

Report of Analysis

Client:	Portal Partners Tri-Venture		Date Collected:	08/11/25	
Project:	Amtrak Sawtooth Bridges 2025		Date Received:	08/12/25	
Client Sample ID:	TG-S03		SDG No.:	Q2832	
Lab Sample ID:	Q2832-05		Matrix:	SOIL	
Analytical Method:	8260D		% Solid:	93.8	
Sample Wt/Vol:	3.86	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:	Date Analyzed	Prep Batch ID
VY023125.D	1	08/15/25 15:11	VY081525

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
75-71-8	Dichlorodifluoromethane	1.60	U	1.60	6.90	ug/Kg
74-87-3	Chloromethane	1.60	U	1.60	6.90	ug/Kg
75-01-4	Vinyl Chloride	1.10	U	1.10	6.90	ug/Kg
74-83-9	Bromomethane	1.50	U	1.50	6.90	ug/Kg
75-00-3	Chloroethane	1.70	U	1.70	6.90	ug/Kg
75-69-4	Trichlorofluoromethane	1.70	U	1.70	6.90	ug/Kg
76-13-1	1,1,2-Trichlorotrifluoroethane	1.50	U	1.50	6.90	ug/Kg
75-35-4	1,1-Dichloroethene	1.40	U	1.40	6.90	ug/Kg
67-64-1	Acetone	6.50	U	6.50	34.5	ug/Kg
75-15-0	Carbon Disulfide	1.50	U	1.50	6.90	ug/Kg
1634-04-4	Methyl tert-butyl Ether	1.00	U	1.00	6.90	ug/Kg
79-20-9	Methyl Acetate	2.10	U	2.10	6.90	ug/Kg
75-09-2	Methylene Chloride	4.90	U	4.90	13.8	ug/Kg
156-60-5	trans-1,2-Dichloroethene	1.20	U	1.20	6.90	ug/Kg
75-34-3	1,1-Dichloroethane	1.10	U	1.10	6.90	ug/Kg
110-82-7	Cyclohexane	1.10	U	1.10	6.90	ug/Kg
78-93-3	2-Butanone	9.00	U	9.00	34.5	ug/Kg
56-23-5	Carbon Tetrachloride	1.30	U	1.30	6.90	ug/Kg
156-59-2	cis-1,2-Dichloroethene	1.00	U	1.00	6.90	ug/Kg
74-97-5	Bromochloromethane	1.60	U	1.60	6.90	ug/Kg
67-66-3	Chloroform	1.20	U	1.20	6.90	ug/Kg
71-55-6	1,1,1-Trichloroethane	1.30	U	1.30	6.90	ug/Kg
108-87-2	Methylcyclohexane	1.30	U	1.30	6.90	ug/Kg
71-43-2	Benzene	1.10	U	1.10	6.90	ug/Kg
107-06-2	1,2-Dichloroethane	1.10	U	1.10	6.90	ug/Kg
79-01-6	Trichloroethene	1.10	U	1.10	6.90	ug/Kg
78-87-5	1,2-Dichloropropane	1.30	U	1.30	6.90	ug/Kg
75-27-4	Bromodichloromethane	1.10	U	1.10	6.90	ug/Kg
108-10-1	4-Methyl-2-Pentanone	4.90	U	4.90	34.5	ug/Kg
108-88-3	Toluene	1.10	U	1.10	6.90	ug/Kg

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CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
10061-02-6	t-1,3-Dichloropropene	0.90	U	0.90	6.90	ug/Kg
10061-01-5	cis-1,3-Dichloropropene	0.86	U	0.86	6.90	ug/Kg
79-00-5	1,1,2-Trichloroethane	1.30	U	1.30	6.90	ug/Kg
591-78-6	2-Hexanone	5.10	U	5.10	34.5	ug/Kg
124-48-1	Dibromochloromethane	1.20	U	1.20	6.90	ug/Kg
106-93-4	1,2-Dibromoethane	1.20	U	1.20	6.90	ug/Kg
127-18-4	Tetrachloroethene	1.50	U	1.50	6.90	ug/Kg
108-90-7	Chlorobenzene	1.30	U	1.30	6.90	ug/Kg
100-41-4	Ethyl Benzene	0.93	U	0.93	6.90	ug/Kg
179601-23-1	m/p-Xylenes	1.70	U	1.70	13.8	ug/Kg
95-47-6	o-Xylene	1.10	U	1.10	6.90	ug/Kg
100-42-5	Styrene	0.98	U	0.98	6.90	ug/Kg
75-25-2	Bromoform	1.20	U	1.20	6.90	ug/Kg
98-82-8	Isopropylbenzene	1.10	U	1.10	6.90	ug/Kg
79-34-5	1,1,2,2-Tetrachloroethane	1.70	U	1.70	6.90	ug/Kg
541-73-1	1,3-Dichlorobenzene	2.40	U	2.40	6.90	ug/Kg
106-46-7	1,4-Dichlorobenzene	2.20	U	2.20	6.90	ug/Kg
95-50-1	1,2-Dichlorobenzene	2.00	U	2.00	6.90	ug/Kg
96-12-8	1,2-Dibromo-3-Chloropropane	2.50	U	2.50	6.90	ug/Kg
120-82-1	1,2,4-Trichlorobenzene	4.10	U	4.10	6.90	ug/Kg
87-61-6	1,2,3-Trichlorobenzene	4.40	U	4.40	6.90	ug/Kg
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	59.2		70 (63) - 130 (155)	118%	SPK: 50
1868-53-7	Dibromofluoromethane	51.4		70 (70) - 130 (134)	103%	SPK: 50
2037-26-5	Toluene-d8	52.7		70 (74) - 130 (123)	105%	SPK: 50
460-00-4	4-Bromofluorobenzene	48.8		70 (17) - 130 (146)	98%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	554000	7.707			
540-36-3	1,4-Difluorobenzene	1040000	8.609			
3114-55-4	Chlorobenzene-d5	947000	11.414			
3855-82-1	1,4-Dichlorobenzene-d4	379000	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