzene          | 0.50   | U         | 0.12     | 0.50 | 1.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane | 0.50   | U         | 0.26     | 0.50 | 1.00       | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene       | 0.50   | U         | 0.19     | 0.50 | 1.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene       | 0.50   | U         | 0.16     | 0.50 | 1.00       | ug/L    |
| <b>SURROGATES</b>         |                           |        |           |          |      |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4     | 54.2   |           | 81 - 118 |      | 108%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane      | 51.5   |           | 80 - 119 |      | 103%       | SPK: 50 |
| 2037-26-5                 | Toluene-d8                | 50.2   |           | 89 - 112 |      | 100%       | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene      | 48.9   |           | 85 - 114 |      | 98%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                           |        |           |          |      |            |         |
| 363-72-4                  | Pentafluorobenzene        | 65800  | 5.544     |          |      |            |         |
| 540-36-3                  | 1,4-Difluorobenzene       | 131000 | 6.757     |          |      |            |         |
| 3114-55-4                 | Chlorobenzene-d5          | 122000 | 10.049    |          |      |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4    | 51400  | 12.018    |          |      |            |         |

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

## Report of Analysis

|                    |                               |           |                 |              |
|--------------------|-------------------------------|-----------|-----------------|--------------|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: |              |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  |              |
| Client Sample ID:  | VX0515WBS01                   |           | SDG No.:        | Q2050        |
| Lab Sample ID:     | VX0515WBS01                   |           | Matrix:         | Water        |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |
| Prep Method :      |                               |           |                 |              |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046206.D        | 1         |           | 05/15/25 12:11 | VX051525      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOD  | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |      |            |       |
| 74-87-3        | Chloromethane                  | 19.5  |           | 0.32  | 0.50 | 1.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 19.1  |           | 0.26  | 0.75 | 1.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 18.9  |           | 1.40  | 3.80 | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 22.5  |           | 0.47  | 0.75 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 20.4  |           | 0.33  | 0.50 | 1.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 19.9  |           | 0.25  | 0.50 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 19.0  |           | 0.23  | 0.75 | 1.00       | ug/L  |
| 67-64-1        | Acetone                        | 100   |           | 1.50  | 3.80 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 16.7  |           | 0.21  | 0.75 | 1.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 19.9  |           | 0.16  | 0.50 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 19.0  |           | 0.28  | 0.50 | 1.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 19.5  |           | 0.23  | 0.50 | 1.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 20.2  |           | 0.23  | 0.50 | 1.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 110   |           | 0.98  | 2.50 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 19.5  |           | 0.25  | 0.50 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 20.2  |           | 0.19  | 0.75 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                     | 21.0  |           | 0.25  | 0.50 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 20.2  |           | 0.20  | 0.50 | 1.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 18.3  |           | 0.16  | 0.50 | 1.00       | ug/L  |
| 71-43-2        | Benzene                        | 20.1  |           | 0.15  | 0.50 | 1.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 20.6  |           | 0.22  | 0.50 | 1.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 19.3  |           | 0.090 | 0.75 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 20.9  |           | 0.20  | 0.50 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 20.6  |           | 0.22  | 0.50 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 110   |           | 0.68  | 2.50 | 5.00       | ug/L  |
| 108-88-3       | Toluene                        | 20.0  |           | 0.14  | 0.50 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene          | 18.5  |           | 0.17  | 0.50 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene        | 19.5  |           | 0.16  | 0.50 | 1.00       | ug/L  |
| 79-00-5        | 1,1,2-Trichloroethane          | 20.7  |           | 0.21  | 0.50 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                     | 110   |           | 0.89  | 2.50 | 5.00       | ug/L  |

## Report of Analysis

|                    |                               |           |                 |              |
|--------------------|-------------------------------|-----------|-----------------|--------------|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: |              |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  |              |
| Client Sample ID:  | VX0515WBS01                   |           | SDG No.:        | Q2050        |
| Lab Sample ID:     | VX0515WBS01                   |           | Matrix:         | Water        |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |
| Prep Method :      |                               |           |                 |              |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046206.D        | 1         |           | 05/15/25 12:11 | VX051525      |

| CAS Number                | Parameter                 | Conc.  | Qualifier | MDL      | LOD  | LOQ / CRQL | Units   |
|---------------------------|---------------------------|--------|-----------|----------|------|------------|---------|
| 124-48-1                  | Dibromochloromethane      | 20.7   |           | 0.18     | 0.50 | 1.00       | ug/L    |
| 127-18-4                  | Tetrachloroethene         | 19.5   |           | 0.23     | 0.50 | 1.00       | ug/L    |
| 108-90-7                  | Chlorobenzene             | 19.5   |           | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-41-4                  | Ethyl Benzene             | 19.8   |           | 0.13     | 0.50 | 1.00       | ug/L    |
| 179601-23-1               | m/p-Xylenes               | 39.4   |           | 0.24     | 1.00 | 2.00       | ug/L    |
| 95-47-6                   | o-Xylene                  | 20.3   |           | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-42-5                  | Styrene                   | 20.5   |           | 0.15     | 0.50 | 1.00       | ug/L    |
| 75-25-2                   | Bromoform                 | 19.1   |           | 0.19     | 0.50 | 1.00       | ug/L    |
| 98-82-8                   | Isopropylbenzene          | 20.1   |           | 0.12     | 0.50 | 1.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane | 19.8   |           | 0.26     | 0.50 | 1.00       | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene       | 19.7   |           | 0.16     | 0.50 | 1.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene       | 19.3   |           | 0.19     | 0.50 | 1.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene       | 20.3   |           | 0.16     | 0.50 | 1.00       | ug/L    |
| <b>SURROGATES</b>         |                           |        |           |          |      |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4     | 50.7   |           | 81 - 118 |      | 101%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane      | 50.0   |           | 80 - 119 |      | 100%       | SPK: 50 |
| 2037-26-5                 | Toluene-d8                | 49.7   |           | 89 - 112 |      | 99%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene      | 48.9   |           | 85 - 114 |      | 98%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                           |        |           |          |      |            |         |
| 363-72-4                  | Pentafluorobenzene        | 89000  | 5.544     |          |      |            |         |
| 540-36-3                  | 1,4-Difluorobenzene       | 158000 | 6.757     |          |      |            |         |
| 3114-55-4                 | Chlorobenzene-d5          | 139000 | 10.049    |          |      |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4    | 65800  | 12.018    |          |      |            |         |

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

## Report of Analysis

|                    |                               |           |                 |              |
|--------------------|-------------------------------|-----------|-----------------|--------------|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: |              |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  |              |
| Client Sample ID:  | VX0515WBSD01                  |           | SDG No.:        | Q2050        |
| Lab Sample ID:     | VX0515WBSD01                  |           | Matrix:         | Water        |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |
| Prep Method :      |                               |           |                 |              |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046207.D        | 1         |           | 05/15/25 12:38 | VX051525      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOD  | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |      |            |       |
| 74-87-3        | Chloromethane                  | 19.3  |           | 0.32  | 0.50 | 1.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 18.2  |           | 0.26  | 0.75 | 1.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 17.8  |           | 1.40  | 3.80 | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 19.8  |           | 0.47  | 0.75 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 19.8  |           | 0.33  | 0.50 | 1.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 19.3  |           | 0.25  | 0.50 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 17.9  |           | 0.23  | 0.75 | 1.00       | ug/L  |
| 67-64-1        | Acetone                        | 100   |           | 1.50  | 3.80 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 16.1  |           | 0.21  | 0.75 | 1.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 20.1  |           | 0.16  | 0.50 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 18.7  |           | 0.28  | 0.50 | 1.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 19.0  |           | 0.23  | 0.50 | 1.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 20.0  |           | 0.23  | 0.50 | 1.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 110   |           | 0.98  | 2.50 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 19.2  |           | 0.25  | 0.50 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 19.3  |           | 0.19  | 0.75 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                     | 20.6  |           | 0.25  | 0.50 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 19.8  |           | 0.20  | 0.50 | 1.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 17.9  |           | 0.16  | 0.50 | 1.00       | ug/L  |
| 71-43-2        | Benzene                        | 19.7  |           | 0.15  | 0.50 | 1.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 20.7  |           | 0.22  | 0.50 | 1.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 18.8  |           | 0.090 | 0.75 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 21.3  |           | 0.20  | 0.50 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 20.3  |           | 0.22  | 0.50 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 110   |           | 0.68  | 2.50 | 5.00       | ug/L  |
| 108-88-3       | Toluene                        | 20.1  |           | 0.14  | 0.50 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene          | 18.8  |           | 0.17  | 0.50 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene        | 19.6  |           | 0.16  | 0.50 | 1.00       | ug/L  |
| 79-00-5        | 1,1,2-Trichloroethane          | 21.5  |           | 0.21  | 0.50 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                     | 110   |           | 0.89  | 2.50 | 5.00       | ug/L  |

## Report of Analysis

|                    |                               |           |                 |              |
|--------------------|-------------------------------|-----------|-----------------|--------------|
| Client:            | Tetra Tech NUS, Inc.          |           | Date Collected: |              |
| Project:           | NWIRP Bethpage 112G08005-WE13 |           | Date Received:  |              |
| Client Sample ID:  | VX0515WBSD01                  |           | SDG No.:        | Q2050        |
| Lab Sample ID:     | VX0515WBSD01                  |           | Matrix:         | Water        |
| Analytical Method: | 8260D                         |           | % Solid:        | 0            |
| Sample Wt/Vol:     | 5                             | Units: mL | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  |                               | uL        | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI                      | ID : 0.18 | Level :         | LOW          |
| Prep Method :      |                               |           |                 |              |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VX046207.D        | 1         |           | 05/15/25 12:38 | VX051525      |

| CAS Number                | Parameter                 | Conc.  | Qualifier | MDL      | LOD  | LOQ / CRQL | Units   |
|---------------------------|---------------------------|--------|-----------|----------|------|------------|---------|
| 124-48-1                  | Dibromochloromethane      | 20.7   |           | 0.18     | 0.50 | 1.00       | ug/L    |
| 127-18-4                  | Tetrachloroethene         | 19.3   |           | 0.23     | 0.50 | 1.00       | ug/L    |
| 108-90-7                  | Chlorobenzene             | 19.5   |           | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-41-4                  | Ethyl Benzene             | 19.7   |           | 0.13     | 0.50 | 1.00       | ug/L    |
| 179601-23-1               | m/p-Xylenes               | 39.6   |           | 0.24     | 1.00 | 2.00       | ug/L    |
| 95-47-6                   | o-Xylene                  | 20.0   |           | 0.12     | 0.50 | 1.00       | ug/L    |
| 100-42-5                  | Styrene                   | 20.5   |           | 0.15     | 0.50 | 1.00       | ug/L    |
| 75-25-2                   | Bromoform                 | 19.9   |           | 0.19     | 0.50 | 1.00       | ug/L    |
| 98-82-8                   | Isopropylbenzene          | 20.1   |           | 0.12     | 0.50 | 1.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane | 20.6   |           | 0.26     | 0.50 | 1.00       | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene       | 19.5   |           | 0.16     | 0.50 | 1.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene       | 18.8   |           | 0.19     | 0.50 | 1.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene       | 20.5   |           | 0.16     | 0.50 | 1.00       | ug/L    |
| <b>SURROGATES</b>         |                           |        |           |          |      |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4     | 49.4   |           | 81 - 118 |      | 99%        | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane      | 49.3   |           | 80 - 119 |      | 99%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                | 48.5   |           | 89 - 112 |      | 97%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene      | 47.9   |           | 85 - 114 |      | 96%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                           |        |           |          |      |            |         |
| 363-72-4                  | Pentafluorobenzene        | 86400  | 5.544     |          |      |            |         |
| 540-36-3                  | 1,4-Difluorobenzene       | 152000 | 6.757     |          |      |            |         |
| 3114-55-4                 | Chlorobenzene-d5          | 133000 | 10.049    |          |      |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4    | 63400  | 12.018    |          |      |            |         |

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



# CALIBRATION SUMMARY

### VOLATILE ORGANICS INITIAL CALIBRATION DATA

|                     |                  |                      |                   |
|---------------------|------------------|----------------------|-------------------|
| Lab Name:           | <u>CHEMTECH</u>  | Contract:            | <u>TETR06</u>     |
| Lab Code:           | <u>CHEM</u>      | Case No.:            | <u>Q2050</u>      |
| Instrument ID:      | <u>MSVOA_X</u>   | SAS No.:             | <u>Q2050</u>      |
| Heated Purge: (Y/N) | <u>N</u>         | SDG No.:             | <u>Q2050</u>      |
| GC Column:          | <u>DB-624UI</u>  | Calibration Date(s): | <u>05/05/2025</u> |
| ID:                 | <u>0.18 (mm)</u> | Calibration Time(s): | <u>11:35</u>      |
|                     |                  |                      | <u>16:27</u>      |

| LAB FILE ID:                   | RRF020 = VX046041.D | RRF050 = VX046042.D | RRF100 = VX046043.D | RRF150 = VX046044.D | RRF005 = VX046046.D | RRF001 = VX046047.D |       |       |
|--------------------------------|---------------------|---------------------|---------------------|---------------------|---------------------|---------------------|-------|-------|
| COMPOUND                       | RRF020              | RRF050              | RRF100              | RRF150              | RRF005              | RRF001              | RRF   | % RSD |
| Chloromethane                  | 0.727               | 0.775               | 0.787               | 0.791               | 0.679               | 0.694               | 0.742 | 6.6   |
| Vinyl Chloride                 | 0.660               | 0.710               | 0.727               | 0.755               | 0.619               | 0.673               | 0.691 | 7.2   |
| Bromomethane                   | 0.296               | 0.326               | 0.340               | 0.334               | 0.305               |                     | 0.320 | 5.8   |
| Chloroethane                   | 0.354               | 0.378               | 0.329               | 0.317               | 0.368               | 0.467               | 0.369 | 14.4  |
| Trichlorofluoromethane         | 1.035               | 1.068               | 0.983               | 0.985               | 0.990               | 1.064               | 1.021 | 3.9   |
| 1,1,2-Trichlorotrifluoroethane | 0.628               | 0.641               | 0.629               | 0.648               | 0.610               | 0.633               | 0.632 | 2.1   |
| 1,1-Dichloroethene             | 0.565               | 0.601               | 0.607               | 0.625               | 0.567               | 0.594               | 0.593 | 3.9   |
| Acetone                        | 0.361               | 0.362               | 0.361               | 0.370               | 0.408               | 0.380               | 0.374 | 4.9   |
| Carbon Disulfide               | 1.295               | 1.455               | 1.522               | 1.597               | 1.141               | 1.423               | 1.406 | 11.7  |
| Methyl tert-butyl Ether        | 2.044               | 2.160               | 2.172               | 2.239               | 1.908               | 1.949               | 2.079 | 6.4   |
| Methylene Chloride             | 0.689               | 0.684               | 0.691               | 0.691               | 0.689               | 0.853               | 0.716 | 9.4   |
| trans-1,2-Dichloroethene       | 0.573               | 0.610               | 0.612               | 0.622               | 0.557               | 0.604               | 0.596 | 4.3   |
| 1,1-Dichloroethane             | 1.233               | 1.263               | 1.263               | 1.286               | 1.154               | 1.116               | 1.219 | 5.6   |
| 2-Butanone                     | 0.540               | 0.555               | 0.558               | 0.569               | 0.539               | 0.495               | 0.543 | 4.8   |
| Carbon Tetrachloride           | 0.528               | 0.558               | 0.552               | 0.577               | 0.505               | 0.541               | 0.544 | 4.6   |
| cis-1,2-Dichloroethene         | 0.716               | 0.737               | 0.738               | 0.755               | 0.642               | 0.719               | 0.718 | 5.5   |
| Chloroform                     | 1.287               | 1.296               | 1.277               | 1.300               | 1.199               | 1.265               | 1.271 | 3     |
| 1,1,1-Trichloroethane          | 1.106               | 1.131               | 1.155               | 1.188               | 1.013               | 1.015               | 1.101 | 6.6   |
| Methylcyclohexane              | 0.596               | 0.641               | 0.627               | 0.658               | 0.587               | 0.627               | 0.623 | 4.3   |
| Benzene                        | 1.426               | 1.474               | 1.441               | 1.477               | 1.337               | 1.348               | 1.417 | 4.3   |
| 1,2-Dichloroethane             | 0.632               | 0.627               | 0.611               | 0.625               | 0.594               | 0.579               | 0.612 | 3.5   |
| Trichloroethene                | 0.344               | 0.355               | 0.345               | 0.362               | 0.315               | 0.324               | 0.341 | 5.3   |
| 1,2-Dichloropropane            | 0.356               | 0.371               | 0.368               | 0.378               | 0.324               | 0.317               | 0.352 | 7.4   |
| Bromodichloromethane           | 0.557               | 0.577               | 0.573               | 0.594               | 0.498               | 0.485               | 0.547 | 8.2   |
| 4-Methyl-2-Pentanone           | 0.620               | 0.634               | 0.630               | 0.631               | 0.555               | 0.561               | 0.605 | 6     |
| Toluene                        | 0.884               | 0.898               | 0.885               | 0.904               | 0.838               | 0.803               | 0.869 | 4.5   |
| t-1,3-Dichloropropene          | 0.468               | 0.528               | 0.555               | 0.591               | 0.406               | 0.371               | 0.487 | 17.9  |
| cis-1,3-Dichloropropene        | 0.531               | 0.578               | 0.602               | 0.623               | 0.469               | 0.423               | 0.538 | 14.6  |
| 1,1,2-Trichloroethane          | 0.349               | 0.354               | 0.351               | 0.356               | 0.337               | 0.308               | 0.343 | 5.3   |
| 2-Hexanone                     | 0.466               | 0.473               | 0.477               | 0.473               | 0.414               | 0.385               | 0.448 | 8.7   |

\* Compounds with required minimum RRF and maximum %RSD values.  
All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.



### VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: TETR06

Lab Code: CHEM Case No.: Q2050 SAS No.: Q2050 SDG No.: Q2050

Instrument ID: MSVOA\_X Calibration Date(s): 05/05/2025 05/05/2025

Heated Purge: (Y/N) N Calibration Time(s): 11:35 16:27

GC Column: DB-624UI ID: 0.18 (mm)

| LAB FILE ID:              |        | RRF020 = VX046041.D |        | RRF050 = VX046042.D |        | RRF100 = VX046043.D |       |       |  |
|---------------------------|--------|---------------------|--------|---------------------|--------|---------------------|-------|-------|--|
|                           |        | RRF150 = VX046044.D |        | RRF005 = VX046046.D |        | RRF001 = VX046047.D |       |       |  |
| COMPOUND                  | RRF020 | RRF050              | RRF100 | RRF150              | RRF005 | RRF001              | RRF   | % RSD |  |
| Dibromochloromethane      | 0.378  | 0.400               | 0.415  | 0.431               | 0.326  | 0.306               | 0.376 | 13.3  |  |
| Tetrachloroethene         | 0.390  | 0.375               | 0.345  | 0.344               | 0.323  | 0.347               | 0.354 | 6.8   |  |
| Chlorobenzene             | 1.093  | 1.098               | 1.085  | 1.114               | 1.046  | 1.131               | 1.094 | 2.7   |  |
| Ethyl Benzene             | 1.919  | 2.022               | 1.979  | 2.036               | 1.816  | 1.803               | 1.929 | 5.2   |  |
| m/p-Xylenes               | 0.706  | 0.740               | 0.721  | 0.740               | 0.678  | 0.648               | 0.706 | 5.2   |  |
| o-Xylene                  | 0.688  | 0.727               | 0.706  | 0.726               | 0.639  | 0.642               | 0.688 | 5.7   |  |
| Styrene                   | 1.135  | 1.219               | 1.214  | 1.230               | 1.012  | 0.951               | 1.127 | 10.6  |  |
| Bromoform                 | 0.270  | 0.304               | 0.312  | 0.327               | 0.236  | 0.234               | 0.281 | 14.2  |  |
| Isopropylbenzene          | 3.843  | 4.130               | 3.876  | 4.156               | 3.562  | 3.789               | 3.893 | 5.7   |  |
| 1,1,2,2-Tetrachloroethane | 1.315  | 1.338               | 1.284  | 1.345               | 1.350  | 1.552               | 1.364 | 7     |  |
| 1,3-Dichlorobenzene       | 1.633  | 1.701               | 1.656  | 1.730               | 1.558  | 1.619               | 1.649 | 3.7   |  |
| 1,4-Dichlorobenzene       | 1.629  | 1.693               | 1.639  | 1.722               | 1.606  | 1.817               | 1.684 | 4.6   |  |
| 1,2-Dichlorobenzene       | 1.613  | 1.696               | 1.634  | 1.702               | 1.577  | 1.710               | 1.655 | 3.3   |  |
| 1,2-Dichloroethane-d4     | 0.953  | 0.910               | 0.930  | 0.932               | 0.935  |                     | 0.932 | 1.6   |  |
| Dibromofluoromethane      | 0.359  | 0.355               | 0.364  | 0.368               | 0.354  |                     | 0.360 | 1.7   |  |
| Toluene-d8                | 1.246  | 1.223               | 1.266  | 1.275               | 1.221  |                     | 1.246 | 2     |  |
| 4-Bromofluorobenzene      | 0.455  | 0.470               | 0.500  | 0.500               | 0.464  |                     | 0.478 | 4.4   |  |

\* Compounds with required minimum RRF and maximum %RSD values.  
All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: TETR06

Lab Code: CHEM Case No.: Q2050 SAS No.: Q2050 SDG No.: Q2050

Instrument ID: MSVOA\_X Calibration Date/Time: 05/15/2025 09:52

Lab File ID: VX046200.D Init. Calib. Date(s): 05/05/2025 05/05/2025

Heated Purge: (Y/N) N Init. Calib. Time(s): 11:35 16:27

GC Column: DB-624UI ID: 0.18 (mm)

| COMPOUND                       | RRF   | RRF050 | MIN<br>RRF | %D     | MAX%D |
|--------------------------------|-------|--------|------------|--------|-------|
| Chloromethane                  | 0.742 | 0.714  | 0.1        | -3.77  | 20    |
| Vinyl Chloride                 | 0.691 | 0.611  |            | -11.58 | 20    |
| Bromomethane                   | 0.320 | 0.287  |            | -10.31 | 20    |
| Chloroethane                   | 0.369 | 0.363  |            | -1.63  | 20    |
| Trichlorofluoromethane         | 1.021 | 0.958  |            | -6.17  | 20    |
| 1,1,2-Trichlorotrifluoroethane | 0.632 | 0.579  |            | -8.39  | 20    |
| 1,1-Dichloroethene             | 0.593 | 0.532  |            | -10.29 | 20    |
| Acetone                        | 0.374 | 0.380  |            | 1.6    | 20    |
| Carbon Disulfide               | 1.406 | 1.195  |            | -15.01 | 20    |
| Methyl tert-butyl Ether        | 2.079 | 2.213  |            | 6.45   | 20    |
| Methylene Chloride             | 0.716 | 0.687  |            | -4.05  | 20    |
| trans-1,2-Dichloroethene       | 0.596 | 0.572  |            | -4.03  | 20    |
| 1,1-Dichloroethane             | 1.219 | 1.221  | 0.1        | 0.16   | 20    |
| 2-Butanone                     | 0.543 | 0.583  |            | 7.37   | 20    |
| Carbon Tetrachloride           | 0.544 | 0.510  |            | -6.25  | 20    |
| cis-1,2-Dichloroethene         | 0.718 | 0.727  |            | 1.25   | 20    |
| Chloroform                     | 1.271 | 1.299  |            | 2.2    | 20    |
| 1,1,1-Trichloroethane          | 1.101 | 1.079  |            | -2     | 20    |
| Methylcyclohexane              | 0.623 | 0.559  |            | -10.27 | 20    |
| Benzene                        | 1.417 | 1.399  |            | -1.27  | 20    |
| 1,2-Dichloroethane             | 0.612 | 0.634  |            | 3.6    | 20    |
| Trichloroethene                | 0.341 | 0.328  |            | -3.81  | 20    |
| 1,2-Dichloropropane            | 0.352 | 0.371  |            | 5.4    | 20    |
| Bromodichloromethane           | 0.547 | 0.585  |            | 6.95   | 20    |
| 4-Methyl-2-Pentanone           | 0.605 | 0.671  |            | 10.91  | 20    |
| Toluene                        | 0.869 | 0.871  |            | 0.23   | 20    |
| t-1,3-Dichloropropene          | 0.487 | 0.535  |            | 9.86   | 20    |
| cis-1,3-Dichloropropene        | 0.538 | 0.583  |            | 8.36   | 20    |
| 1,1,2-Trichloroethane          | 0.343 | 0.367  |            | 7      | 20    |
| 2-Hexanone                     | 0.448 | 0.501  |            | 11.83  | 20    |
| Dibromochloromethane           | 0.376 | 0.418  |            | 11.17  | 20    |
| Tetrachloroethene              | 0.354 | 0.326  |            | -7.91  | 20    |
| Chlorobenzene                  | 1.094 | 1.065  | 0.3        | -2.65  | 20    |
| Ethyl Benzene                  | 1.929 | 1.871  |            | -3.01  | 20    |
| m/p-Xylenes                    | 0.706 | 0.693  |            | -1.84  | 20    |
| o-Xylene                       | 0.688 | 0.686  |            | -0.29  | 20    |
| Styrene                        | 1.127 | 1.188  |            | 5.41   | 20    |
| Bromoform                      | 0.281 | 0.301  | 0.1        | 7.12   | 20    |

All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: TETR06

Lab Code: CHEM Case No.: Q2050 SAS No.: Q2050 SDG No.: Q2050

Instrument ID: MSVOA\_X Calibration Date/Time: 05/15/2025 09:52

Lab File ID: VX046200.D Init. Calib. Date(s): 05/05/2025 05/05/2025

Heated Purge: (Y/N) N Init. Calib. Time(s): 11:35 16:27

GC Column: DB-624UI ID: 0.18 (mm)

| COMPOUND                  | RRF   | RRF050 | MIN<br>RRF | %D    | MAX%D |
|---------------------------|-------|--------|------------|-------|-------|
| Isopropylbenzene          | 3.893 | 3.674  |            | -5.63 | 20    |
| 1,1,2,2-Tetrachloroethane | 1.364 | 1.329  | 0.3        | -2.57 | 20    |
| 1,3-Dichlorobenzene       | 1.649 | 1.630  |            | -1.15 | 20    |
| 1,4-Dichlorobenzene       | 1.684 | 1.628  |            | -3.33 | 20    |
| 1,2-Dichlorobenzene       | 1.655 | 1.656  |            | 0.06  | 20    |
| 1,2-Dichloroethane-d4     | 0.932 | 0.941  |            | 0.97  | 20    |
| Dibromofluoromethane      | 0.360 | 0.366  |            | 1.67  | 20    |
| Toluene-d8                | 1.246 | 1.215  |            | -2.49 | 20    |
| 4-Bromofluorobenzene      | 0.478 | 0.492  |            | 2.93  | 20    |

All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: TETR06

Lab Code: CHEM Case No.: Q2050 SAS No.: Q2050 SDG No.: Q2050

Instrument ID: MSVOA\_X Calibration Date/Time: 05/15/2025 20:01

Lab File ID: VX046226.D Init. Calib. Date(s): 05/05/2025 05/05/2025

Heated Purge: (Y/N) N Init. Calib. Time(s): 11:35 16:27

GC Column: DB-624UI ID: 0.18 (mm)

| COMPOUND                       | RRF   | RRF050 | MIN<br>RRF | %D    | MAX%D |
|--------------------------------|-------|--------|------------|-------|-------|
| Chloromethane                  | 0.742 | 0.751  | 0.1        | 1.21  | 50    |
| Vinyl Chloride                 | 0.691 | 0.688  |            | -0.43 | 50    |
| Bromomethane                   | 0.320 | 0.304  |            | -5    | 50    |
| Chloroethane                   | 0.369 | 0.380  |            | 2.98  | 50    |
| Trichlorofluoromethane         | 1.021 | 1.042  |            | 2.06  | 50    |
| 1,1,2-Trichlorotrifluoroethane | 0.632 | 0.632  |            | 0     | 50    |
| 1,1-Dichloroethene             | 0.593 | 0.583  |            | -1.69 | 50    |
| Acetone                        | 0.374 | 0.390  |            | 4.28  | 50    |
| Carbon Disulfide               | 1.406 | 1.298  |            | -7.68 | 50    |
| Methyl tert-butyl Ether        | 2.079 | 2.200  |            | 5.82  | 50    |
| Methylene Chloride             | 0.716 | 0.691  |            | -3.49 | 50    |
| trans-1,2-Dichloroethene       | 0.596 | 0.592  |            | -0.67 | 50    |
| 1,1-Dichloroethane             | 1.219 | 1.261  | 0.1        | 3.44  | 50    |
| 2-Butanone                     | 0.543 | 0.593  |            | 9.21  | 50    |
| Carbon Tetrachloride           | 0.544 | 0.552  |            | 1.47  | 50    |
| cis-1,2-Dichloroethene         | 0.718 | 0.724  |            | 0.84  | 50    |
| Chloroform                     | 1.271 | 1.306  |            | 2.75  | 50    |
| 1,1,1-Trichloroethane          | 1.101 | 1.149  |            | 4.36  | 50    |
| Methylcyclohexane              | 0.623 | 0.600  |            | -3.69 | 50    |
| Benzene                        | 1.417 | 1.453  |            | 2.54  | 50    |
| 1,2-Dichloroethane             | 0.612 | 0.624  |            | 1.96  | 50    |
| Trichloroethene                | 0.341 | 0.339  |            | -0.59 | 50    |
| 1,2-Dichloropropane            | 0.352 | 0.379  |            | 7.67  | 50    |
| Bromodichloromethane           | 0.547 | 0.579  |            | 5.85  | 50    |
| 4-Methyl-2-Pentanone           | 0.605 | 0.676  |            | 11.74 | 50    |
| Toluene                        | 0.869 | 0.895  |            | 2.99  | 50    |
| t-1,3-Dichloropropene          | 0.487 | 0.511  |            | 4.93  | 50    |
| cis-1,3-Dichloropropene        | 0.538 | 0.563  |            | 4.65  | 50    |
| 1,1,2-Trichloroethane          | 0.343 | 0.367  |            | 7     | 50    |
| 2-Hexanone                     | 0.448 | 0.498  |            | 11.16 | 50    |
| Dibromochloromethane           | 0.376 | 0.408  |            | 8.51  | 50    |
| Tetrachloroethene              | 0.354 | 0.324  |            | -8.48 | 50    |
| Chlorobenzene                  | 1.094 | 1.101  | 0.3        | 0.64  | 50    |
| Ethyl Benzene                  | 1.929 | 2.000  |            | 3.68  | 50    |
| m/p-Xylenes                    | 0.706 | 0.733  |            | 3.82  | 50    |
| o-Xylene                       | 0.688 | 0.721  |            | 4.8   | 50    |
| Styrene                        | 1.127 | 1.224  |            | 8.61  | 50    |
| Bromoform                      | 0.281 | 0.290  | 0.1        | 3.2   | 50    |

All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: TETR06

Lab Code: CHEM Case No.: Q2050 SAS No.: Q2050 SDG No.: Q2050

Instrument ID: MSVOA\_X Calibration Date/Time: 05/15/2025 20:01

Lab File ID: VX046226.D Init. Calib. Date(s): 05/05/2025 05/05/2025

Heated Purge: (Y/N) N Init. Calib. Time(s): 11:35 16:27

GC Column: DB-624UI ID: 0.18 (mm)

| COMPOUND                  | RRF   | RRF050 | MIN<br>RRF | %D    | MAX%D |
|---------------------------|-------|--------|------------|-------|-------|
| Isopropylbenzene          | 3.893 | 4.011  |            | 3.03  | 50    |
| 1,1,2,2-Tetrachloroethane | 1.364 | 1.352  | 0.3        | -0.88 | 50    |
| 1,3-Dichlorobenzene       | 1.649 | 1.651  |            | 0.12  | 50    |
| 1,4-Dichlorobenzene       | 1.684 | 1.597  |            | -5.17 | 50    |
| 1,2-Dichlorobenzene       | 1.655 | 1.632  |            | -1.39 | 50    |
| 1,2-Dichloroethane-d4     | 0.932 | 0.923  |            | -0.97 | 50    |
| Dibromofluoromethane      | 0.360 | 0.369  |            | 2.5   | 50    |
| Toluene-d8                | 1.246 | 1.242  |            | -0.32 | 50    |
| 4-Bromofluorobenzene      | 0.478 | 0.482  |            | 0.84  | 50    |

All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.

## LAB CHRONICLE

|                 |                      |                   |                               |
|-----------------|----------------------|-------------------|-------------------------------|
| <b>OrderID:</b> | Q2050                | <b>OrderDate:</b> | 5/15/2025 9:46:00 AM          |
| <b>Client:</b>  | Tetra Tech NUS, Inc. | <b>Project:</b>   | NWIRP Bethpage 112G08005-WE13 |
| <b>Contact:</b> | Ernie Wu             | <b>Location:</b>  | L31,VOA Ref. #3 Water         |

| LabID           | ClientID                           | Matrix       | Test           | Method        | Sample Date     | Prep Date | Anal Date | Received        |
|-----------------|------------------------------------|--------------|----------------|---------------|-----------------|-----------|-----------|-----------------|
| <b>Q2050-02</b> | <b>BP-VPB-182-GW-60-6<br/>2</b>    | <b>Water</b> |                |               | <b>05/12/25</b> |           |           | <b>05/14/25</b> |
|                 |                                    |              | SVOC-SIMGroup1 | 8270-Modified |                 | 05/16/25  | 05/19/25  |                 |
| <b>Q2050-04</b> | <b>BP-VPB-182-GW-150-<br/>152</b>  | <b>Water</b> |                |               | <b>05/13/25</b> |           |           | <b>05/14/25</b> |
|                 |                                    |              | SVOC-SIMGroup1 | 8270-Modified |                 | 05/16/25  | 05/19/25  |                 |
| <b>Q2050-06</b> | <b>BP-VPB-182-GW-220-<br/>222</b>  | <b>Water</b> |                |               | <b>05/14/25</b> |           |           | <b>05/14/25</b> |
|                 |                                    |              | SVOC-SIMGroup1 | 8270-Modified |                 | 05/16/25  | 05/19/25  |                 |
| <b>Q2050-08</b> | <b>BP-VPB-182-EB-2025<br/>0514</b> | <b>Water</b> |                |               | <b>05/14/25</b> |           |           | <b>05/14/25</b> |
|                 |                                    |              | SVOC-SIMGroup1 | 8270-Modified |                 | 05/16/25  | 05/19/25  |                 |



284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,  
Fax : 908 789 8922

### Hit Summary Sheet SW-846

**SDG No.:** Q2050  
**Client:** Tetra Tech NUS, Inc.

| Sample ID          | Client ID                    | Parameter                   | Concentration | C           | MDL  | LOD  | RDL  | Units |
|--------------------|------------------------------|-----------------------------|---------------|-------------|------|------|------|-------|
| <b>Client ID :</b> | <b>BP-VPB-182-GW-220-222</b> |                             |               |             |      |      |      |       |
| Q2050-06           | BP-VPB-182-GW-220-22 WATER   | 1,4-Dioxane                 | 0.130         | J           | 0.08 | 0.23 | 0.23 | ug/L  |
|                    |                              | <b>Total Svoc :</b>         |               | <b>0.13</b> |      |      |      |       |
|                    |                              | <b>Total Concentration:</b> |               | <b>0.13</b> |      |      |      |       |



# SAMPLE DATA



## Report of Analysis

|                    |                               |                 |                |
|--------------------|-------------------------------|-----------------|----------------|
| Client:            | Tetra Tech NUS, Inc.          | Date Collected: | 05/12/25       |
| Project:           | NWIRP Bethpage 112G08005-WE13 | Date Received:  | 05/14/25       |
| Client Sample ID:  | BP-VPB-182-GW-60-62           | SDG No.:        | Q2050          |
| Lab Sample ID:     | Q2050-02                      | Matrix:         | Water          |
| Analytical Method: | SW8270ESIM                    | % Solid:        | 0              |
| Sample Wt/Vol:     | 550 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 :      |                               |                 |                |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BN037048.D        | 1         | 05/16/25 08:42 | 05/19/25 16:51 | PB168032      |

| CAS Number                | Parameter               | Conc. | Qualifier | MDL      | LOD  | LOQ / CRQL | Units    |
|---------------------------|-------------------------|-------|-----------|----------|------|------------|----------|
| <b>TARGETS</b>            |                         |       |           |          |      |            |          |
| 123-91-1                  | 1,4-Dioxane             | 0.36  | U         | 0.12     | 0.36 | 0.36       | ug/L     |
| <b>SURROGATES</b>         |                         |       |           |          |      |            |          |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.24  |           | 30 - 150 |      | 60%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.39  |           | 30 - 150 |      | 98%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.26  |           | 55 - 111 |      | 65%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.28  |           | 53 - 106 |      | 71%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.52  |           | 58 - 132 |      | 129%       | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 2090  | 7.611     |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 5700  | 10.394    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 3310  | 14.256    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 6660  | 17.009    |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 6100  | 21.206    |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 5200  | 23.409    |          |      |            |          |

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

## Report of Analysis

|                    |                               |                 |                |
|--------------------|-------------------------------|-----------------|----------------|
| Client:            | Tetra Tech NUS, Inc.          | Date Collected: | 05/13/25       |
| Project:           | NWIRP Bethpage 112G08005-WE13 | Date Received:  | 05/14/25       |
| Client Sample ID:  | BP-VPB-182-GW-150-152         | SDG No.:        | Q2050          |
| Lab Sample ID:     | Q2050-04                      | Matrix:         | Water          |
| Analytical Method: | SW8270ESIM                    | % Solid:        | 0              |
| Sample Wt/Vol:     | 910 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 :      |                               |                 |                |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BN037049.D        | 1         | 05/16/25 08:42 | 05/19/25 17:27 | PB168032      |

| CAS Number                | Parameter               | Conc. | Qualifier | MDL      | LOD  | LOQ / CRQL | Units    |
|---------------------------|-------------------------|-------|-----------|----------|------|------------|----------|
| <b>TARGETS</b>            |                         |       |           |          |      |            |          |
| 123-91-1                  | 1,4-Dioxane             | 0.22  | U         | 0.070    | 0.22 | 0.22       | ug/L     |
| <b>SURROGATES</b>         |                         |       |           |          |      |            |          |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.28  |           | 30 - 150 |      | 69%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.37  |           | 30 - 150 |      | 93%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.28  |           | 55 - 111 |      | 71%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.33  |           | 53 - 106 |      | 83%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.58  | *         | 58 - 132 |      | 145%       | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 2170  |           | 7.611    |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 5840  |           | 10.394   |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 3380  |           | 14.267   |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 6630  |           | 17.009   |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 5280  |           | 21.207   |      |            |          |
| 1520-96-3                 | Perylene-d12            | 4530  |           | 23.407   |      |            |          |

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\* = Values outside of QC limits

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() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                               |                 |                |
|--------------------|-------------------------------|-----------------|----------------|
| Client:            | Tetra Tech NUS, Inc.          | Date Collected: | 05/14/25       |
| Project:           | NWIRP Bethpage 112G08005-WE13 | Date Received:  | 05/14/25       |
| Client Sample ID:  | BP-VPB-182-GW-220-222         | SDG No.:        | Q2050          |
| Lab Sample ID:     | Q2050-06                      | Matrix:         | Water          |
| Analytical Method: | SW8270ESIM                    | % Solid:        | 0              |
| Sample Wt/Vol:     | 870 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 :      |                               |                 |                |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BN037050.D        | 1         | 05/16/25 08:42 | 05/19/25 18:03 | PB168032      |

| CAS Number                | Parameter               | Conc. | Qualifier | MDL      | LOD  | LOQ / CRQL | Units    |
|---------------------------|-------------------------|-------|-----------|----------|------|------------|----------|
| <b>TARGETS</b>            |                         |       |           |          |      |            |          |
| 123-91-1                  | 1,4-Dioxane             | 0.13  | J         | 0.080    | 0.23 | 0.23       | ug/L     |
| <b>SURROGATES</b>         |                         |       |           |          |      |            |          |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.27  |           | 30 - 150 |      | 68%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.36  |           | 30 - 150 |      | 90%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.27  |           | 55 - 111 |      | 68%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.32  |           | 53 - 106 |      | 80%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.49  |           | 58 - 132 |      | 122%       | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 1930  | 7.611     |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 5280  | 10.394    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 3090  | 14.256    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 6260  | 17.009    |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 4980  | 21.207    |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 4280  | 23.41     |          |      |            |          |

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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

## Report of Analysis

|                    |                               |                 |                |
|--------------------|-------------------------------|-----------------|----------------|
| Client:            | Tetra Tech NUS, Inc.          | Date Collected: | 05/14/25       |
| Project:           | NWIRP Bethpage 112G08005-WE13 | Date Received:  | 05/14/25       |
| Client Sample ID:  | BP-VPB-182-EB-20250514        | SDG No.:        | Q2050          |
| Lab Sample ID:     | Q2050-08                      | Matrix:         | Water          |
| Analytical Method: | SW8270ESIM                    | % Solid:        | 0              |
| Sample Wt/Vol:     | 950 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 :      |                               |                 |                |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BN037051.D        | 1         | 05/16/25 08:42 | 05/19/25 18:40 | PB168032      |

| 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 | 0.27  |           | 30 - 150 |      | 68%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.36  |           | 30 - 150 |      | 90%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.28  |           | 55 - 111 |      | 69%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.34  |           | 53 - 106 |      | 84%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.54  | *         | 58 - 132 |      | 134%       | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 1900  |           | 7.618    |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 5250  |           | 10.394   |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 3120  |           | 14.267   |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 6080  |           | 17.009   |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 4710  |           | 21.207   |      |            |          |
| 1520-96-3                 | Perylene-d12            | 4050  |           | 23.412   |      |            |          |

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# QC SUMMARY

## Surrogate Summary

SW-846

SDG No.: Q2050

Client: Tetra Tech NUS, Inc.

Analytical Method: 8270-Modified

| Lab Sample ID | Client ID              | Parameter               | Spike (PPM) | Result (PPM) | Recovery (%) | Qual | Limits (%) |      |
|---------------|------------------------|-------------------------|-------------|--------------|--------------|------|------------|------|
|               |                        |                         |             |              |              |      | Low        | High |
| PB168032BL    | PB168032BL             | 2-Methylnaphthalene-d10 | 0.4         | 0.36         | 90           |      | 30         | 150  |
|               |                        | Fluoranthene-d10        | 0.4         | 0.34         | 84           |      | 30         | 150  |
|               |                        | Nitrobenzene-d5         | 0.4         | 0.32         | 81           |      | 55         | 111  |
|               |                        | 2-Fluorobiphenyl        | 0.4         | 0.34         | 84           |      | 53         | 106  |
|               |                        | Terphenyl-d14           | 0.4         | 0.42         | 104          |      | 58         | 132  |
| PB168032BS    | PB168032BS             | 2-Methylnaphthalene-d10 | 0.4         | 0.39         | 96           |      | 30         | 150  |
|               |                        | Fluoranthene-d10        | 0.4         | 0.33         | 81           |      | 30         | 150  |
|               |                        | Nitrobenzene-d5         | 0.4         | 0.32         | 81           |      | 55         | 111  |
|               |                        | 2-Fluorobiphenyl        | 0.4         | 0.33         | 83           |      | 53         | 106  |
|               |                        | Terphenyl-d14           | 0.4         | 0.39         | 97           |      | 58         | 132  |
| PB168032BSD   | PB168032BSD            | 2-Methylnaphthalene-d10 | 0.4         | 0.39         | 98           |      | 30         | 150  |
|               |                        | Fluoranthene-d10        | 0.4         | 0.33         | 83           |      | 30         | 150  |
|               |                        | Nitrobenzene-d5         | 0.4         | 0.35         | 88           |      | 55         | 111  |
|               |                        | 2-Fluorobiphenyl        | 0.4         | 0.33         | 83           |      | 53         | 106  |
|               |                        | Terphenyl-d14           | 0.4         | 0.42         | 104          |      | 58         | 132  |
| Q2050-02      | BP-VPB-182-GW-60-62    | 2-Methylnaphthalene-d10 | 0.4         | 0.24         | 60           |      | 30         | 150  |
|               |                        | Fluoranthene-d10        | 0.4         | 0.39         | 98           |      | 30         | 150  |
|               |                        | Nitrobenzene-d5         | 0.4         | 0.26         | 65           |      | 55         | 111  |
|               |                        | 2-Fluorobiphenyl        | 0.4         | 0.28         | 71           |      | 53         | 106  |
|               |                        | Terphenyl-d14           | 0.4         | 0.52         | 129          |      | 58         | 132  |
| Q2050-04      | BP-VPB-182-GW-150-152  | 2-Methylnaphthalene-d10 | 0.4         | 0.28         | 69           |      | 30         | 150  |
|               |                        | Fluoranthene-d10        | 0.4         | 0.37         | 93           |      | 30         | 150  |
|               |                        | Nitrobenzene-d5         | 0.4         | 0.28         | 71           |      | 55         | 111  |
|               |                        | 2-Fluorobiphenyl        | 0.4         | 0.33         | 83           |      | 53         | 106  |
|               |                        | Terphenyl-d14           | 0.4         | 0.58         | 145          | *    | 58         | 132  |
| Q2050-06      | BP-VPB-182-GW-220-222  | 2-Methylnaphthalene-d10 | 0.4         | 0.27         | 68           |      | 30         | 150  |
|               |                        | Fluoranthene-d10        | 0.4         | 0.36         | 90           |      | 30         | 150  |
|               |                        | Nitrobenzene-d5         | 0.4         | 0.27         | 68           |      | 55         | 111  |
|               |                        | 2-Fluorobiphenyl        | 0.4         | 0.32         | 80           |      | 53         | 106  |
|               |                        | Terphenyl-d14           | 0.4         | 0.49         | 122          |      | 58         | 132  |
| Q2050-08      | BP-VPB-182-EB-20250514 | 2-Methylnaphthalene-d10 | 0.4         | 0.27         | 68           |      | 30         | 150  |
|               |                        | Fluoranthene-d10        | 0.4         | 0.36         | 90           |      | 30         | 150  |
|               |                        | Nitrobenzene-d5         | 0.4         | 0.28         | 69           |      | 55         | 111  |
|               |                        | 2-Fluorobiphenyl        | 0.4         | 0.34         | 84           |      | 53         | 106  |
|               |                        | Terphenyl-d14           | 0.4         | 0.54         | 134          | *    | 58         | 132  |

**Laboratory Control Sample/Laboratory Control Sample Duplicate Summary**

SW-846

SDG No.: Q2050

Client: Tetra Tech NUS, Inc.

Analytical Method: 8270-Modified DataFile: BN037080.D

| Lab Sample ID | Parameter   | Spike | Result | Unit | Rec | RPD | Qual | RPD  | Low | Limits | RPD |
|---------------|-------------|-------|--------|------|-----|-----|------|------|-----|--------|-----|
|               |             |       |        |      |     |     |      | Qual |     | High   |     |
| PB168032BS    | 1,4-Dioxane | 0.4   | 0.31   | ug/L | 78  |     |      |      | 70  | 130    |     |

**Laboratory Control Sample/Laboratory Control Sample Duplicate Summary**

SW-846

**SDG No.:** Q2050

**Client:** Tetra Tech NUS, Inc.

**Analytical Method:** 8270-Modified **DataFile:** BN037081.D

| Lab Sample ID | Parameter   | Spike | Result | Unit | Rec | RPD | Qual | RPD  | Limits |      | RPD |
|---------------|-------------|-------|--------|------|-----|-----|------|------|--------|------|-----|
|               |             |       |        |      |     |     |      | Qual | Low    | High |     |
| PB168032BSD   | 1,4-Dioxane | 0.4   | 0.32   | ug/L | 80  | 3   |      |      | 70     | 130  | 20  |



4B

SEMIVOLATILE METHOD BLANK SUMMARY

EPA SAMPLE NO.

PB168032BL

Lab Name: CHEMTECH

Contract: TETR06

Lab Code: CHEM Case No.: Q2050

SAS No.: Q2050 SDG NO.: Q2050

Lab File ID: BN037075.D

Lab Sample ID: PB168032BL

Instrument ID: BNA\_N

Date Extracted: 05/16/2025

Matrix: (soil/water) Water

Date Analyzed: 05/20/2025

Level: (low/med) LOW

Time Analyzed: 15:45

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

| EPA<br>SAMPLE NO.      | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED |
|------------------------|------------------|----------------|------------------|
| PB168032BS             | PB168032BS       | BN037080.D     | 05/20/2025       |
| PB168032BSD            | PB168032BSD      | BN037081.D     | 05/20/2025       |
| BP-VPB-182-GW-60-62    | Q2050-02         | BN037048.D     | 05/19/2025       |
| BP-VPB-182-GW-150-152  | Q2050-04         | BN037049.D     | 05/19/2025       |
| BP-VPB-182-GW-220-222  | Q2050-06         | BN037050.D     | 05/19/2025       |
| BP-VPB-182-EB-20250514 | Q2050-08         | BN037051.D     | 05/19/2025       |

COMMENTS: \_\_\_\_\_

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: CHEMTECH

Contract: TETR06

Lab Code: CHEM

SAS No.: Q2050

SDG NO.: Q2050

Lab File ID: BN036998.D

DFTPP Injection Date: 05/13/2025

Instrument ID: BNA\_N

DFTPP Injection Time: 17:02

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 51  | 10.0 - 80.0% of mass 198           | 62.8                 |
| 68  | Less than 2.0% of mass 69          | 0.8 ( 1.4 ) 1        |
| 69  | Mass 69 relative abundance         | 55.6                 |
| 70  | Less than 2.0% of mass 69          | 0.3 ( 0.6 ) 1        |
| 127 | 10.0 - 80.0% of mass 198           | 52.7                 |
| 197 | Less than 2.0% of mass 198         | 0.0                  |
| 198 | Base Peak, 100% relative abundance | 100                  |
| 199 | 5.0 to 9.0% of mass 198            | 6.9                  |
| 275 | 10.0 - 60.0% of mass 198           | 23.8                 |
| 365 | Greater than 1% of mass 198        | 3.9                  |
| 441 | Present, but less than mass 443    | 8.7                  |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 10.4 ( 19 ) 2        |

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED | TIME<br>ANALYZED |
|-------------------|------------------|----------------|------------------|------------------|
| SSTDICC0.1        | SSTDICC0.1       | BN036999.D     | 05/13/2025       | 17:41            |
| SSTDICC0.2        | SSTDICC0.2       | BN037000.D     | 05/13/2025       | 18:17            |
| SSTDICCC0.4       | SSTDICCC0.4      | BN037001.D     | 05/13/2025       | 18:53            |
| SSTDICC0.8        | SSTDICC0.8       | BN037002.D     | 05/13/2025       | 19:29            |
| SSTDICC1.6        | SSTDICC1.6       | BN037003.D     | 05/13/2025       | 20:05            |
| SSTDICC3.2        | SSTDICC3.2       | BN037004.D     | 05/13/2025       | 20:41            |
| SSTDICC5.0        | SSTDICC5.0       | BN037005.D     | 05/13/2025       | 21:17            |

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: CHEMTECH

Contract: TETR06

Lab Code: CHEM

SAS No.: Q2050

SDG NO.: Q2050

Lab File ID: BN037037.D

DFTPP Injection Date: 05/19/2025

Instrument ID: BNA\_N

DFTPP Injection Time: 10:08

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 51  | 10.0 - 80.0% of mass 198           | 62.3                 |
| 68  | Less than 2.0% of mass 69          | 0.8 ( 1.4 ) 1        |
| 69  | Mass 69 relative abundance         | 55.6                 |
| 70  | Less than 2.0% of mass 69          | 0.4 ( 0.7 ) 1        |
| 127 | 10.0 - 80.0% of mass 198           | 51.1                 |
| 197 | Less than 2.0% of mass 198         | 0.0                  |
| 198 | Base Peak, 100% relative abundance | 100                  |
| 199 | 5.0 to 9.0% of mass 198            | 6.9                  |
| 275 | 10.0 - 60.0% of mass 198           | 24.7                 |
| 365 | Greater than 1% of mass 198        | 3.8                  |
| 441 | Present, but less than mass 443    | 8.8                  |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 10.1 ( 18 ) 2        |

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| EPA<br>SAMPLE NO.      | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED | TIME<br>ANALYZED |
|------------------------|------------------|----------------|------------------|------------------|
| SSTDCCC0.4             | SSTDCCC0.4       | BN037038.D     | 05/19/2025       | 10:47            |
| BP-VPB-182-GW-60-62    | Q2050-02         | BN037048.D     | 05/19/2025       | 16:51            |
| BP-VPB-182-GW-150-152  | Q2050-04         | BN037049.D     | 05/19/2025       | 17:27            |
| BP-VPB-182-GW-220-222  | Q2050-06         | BN037050.D     | 05/19/2025       | 18:03            |
| BP-VPB-182-EB-20250514 | Q2050-08         | BN037051.D     | 05/19/2025       | 18:40            |
| SSTDCCC0.4EC           | SSTDCCC0.4       | BN037055.D     | 05/19/2025       | 21:05            |

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: CHEMTECH

Contract: TETR06

Lab Code: CHEM

SAS No.: Q2050

SDG NO.: Q2050

Lab File ID: BN037073.D

DFTPP Injection Date: 05/20/2025

Instrument ID: BNA\_N

DFTPP Injection Time: 14:30

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 51  | 10.0 - 80.0% of mass 198           | 69.3                 |
| 68  | Less than 2.0% of mass 69          | 0.0 ( 0.0 ) 1        |
| 69  | Mass 69 relative abundance         | 59.3                 |
| 70  | Less than 2.0% of mass 69          | 0.3 ( 0.4 ) 1        |
| 127 | 10.0 - 80.0% of mass 198           | 52.8                 |
| 197 | Less than 2.0% of mass 198         | 0.2                  |
| 198 | Base Peak, 100% relative abundance | 100                  |
| 199 | 5.0 to 9.0% of mass 198            | 6.9                  |
| 275 | 10.0 - 60.0% of mass 198           | 25.4                 |
| 365 | Greater than 1% of mass 198        | 3.7                  |
| 441 | Present, but less than mass 443    | 8.8                  |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 10.5 (19.4) 2        |

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED | TIME<br>ANALYZED |
|-------------------|------------------|----------------|------------------|------------------|
| SSTDCCC0.4        | SSTDCCC0.4       | BN037074.D     | 05/20/2025       | 15:09            |
| PB168032BL        | PB168032BL       | BN037075.D     | 05/20/2025       | 15:45            |
| PB168032BS        | PB168032BS       | BN037080.D     | 05/20/2025       | 18:47            |
| PB168032BSD       | PB168032BSD      | BN037081.D     | 05/20/2025       | 19:23            |
| SSTDCCC0.4EC      | SSTDCCC0.4       | BN037082.D     | 05/20/2025       | 19:59            |

8B

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH

Lab Code: CHEM Case No.: Q2050 SAS No.: Q2050 SDG NO.: Q2050

EPA Sample No.: SSTDCCC0.4 Date Analyzed: 05/19/2025

Lab File ID: BN037038.D Time Analyzed: 10:47

Instrument ID: BNA\_N GC Column: ZB-GR ID: 0.25 (mm)

|                           | IS1 (DCB)<br>AREA # | RT #  | IS2 (NPT)<br>AREA # | RT #   | IS3 (ANT)<br>AREA # | RT #   |
|---------------------------|---------------------|-------|---------------------|--------|---------------------|--------|
| 12 HOUR STD               | 2152                | 7.611 | 5916                | 10.39  | 3512                | 14.26  |
| UPPER LIMIT               | 4304                | 8.111 | 11832               | 10.894 | 7024                | 14.756 |
| LOWER LIMIT               | 1076                | 7.111 | 2958                | 9.894  | 1756                | 13.756 |
| EPA SAMPLE NO.            |                     |       |                     |        |                     |        |
| 01 BP-VPB-182-GW-60-62    | 2086                | 7.61  | 5701                | 10.39  | 3309                | 14.26  |
| 02 BP-VPB-182-GW-150-152  | 2171                | 7.61  | 5837                | 10.39  | 3378                | 14.27  |
| 03 BP-VPB-182-GW-220-222  | 1931                | 7.61  | 5282                | 10.39  | 3090                | 14.26  |
| 04 BP-VPB-182-EB-20250514 | 1904                | 7.62  | 5247                | 10.39  | 3115                | 14.27  |

IS1 (DCB) = 1,4-Dichlorobenzene-d4

IS2 (NPT) = Naphthalene-d8

IS3 (ANT) = Acenaphthene-d10

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

8C

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH

Lab Code: CHEM Case No.: Q2050 SAS No.: Q2050 SDG NO.: Q2050

EPA Sample No.: SSTDCCC0.4 Date Analyzed: 05/19/2025

Lab File ID: BN037038.D Time Analyzed: 10:47

Instrument ID: BNA\_N GC Column: ZB-GR ID: 0.25 (mm)

|                           | IS4 (PHN)<br>AREA # | RT #   | IS5 (CRY)<br>AREA # | RT #   | IS6 (PRY)<br>AREA # | RT #  |
|---------------------------|---------------------|--------|---------------------|--------|---------------------|-------|
| 12 HOUR STD               | 7043                | 17.009 | 5607                | 21.207 | 4830                | 23.41 |
| UPPER LIMIT               | 14086               | 17.509 | 11214               | 21.707 | 9660                | 23.91 |
| LOWER LIMIT               | 3521.5              | 16.509 | 2803.5              | 20.707 | 2415                | 22.91 |
| EPA SAMPLE NO.            |                     |        |                     |        |                     |       |
| 01 BP-VPB-182-GW-60-62    | 6655                | 17.01  | 6097                | 21.21  | 5198                | 23.41 |
| 02 BP-VPB-182-GW-150-152  | 6626                | 17.01  | 5279                | 21.21  | 4526                | 23.41 |
| 03 BP-VPB-182-GW-220-222  | 6262                | 17.01  | 4980                | 21.21  | 4275                | 23.41 |
| 04 BP-VPB-182-EB-20250514 | 6080                | 17.01  | 4710                | 21.21  | 4050                | 23.41 |

IS4 (PHN) = Phenanthrene-d10

IS5 (CRY) = Chrysene-d12

IS6 (PRY) = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

8B

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH

Lab Code: CHEM Case No.: Q2050 SAS No.: Q2050 SDG NO.: Q2050

EPA Sample No.: SSTDCCC0.4 Date Analyzed: 05/20/2025

Lab File ID: BN037074.D Time Analyzed: 15:09

Instrument ID: BNA\_N GC Column: ZB-GR ID: 0.25 (mm)

|                | IS1 (DCB)<br>AREA # | RT #  | IS2 (NPT)<br>AREA # | RT #   | IS3 (ANT)<br>AREA # | RT #   |
|----------------|---------------------|-------|---------------------|--------|---------------------|--------|
| 12 HOUR STD    | 2240                | 7.611 | 6230                | 10.39  | 3461                | 14.26  |
| UPPER LIMIT    | 4480                | 8.111 | 12460               | 10.894 | 6922                | 14.756 |
| LOWER LIMIT    | 1120                | 7.111 | 3115                | 9.894  | 1730.5              | 13.756 |
| EPA SAMPLE NO. |                     |       |                     |        |                     |        |
| 01 PB168032BL  | 1695                | 7.61  | 4260                | 10.39  | 2468                | 14.26  |
| 02 PB168032BS  | 1842                | 7.61  | 4702                | 10.39  | 2635                | 14.26  |
| 03 PB168032BSD | 1841                | 7.61  | 4693                | 10.39  | 2613                | 14.26  |

IS1 (DCB) = 1,4-Dichlorobenzene-d4

IS2 (NPT) = Naphthalene-d8

IS3 (ANT) = Acenaphthene-d10

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

8C

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH

Lab Code: CHEM Case No.: Q2050 SAS No.: Q2050 SDG NO.: Q2050

EPA Sample No.: SSTDCCC0.4 Date Analyzed: 05/20/2025

Lab File ID: BN037074.D Time Analyzed: 15:09

Instrument ID: BNA\_N GC Column: ZB-GR ID: 0.25 (mm)

|                | IS4 (PHN)<br>AREA # | RT #   | IS5 (CRY)<br>AREA # | RT #   | IS6 (PRY)<br>AREA # | RT #   |
|----------------|---------------------|--------|---------------------|--------|---------------------|--------|
| 12 HOUR STD    | 6762                | 17.009 | 4645                | 21.207 | 3820                | 23.407 |
| UPPER LIMIT    | 13524               | 17.509 | 9290                | 21.707 | 7640                | 23.907 |
| LOWER LIMIT    | 3381                | 16.509 | 2322.5              | 20.707 | 1910                | 22.907 |
| EPA SAMPLE NO. |                     |        |                     |        |                     |        |
| 01 PB168032BL  | 5072                | 17.01  | 3622                | 21.21  | 3128                | 23.41  |
| 02 PB168032BS  | 5352                | 17.01  | 4047                | 21.21  | 3455                | 23.40  |
| 03 PB168032BSD | 5111                | 17.01  | 3649                | 21.20  | 3038                | 23.40  |

IS4 (PHN) = Phenanthrene-d10

IS5 (CRY) = Chrysene-d12

IS6 (PRY) = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.





# QC SAMPLE DATA

## Report of Analysis

|                    |                               |                  |                 |                |
|--------------------|-------------------------------|------------------|-----------------|----------------|
| Client:            | Tetra Tech NUS, Inc.          |                  | Date Collected: |                |
| Project:           | NWIRP Bethpage 112G08005-WE13 |                  | Date Received:  |                |
| Client Sample ID:  | PB168032BL                    |                  | SDG No.:        | Q2050          |
| Lab Sample ID:     | PB168032BL                    |                  | Matrix:         | Water          |
| Analytical Method: | SW8270ESIM                    |                  | % Solid:        | 0              |
| Sample Wt/Vol:     | 1000                          | 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 :      |                               |                  |                 |                |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BN037075.D        | 1         | 05/16/25 08:42 | 05/20/25 15:45 | PB168032      |

| CAS Number                | Parameter               | Conc. | Qualifier | MDL      | LOD  | LOQ / CRQL | Units    |
|---------------------------|-------------------------|-------|-----------|----------|------|------------|----------|
| <b>TARGETS</b>            |                         |       |           |          |      |            |          |
| 123-91-1                  | 1,4-Dioxane             | 0.20  | U         | 0.070    | 0.20 | 0.20       | ug/L     |
| <b>SURROGATES</b>         |                         |       |           |          |      |            |          |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.36  |           | 30 - 150 |      | 90%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.34  |           | 30 - 150 |      | 84%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.32  |           | 55 - 111 |      | 81%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.34  |           | 53 - 106 |      | 84%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.42  |           | 58 - 132 |      | 104%       | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 1700  | 7.611     |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 4260  | 10.393    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 2470  | 14.256    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 5070  | 17.008    |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 3620  | 21.206    |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 3130  | 23.409    |          |      |            |          |

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

## Report of Analysis

|                    |                               |                  |                 |                |
|--------------------|-------------------------------|------------------|-----------------|----------------|
| Client:            | Tetra Tech NUS, Inc.          |                  | Date Collected: |                |
| Project:           | NWIRP Bethpage 112G08005-WE13 |                  | Date Received:  |                |
| Client Sample ID:  | PB168032BS                    |                  | SDG No.:        | Q2050          |
| Lab Sample ID:     | PB168032BS                    |                  | Matrix:         | Water          |
| Analytical Method: | SW8270ESIM                    |                  | % Solid:        | 0              |
| Sample Wt/Vol:     | 1000                          | 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 :      |                               |                  |                 |                |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BN037080.D        | 1         | 05/16/25 08:42 | 05/20/25 18:47 | PB168032      |

| CAS Number                | Parameter               | Conc. | Qualifier | MDL      | LOD  | LOQ / CRQL | Units    |
|---------------------------|-------------------------|-------|-----------|----------|------|------------|----------|
| <b>TARGETS</b>            |                         |       |           |          |      |            |          |
| 123-91-1                  | 1,4-Dioxane             | 0.31  |           | 0.070    | 0.20 | 0.20       | ug/L     |
| <b>SURROGATES</b>         |                         |       |           |          |      |            |          |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.39  |           | 30 - 150 |      | 96%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.33  |           | 30 - 150 |      | 81%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.32  |           | 55 - 111 |      | 81%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.33  |           | 53 - 106 |      | 83%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.39  |           | 58 - 132 |      | 97%        | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 1840  | 7.611     |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 4700  | 10.394    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 2640  | 14.256    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 5350  | 17.009    |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 4050  | 21.206    |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 3460  | 23.404    |          |      |            |          |

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

## Report of Analysis

|                    |                               |                  |                 |                |
|--------------------|-------------------------------|------------------|-----------------|----------------|
| Client:            | Tetra Tech NUS, Inc.          |                  | Date Collected: |                |
| Project:           | NWIRP Bethpage 112G08005-WE13 |                  | Date Received:  |                |
| Client Sample ID:  | PB168032BSD                   |                  | SDG No.:        | Q2050          |
| Lab Sample ID:     | PB168032BSD                   |                  | Matrix:         | Water          |
| Analytical Method: | SW8270ESIM                    |                  | % Solid:        | 0              |
| Sample Wt/Vol:     | 1000                          | 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 :      |                               |                  |                 |                |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BN037081.D        | 1         | 05/16/25 08:42 | 05/20/25 19:23 | PB168032      |

| CAS Number                | Parameter               | Conc. | Qualifier | MDL      | LOD  | LOQ / CRQL | Units    |
|---------------------------|-------------------------|-------|-----------|----------|------|------------|----------|
| <b>TARGETS</b>            |                         |       |           |          |      |            |          |
| 123-91-1                  | 1,4-Dioxane             | 0.32  |           | 0.070    | 0.20 | 0.20       | ug/L     |
| <b>SURROGATES</b>         |                         |       |           |          |      |            |          |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.39  |           | 30 - 150 |      | 98%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.33  |           | 30 - 150 |      | 83%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.35  |           | 55 - 111 |      | 88%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.33  |           | 53 - 106 |      | 83%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.42  |           | 58 - 132 |      | 104%       | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 1840  | 7.611     |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 4690  | 10.394    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 2610  | 14.256    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 5110  | 17.009    |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 3650  | 21.198    |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 3040  | 23.401    |          |      |            |          |

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



# CALIBRATION SUMMARY

Method Path : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\  
 Method File : 8270-SIM-BN051425.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 Last Update : Wed May 14 11:26:32 2025  
 Response Via : Initial Calibration

## Calibration Files

0.1 =BN036999.D 0.2 =BN037000.D 0.4 =BN037001.D 0.8 =BN037002.D 1.6 =BN037003.D 3.2 =BN037004.D 5.0 =BN037005.D

|           | Compound              | 0.1            | 0.2   | 0.4   | 0.8   | 1.6   | 3.2   | 5.0   | Avg   | %RSD  |
|-----------|-----------------------|----------------|-------|-------|-------|-------|-------|-------|-------|-------|
| -----     |                       |                |       |       |       |       |       |       |       |       |
| 1) I      | 1,4-Dichlorobenzen... | -----ISTD----- |       |       |       |       |       |       |       |       |
| 2)        | 1,4-Dioxane           | 0.510          | 0.512 | 0.487 | 0.514 | 0.467 | 0.454 | 0.491 | 5.25  |       |
| 3)        | n-Nitrosodimet...     | 1.465          | 0.974 | 0.980 | 0.971 | 1.075 | 0.967 | 0.950 | 1.054 | 17.59 |
| 4) S      | 2-Fluorophenol        | 1.101          | 1.134 | 1.024 | 1.093 | 0.964 | 0.971 | 1.048 | 6.87  |       |
| 5) S      | Phenol-d6             | 1.304          | 1.385 | 1.236 | 1.392 | 1.259 | 1.292 | 1.311 | 4.91  |       |
| 6)        | bis(2-Chloroet...     | 1.441          | 1.163 | 1.153 | 1.135 | 1.240 | 1.168 | 1.148 | 1.207 | 9.02  |
|           |                       |                |       |       |       |       |       |       |       |       |
| 7) I      | Naphthalene-d8        | -----ISTD----- |       |       |       |       |       |       |       |       |
| 8) S      | Nitrobenzene-d5       | 0.546          | 0.383 | 0.398 | 0.400 | 0.452 | 0.426 | 0.442 | 0.436 | 12.60 |
| 9)        | Naphthalene           | 1.326          | 1.140 | 1.144 | 1.122 | 1.226 | 1.152 | 1.165 | 1.182 | 6.05  |
| 10)       | Hexachlorobuta...     | 0.286          | 0.248 | 0.244 | 0.235 | 0.256 | 0.236 | 0.233 | 0.248 | 7.47  |
| 11) SURR2 | Methylnaphth...       | 0.529          | 0.547 | 0.552 | 0.548 | 0.603 | 0.574 | 0.588 | 0.563 | 4.65  |
| 12)       | 2-Methylnaphth...     | 0.754          | 0.724 | 0.736 | 0.733 | 0.814 | 0.770 | 0.790 | 0.760 | 4.34  |
|           |                       |                |       |       |       |       |       |       |       |       |
| 13) I     | Acenaphthene-d10      | -----ISTD----- |       |       |       |       |       |       |       |       |
| 14) S     | 2,4,6-Tribromo...     | 0.174          | 0.168 | 0.178 | 0.160 | 0.186 | 0.175 | 0.189 | 0.176 | 5.77  |
| 15) S     | 2-Fluorobiphenyl      | 1.912          | 1.801 | 1.901 | 1.802 | 1.927 | 1.672 | 1.807 | 1.832 | 4.90  |
| 16)       | Acenaphthylene        | 1.906          | 1.838 | 1.894 | 1.849 | 2.071 | 1.997 | 2.075 | 1.947 | 5.14  |
| 17)       | Acenaphthene          | 1.255          | 1.229 | 1.243 | 1.217 | 1.350 | 1.298 | 1.315 | 1.272 | 3.89  |
| 18)       | Fluorene              | 1.602          | 1.581 | 1.635 | 1.611 | 1.779 | 1.721 | 1.752 | 1.669 | 4.80  |
|           |                       |                |       |       |       |       |       |       |       |       |
| 19) I     | Phenanthrene-d10      | -----ISTD----- |       |       |       |       |       |       |       |       |
| 20)       | 4,6-Dinitro-2-...     | 0.060          | 0.073 | 0.079 | 0.102 | 0.103 | 0.124 | 0.090 | 26.02 |       |
| 21)       | 4-Bromophenyl-...     | 0.243          | 0.246 | 0.250 | 0.247 | 0.262 | 0.261 | 0.259 | 0.253 | 3.13  |
| 22)       | Hexachlorobenzene     | 0.267          | 0.269 | 0.281 | 0.259 | 0.281 | 0.270 | 0.267 | 0.270 | 3.03  |
| 23)       | Atrazine              | 0.199          | 0.207 | 0.211 | 0.213 | 0.237 | 0.234 | 0.242 | 0.220 | 7.64  |
| 24)       | Pentachlorophenol     | 0.133          | 0.134 | 0.141 | 0.137 | 0.159 | 0.162 | 0.177 | 0.149 | 11.45 |
| 25)       | Phenanthrene          | 1.259          | 1.272 | 1.292 | 1.263 | 1.367 | 1.337 | 1.361 | 1.307 | 3.56  |
| 26)       | Anthracene            | 1.099          | 1.104 | 1.166 | 1.130 | 1.269 | 1.259 | 1.300 | 1.190 | 7.13  |
| 27) SURR  | Fluoranthene-d10      | 1.033          | 1.033 | 1.078 | 1.042 | 1.153 | 1.161 | 1.178 | 1.097 | 5.95  |
| 28)       | Fluoranthene          | 1.461          | 1.439 | 1.500 | 1.496 | 1.670 | 1.672 | 1.693 | 1.562 | 7.13  |
|           |                       |                |       |       |       |       |       |       |       |       |
| 29) I     | Chrysene-d12          | -----ISTD----- |       |       |       |       |       |       |       |       |
| 30)       | Pyrene                | 1.744          | 1.708 | 1.727 | 1.656 | 1.790 | 1.641 | 1.711 | 1.711 | 2.96  |
| 31) S     | Terphenyl-d14         | 0.897          | 0.844 | 0.871 | 0.822 | 0.891 | 0.816 | 0.848 | 0.856 | 3.73  |
| 32)       | Benzo(a)anthra...     | 1.463          | 1.432 | 1.485 | 1.438 | 1.594 | 1.521 | 1.609 | 1.506 | 4.77  |
| 33)       | Chrysene              | 1.655          | 1.559 | 1.616 | 1.532 | 1.653 | 1.560 | 1.576 | 1.593 | 3.05  |
| 34)       | Bis(2-ethylhex...     | 0.955          | 0.919 | 0.906 | 0.855 | 0.941 | 0.903 | 1.011 | 0.927 | 5.27  |
|           |                       |                |       |       |       |       |       |       |       |       |
| 35) I     | Perylene-d12          | -----ISTD----- |       |       |       |       |       |       |       |       |

Method Path : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\

Method File : 8270-SIM-BN051425.M

|       |                   |       |       |       |       |       |       |       |       |      |
|-------|-------------------|-------|-------|-------|-------|-------|-------|-------|-------|------|
| 36)   | Indeno(1,2,3-c... | 1.511 | 1.613 | 1.645 | 1.568 | 1.687 | 1.732 | 1.680 | 1.634 | 4.65 |
| 37)   | Benzo(b)fluora... | 1.631 | 1.570 | 1.602 | 1.599 | 1.749 | 1.698 | 1.765 | 1.659 | 4.71 |
| 38)   | Benzo(k)fluora... | 1.539 | 1.538 | 1.642 | 1.601 | 1.770 | 1.661 | 1.719 | 1.639 | 5.34 |
| 39) C | Benzo(a)pyrene    | 1.380 | 1.343 | 1.381 | 1.331 | 1.486 | 1.444 | 1.486 | 1.407 | 4.59 |
| 40)   | Dibenzo(a,h)an... | 1.116 | 1.232 | 1.273 | 1.237 | 1.340 | 1.376 | 1.334 | 1.272 | 6.90 |
| 41)   | Benzo(g,h,i)pe... | 1.299 | 1.407 | 1.424 | 1.330 | 1.403 | 1.439 | 1.376 | 1.383 | 3.72 |

-----  
(#) = Out of Range

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: TETR06

Lab Code: CHEM Case No.: Q2050 SAS No.: Q2050 SDG No.: Q2050

Instrument ID: BNA\_N Calibration Date/Time: 05/19/2025 10:47

Lab File ID: BN037038.D Init. Calib. Date(s): 05/13/2025 05/13/2025

EPA Sample No.: SSTDCCC0.4 Init. Calib. Time(s): 17:41 21:17

GC Column: ZB-GR ID: 0.25 (mm)

| COMPOUND                | RRF   | RRF0.4 | MIN<br>RRF | %D    | MAX%D |
|-------------------------|-------|--------|------------|-------|-------|
| 2-Methylnaphthalene-d10 | 0.563 | 0.598  |            | 6.2   | 20.0  |
| Fluoranthene-d10        | 1.097 | 1.086  |            | -1.0  | 20.0  |
| 2-Fluorophenol          | 1.048 | 0.944  |            | -9.9  | 20.0  |
| Phenol-d6               | 1.311 | 1.169  |            | -10.8 | 20.0  |
| Nitrobenzene-d5         | 0.436 | 0.392  |            | -10.1 | 20.0  |
| 2-Fluorobiphenyl        | 1.832 | 1.706  |            | -6.9  | 20.0  |
| 2,4,6-Tribromophenol    | 0.176 | 0.147  |            | -16.5 | 20.0  |
| Terphenyl-d14           | 0.856 | 0.915  |            | 6.9   | 20.0  |
| 1,4-Dioxane             | 0.491 | 0.461  |            | -6.1  | 20.0  |

All other compounds must meet a minimum RRF of 0.010.



7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: TETR06

Lab Code: CHEM Case No.: Q2050 SAS No.: Q2050 SDG No.: Q2050

Instrument ID: BNA\_N Calibration Date/Time: 05/19/2025 21:05

Lab File ID: BN037055.D Init. Calib. Date(s): 05/13/2025 05/13/2025

EPA Sample No.: SSTDCCC0.4EC Init. Calib. Time(s): 17:41 21:17

GC Column: ZB-GR ID: 0.25 (mm)

| COMPOUND                | RRF   | RRF0.4 | MIN<br>RRF | %D    | MAX%D |
|-------------------------|-------|--------|------------|-------|-------|
| 2-Methylnaphthalene-d10 | 0.563 | 0.581  |            | 3.2   | 50.0  |
| Fluoranthene-d10        | 1.097 | 1.058  |            | -3.6  | 50.0  |
| 2-Fluorophenol          | 1.048 | 0.956  |            | -8.8  | 50.0  |
| Phenol-d6               | 1.311 | 1.171  |            | -10.7 | 50.0  |
| Nitrobenzene-d5         | 0.436 | 0.406  |            | -6.9  | 50.0  |
| 2-Fluorobiphenyl        | 1.832 | 1.804  |            | -1.5  | 50.0  |
| 2,4,6-Tribromophenol    | 0.176 | 0.160  |            | -9.1  | 50.0  |
| Terphenyl-d14           | 0.856 | 0.927  |            | 8.3   | 50.0  |
| 1,4-Dioxane             | 0.491 | 0.456  |            | -7.1  | 50.0  |

All other compounds must meet a minimum RRF of 0.010.

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: TETR06

Lab Code: CHEM Case No.: Q2050 SAS No.: Q2050 SDG No.: Q2050

Instrument ID: BNA\_N Calibration Date/Time: 05/20/2025 15:09

Lab File ID: BN037074.D Init. Calib. Date(s): 05/13/2025 05/13/2025

EPA Sample No.: SSTDCCC0.4 Init. Calib. Time(s): 17:41 21:17

GC Column: ZB-GR ID: 0.25 (mm)

| COMPOUND                | RRF   | RRF0.4 | MIN<br>RRF | %D    | MAX%D |
|-------------------------|-------|--------|------------|-------|-------|
| 2-Methylnaphthalene-d10 | 0.563 | 0.573  |            | 1.8   | 20.0  |
| Fluoranthene-d10        | 1.097 | 1.010  |            | -7.9  | 20.0  |
| 2-Fluorophenol          | 1.048 | 0.964  |            | -8.0  | 20.0  |
| Phenol-d6               | 1.311 | 1.198  |            | -8.6  | 20.0  |
| Nitrobenzene-d5         | 0.436 | 0.402  |            | -7.8  | 20.0  |
| 2-Fluorobiphenyl        | 1.832 | 1.798  |            | -1.9  | 20.0  |
| 2,4,6-Tribromophenol    | 0.176 | 0.153  |            | -13.1 | 20.0  |
| Terphenyl-d14           | 0.856 | 0.940  |            | 9.8   | 20.0  |
| 1,4-Dioxane             | 0.491 | 0.488  |            | -0.6  | 20.0  |

All other compounds must meet a minimum RRF of 0.010.

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: TETR06

Lab Code: CHEM Case No.: Q2050 SAS No.: Q2050 SDG No.: Q2050

Instrument ID: BNA\_N Calibration Date/Time: 05/20/2025 19:59

Lab File ID: BN037082.D Init. Calib. Date(s): 05/13/2025 05/13/2025

EPA Sample No.: SSTDCCC0.4EC Init. Calib. Time(s): 17:41 21:17

GC Column: ZB-GR ID: 0.25 (mm)

| COMPOUND                | RRF   | RRF0.4 | MIN<br>RRF | %D    | MAX%D |
|-------------------------|-------|--------|------------|-------|-------|
| 2-Methylnaphthalene-d10 | 0.563 | 0.581  |            | 3.2   | 50.0  |
| Fluoranthene-d10        | 1.097 | 1.017  |            | -7.3  | 50.0  |
| 2-Fluorophenol          | 1.048 | 0.946  |            | -9.7  | 50.0  |
| Phenol-d6               | 1.311 | 1.150  |            | -12.3 | 50.0  |
| Nitrobenzene-d5         | 0.436 | 0.417  |            | -4.4  | 50.0  |
| 2-Fluorobiphenyl        | 1.832 | 1.796  |            | -2.0  | 50.0  |
| 2,4,6-Tribromophenol    | 0.176 | 0.165  |            | -6.3  | 50.0  |
| Terphenyl-d14           | 0.856 | 0.957  |            | 11.8  | 50.0  |
| 1,4-Dioxane             | 0.491 | 0.473  |            | -3.7  | 50.0  |

All other compounds must meet a minimum RRF of 0.010.