

## Report of Analysis

|                    |                              |           |                 |               |
|--------------------|------------------------------|-----------|-----------------|---------------|
| Client:            | Portal Partners Tri-Venture  |           | Date Collected: |               |
| Project:           | Amtrak Sawtooth Bridges 2024 |           | Date Received:  |               |
| Client Sample ID:  | VN1101WBS01                  |           | SDG No.:        | P4658         |
| Lab Sample ID:     | VN1101WBS01                  |           | Matrix:         | Water         |
| Analytical Method: | SW8260                       |           | % Solid:        | 0             |
| Sample Wt/Vol:     | 5                            | Units: mL | Final Vol:      | 5000 uL       |
| Soil Aliquot Vol:  |                              | uL        | Test:           | VOC-TCLVOA-10 |
| GC Column:         | RXI-624                      | ID : 0.25 | Level :         | LOW           |
| Prep Method :      |                              |           |                 |               |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VN084631.D        | 1         |           | 11/01/24 12:08 | VN110124      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                                |       |           |      |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 19.2  |           | 0.21 | 1.00       | ug/L  |
| 74-87-3        | Chloromethane                  | 16.4  |           | 0.35 | 1.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 18.9  |           | 0.34 | 1.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 18.0  |           | 1.40 | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 18.2  |           | 0.56 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 18.8  |           | 0.34 | 1.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 19.3  |           | 0.25 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 17.7  |           | 0.26 | 1.00       | ug/L  |
| 67-64-1        | Acetone                        | 96.0  |           | 1.40 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 16.7  |           | 0.32 | 1.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 19.1  |           | 0.16 | 1.00       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 16.1  |           | 0.60 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 19.0  |           | 0.32 | 1.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 18.5  |           | 0.25 | 1.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 19.1  |           | 0.23 | 1.00       | ug/L  |
| 110-82-7       | Cyclohexane                    | 18.3  |           | 1.60 | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 96.7  |           | 1.30 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 19.7  |           | 0.25 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 18.4  |           | 0.25 | 1.00       | ug/L  |
| 74-97-5        | Bromochloromethane             | 21.5  |           | 0.18 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                     | 19.6  |           | 0.26 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 19.4  |           | 0.19 | 1.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 18.4  |           | 0.19 | 1.00       | ug/L  |
| 71-43-2        | Benzene                        | 18.8  |           | 0.16 | 1.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 19.5  |           | 0.24 | 1.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 18.4  |           | 0.32 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 19.1  |           | 0.19 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 19.2  |           | 0.24 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 100   |           | 0.75 | 5.00       | ug/L  |
| 108-88-3       | Toluene                        | 20.1  |           | 0.18 | 1.00       | ug/L  |

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| Client Sample ID:  | VN1101WBS01                  |           | SDG No.:        | P4658         |
| Lab Sample ID:     | VN1101WBS01                  |           | Matrix:         | Water         |
| Analytical Method: | SW8260                       |           | % Solid:        | 0             |
| Sample Wt/Vol:     | 5                            | Units: mL | Final Vol:      | 5000 uL       |
| Soil Aliquot Vol:  |                              | uL        | Test:           | VOC-TCLVOA-10 |
| GC Column:         | RXI-624                      | ID : 0.25 | Level :         | LOW           |
| Prep Method :      |                              |           |                 |               |

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|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VN084631.D        | 1         |           | 11/01/24 12:08 | VN110124      |

| CAS Number                | Parameter                   | Conc.  | Qualifier | MDL                 | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|--------|-----------|---------------------|------------|---------|
| 10061-02-6                | t-1,3-Dichloropropene       | 17.7   |           | 0.21                | 1.00       | ug/L    |
| 10061-01-5                | cis-1,3-Dichloropropene     | 18.7   |           | 0.18                | 1.00       | ug/L    |
| 79-00-5                   | 1,1,2-Trichloroethane       | 20.0   |           | 0.21                | 1.00       | ug/L    |
| 591-78-6                  | 2-Hexanone                  | 110    |           | 1.10                | 5.00       | ug/L    |
| 124-48-1                  | Dibromochloromethane        | 20.3   |           | 0.18                | 1.00       | ug/L    |
| 106-93-4                  | 1,2-Dibromoethane           | 19.0   |           | 0.16                | 1.00       | ug/L    |
| 127-18-4                  | Tetrachloroethene           | 18.6   |           | 0.25                | 1.00       | ug/L    |
| 108-90-7                  | Chlorobenzene               | 18.3   |           | 0.13                | 1.00       | ug/L    |
| 100-41-4                  | Ethyl Benzene               | 18.8   |           | 0.16                | 1.00       | ug/L    |
| 179601-23-1               | m/p-Xylenes                 | 38.9   |           | 0.31                | 2.00       | ug/L    |
| 95-47-6                   | o-Xylene                    | 20.0   |           | 0.14                | 1.00       | ug/L    |
| 100-42-5                  | Styrene                     | 19.4   |           | 0.16                | 1.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 19.0   |           | 0.21                | 1.00       | ug/L    |
| 98-82-8                   | Isopropylbenzene            | 18.5   |           | 0.13                | 1.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 18.1   |           | 0.27                | 1.00       | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene         | 16.8   |           | 0.24                | 1.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 17.3   |           | 0.27                | 1.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 17.0   |           | 0.19                | 1.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 17.9   |           | 0.46                | 1.00       | ug/L    |
| 120-82-1                  | 1,2,4-Trichlorobenzene      | 14.7   |           | 0.42                | 1.00       | ug/L    |
| 87-61-6                   | 1,2,3-Trichlorobenzene      | 15.3   |           | 0.51                | 1.00       | ug/L    |
| <b>SURROGATES</b>         |                             |        |           |                     |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 44.3   |           | 70 (74) - 130 (125) | 89%        | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 45.1   |           | 70 (75) - 130 (124) | 90%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 46.1   |           | 70 (86) - 130 (113) | 92%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 47.0   |           | 70 (77) - 130 (121) | 94%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |        |           |                     |            |         |
| 363-72-4                  | Pentafluorobenzene          | 191000 | 8.224     |                     |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 320000 | 9.1       |                     |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 288000 | 11.865    |                     |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 147000 | 13.794    |                     |            |         |

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| CAS Number | Parameter | Conc. | Qualifier | MDL | 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