

Report of Analysis

Client:	Portal Partners Tri-Venture		Date Collected:	06/01/25	
Project:	Amtrak Sawtooth Bridges 2025		Date Received:	06/03/25	
Client Sample ID:	B-207-SB02		SDG No.:	Q2198	
Lab Sample ID:	Q2198-03		Matrix:	SOIL	
Analytical Method:	8260D		% Solid:	52.1	
Sample Wt/Vol:	4.87	Units: g	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
VY022553.D	1		06/04/25 16:11	VY060425

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
75-71-8	Dichlorodifluoromethane	2.20	U	2.20	9.90	ug/Kg
74-87-3	Chloromethane	2.20	U	2.20	9.90	ug/Kg
75-01-4	Vinyl Chloride	1.60	U	1.60	9.90	ug/Kg
74-83-9	Bromomethane	2.10	U	2.10	9.90	ug/Kg
75-00-3	Chloroethane	2.50	U	2.50	9.90	ug/Kg
75-69-4	Trichlorofluoromethane	2.40	U	2.40	9.90	ug/Kg
76-13-1	1,1,2-Trichlorotrifluoroethane	2.10	U	2.10	9.90	ug/Kg
75-35-4	1,1-Dichloroethene	2.00	U	2.00	9.90	ug/Kg
67-64-1	Acetone	71.8		9.30	49.3	ug/Kg
75-15-0	Carbon Disulfide	3.40	J	2.10	9.90	ug/Kg
1634-04-4	Methyl tert-butyl Ether	1.40	U	1.40	9.90	ug/Kg
79-20-9	Methyl Acetate	3.00	U	3.00	9.90	ug/Kg
75-09-2	Methylene Chloride	7.00	U	7.00	19.7	ug/Kg
156-60-5	trans-1,2-Dichloroethene	1.70	U	1.70	9.90	ug/Kg
75-34-3	1,1-Dichloroethane	1.60	U	1.60	9.90	ug/Kg
110-82-7	Cyclohexane	1.60	U	1.60	9.90	ug/Kg
78-93-3	2-Butanone	14.8	J	12.9	49.3	ug/Kg
56-23-5	Carbon Tetrachloride	1.90	U	1.90	9.90	ug/Kg
156-59-2	cis-1,2-Dichloroethene	1.50	U	1.50	9.90	ug/Kg
74-97-5	Bromochloromethane	2.30	U	2.30	9.90	ug/Kg
67-66-3	Chloroform	1.70	U	1.70	9.90	ug/Kg
71-55-6	1,1,1-Trichloroethane	1.80	U	1.80	9.90	ug/Kg
108-87-2	Methylcyclohexane	1.80	U	1.80	9.90	ug/Kg
71-43-2	Benzene	1.60	U	1.60	9.90	ug/Kg
107-06-2	1,2-Dichloroethane	1.60	U	1.60	9.90	ug/Kg
79-01-6	Trichloroethene	1.60	U	1.60	9.90	ug/Kg
78-87-5	1,2-Dichloropropane	1.80	U	1.80	9.90	ug/Kg
75-27-4	Bromodichloromethane	1.50	U	1.50	9.90	ug/Kg
108-10-1	4-Methyl-2-Pentanone	7.10	U	7.10	49.3	ug/Kg
108-88-3	Toluene	1.50	U	1.50	9.90	ug/Kg

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10061-02-6	t-1,3-Dichloropropene	1.30	U	1.30	9.90	ug/Kg
10061-01-5	cis-1,3-Dichloropropene	1.20	U	1.20	9.90	ug/Kg
79-00-5	1,1,2-Trichloroethane	1.80	U	1.80	9.90	ug/Kg
591-78-6	2-Hexanone	7.30	U	7.30	49.3	ug/Kg
124-48-1	Dibromochloromethane	1.70	U	1.70	9.90	ug/Kg
106-93-4	1,2-Dibromoethane	1.70	U	1.70	9.90	ug/Kg
127-18-4	Tetrachloroethene	2.10	U	2.10	9.90	ug/Kg
108-90-7	Chlorobenzene	1.80	U	1.80	9.90	ug/Kg
100-41-4	Ethyl Benzene	1.30	U	1.30	9.90	ug/Kg
179601-23-1	m/p-Xylenes	2.40	U	2.40	19.7	ug/Kg
95-47-6	o-Xylene	1.60	U	1.60	9.90	ug/Kg
100-42-5	Styrene	1.40	U	1.40	9.90	ug/Kg
75-25-2	Bromoform	1.70	U	1.70	9.90	ug/Kg
98-82-8	Isopropylbenzene	1.50	U	1.50	9.90	ug/Kg
79-34-5	1,1,2,2-Tetrachloroethane	2.40	U	2.40	9.90	ug/Kg
541-73-1	1,3-Dichlorobenzene	3.40	U	3.40	9.90	ug/Kg
106-46-7	1,4-Dichlorobenzene	3.10	U	3.10	9.90	ug/Kg
95-50-1	1,2-Dichlorobenzene	2.90	U	2.90	9.90	ug/Kg
96-12-8	1,2-Dibromo-3-Chloropropane	3.60	U	3.60	9.90	ug/Kg
120-82-1	1,2,4-Trichlorobenzene	5.90	U	5.90	9.90	ug/Kg
87-61-6	1,2,3-Trichlorobenzene	6.30	U	6.30	9.90	ug/Kg
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	51.4		70 (63) - 130 (155)	103%	SPK: 50
1868-53-7	Dibromofluoromethane	50.7		70 (70) - 130 (134)	101%	SPK: 50
2037-26-5	Toluene-d8	49.5		70 (74) - 130 (123)	99%	SPK: 50
460-00-4	4-Bromofluorobenzene	37.6		70 (17) - 130 (146)	75%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	232000	7.707			
540-36-3	1,4-Difluorobenzene	423000	8.609			
3114-55-4	Chlorobenzene-d5	325000	11.414			
3855-82-1	1,4-Dichlorobenzene-d4	102000	13.34			

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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