

Report of Analysis

Client:	Portal Partners Tri-Venture		Date Collected:	
Project:	Amtrak Sawtooth Bridges 2025		Date Received:	
Client Sample ID:	PB166982BSD		SDG No.:	Q1488
Lab Sample ID:	PB166982BSD		Matrix:	Water
Analytical Method:	SW8270		% Solid:	0
Sample Wt/Vol:	1000	Units: mL	Final Vol:	1000 uL
Soil Aliquot Vol:		uL	Test:	SVOC-TCL BNA -20
Extraction Type :		Decanted : N	Level :	LOW
Injection Volume :		GPC Factor : 1.0	GPC Cleanup :	N PH :
Prep Method :	SW3510C			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BF141863.D	1	03/05/25 08:35	03/06/25 12:21	PB166982

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
100-52-7	Benzaldehyde	23.1		4.00	10.0	ug/L
108-95-2	Phenol	41.6		0.93	5.00	ug/L
111-44-4	bis(2-Chloroethyl)ether	41.1		1.20	5.00	ug/L
95-57-8	2-Chlorophenol	44.8		0.71	5.00	ug/L
95-48-7	2-Methylphenol	43.2		1.10	5.00	ug/L
108-60-1	2,2-oxybis(1-Chloropropane)	35.1		1.40	5.00	ug/L
98-86-2	Acetophenone	46.8		1.10	5.00	ug/L
65794-96-9	3+4-Methylphenols	42.8		1.20	10.0	ug/L
621-64-7	n-Nitroso-di-n-propylamine	38.5		1.50	2.50	ug/L
67-72-1	Hexachloroethane	44.7		1.00	5.00	ug/L
98-95-3	Nitrobenzene	48.5		1.30	5.00	ug/L
78-59-1	Isophorone	43.1		1.10	5.00	ug/L
88-75-5	2-Nitrophenol	53.4		2.00	5.00	ug/L
105-67-9	2,4-Dimethylphenol	53.9		1.50	5.00	ug/L
111-91-1	bis(2-Chloroethoxy)methane	40.1		1.00	5.00	ug/L
120-83-2	2,4-Dichlorophenol	44.6		0.88	5.00	ug/L
91-20-3	Naphthalene	43.1		1.00	5.00	ug/L
106-47-8	4-Chloroaniline	20.1		1.30	5.00	ug/L
87-68-3	Hexachlorobutadiene	46.6		1.30	5.00	ug/L
105-60-2	Caprolactam	50.5		1.70	10.0	ug/L
59-50-7	4-Chloro-3-methylphenol	43.9		0.84	5.00	ug/L
91-57-6	2-Methylnaphthalene	41.2		1.10	5.00	ug/L
77-47-4	Hexachlorocyclopentadiene	160	E	5.00	10.0	ug/L
88-06-2	2,4,6-Trichlorophenol	47.0		0.89	5.00	ug/L
95-95-4	2,4,5-Trichlorophenol	48.8		1.00	5.00	ug/L
92-52-4	1,1-Biphenyl	48.0		0.91	5.00	ug/L
91-58-7	2-Chloronaphthalene	44.3		0.97	5.00	ug/L
88-74-4	2-Nitroaniline	51.4		1.40	5.00	ug/L
131-11-3	Dimethylphthalate	44.6		0.93	5.00	ug/L

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208-96-8	Acenaphthylene	46.3		1.00	5.00	ug/L
606-20-2	2,6-Dinitrotoluene	51.6		1.20	5.00	ug/L
99-09-2	3-Nitroaniline	34.7		1.40	5.00	ug/L
83-32-9	Acenaphthene	51.0		0.81	5.00	ug/L
51-28-5	2,4-Dinitrophenol	120	E	6.40	10.0	ug/L
100-02-7	4-Nitrophenol	120	E	2.00	10.0	ug/L
132-64-9	Dibenzofuran	42.4		0.93	5.00	ug/L
121-14-2	2,4-Dinitrotoluene	57.5		1.50	5.00	ug/L
84-66-2	Diethylphthalate	43.2		1.00	5.00	ug/L
7005-72-3	4-Chlorophenyl-phenylether	44.2		0.98	5.00	ug/L
86-73-7	Fluorene	44.5		0.96	5.00	ug/L
100-01-6	4-Nitroaniline	61.0		2.00	5.00	ug/L
534-52-1	4,6-Dinitro-2-methylphenol	59.5		3.10	10.0	ug/L
86-30-6	n-Nitrosodiphenylamine	43.0		0.89	5.00	ug/L
101-55-3	4-Bromophenyl-phenylether	43.4		0.95	5.00	ug/L
118-74-1	Hexachlorobenzene	46.8		1.10	5.00	ug/L
1912-24-9	Atrazine	63.7		1.30	5.00	ug/L
87-86-5	Pentachlorophenol	97.3	E	1.90	10.0	ug/L
85-01-8	Phenanthrene	45.5		0.89	5.00	ug/L
120-12-7	Anthracene	46.1		1.10	5.00	ug/L
86-74-8	Carbazole	46.4		1.20	5.00	ug/L
84-74-2	Di-n-butylphthalate	42.7		1.50	5.00	ug/L
206-44-0	Fluoranthene	47.4		1.30	5.00	ug/L
129-00-0	Pyrene	42.3		1.10	5.00	ug/L
85-68-7	Butylbenzylphthalate	45.7		2.10	5.00	ug/L
91-94-1	3,3-Dichlorobenzidine	29.7		1.30	10.0	ug/L
56-55-3	Benzo(a)anthracene	44.6		0.94	5.00	ug/L
218-01-9	Chrysene	44.7		0.86	5.00	ug/L
117-81-7	Bis(2-ethylhexyl)phthalate	46.9		1.90	5.00	ug/L
117-84-0	Di-n-octyl phthalate	45.8		2.50	10.0	ug/L
205-99-2	Benzo(b)fluoranthene	44.0		1.10	5.00	ug/L

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207-08-9	Benzo(k)fluoranthene	45.2		1.20	5.00	ug/L
50-32-8	Benzo(a)pyrene	47.5		1.70	5.00	ug/L
193-39-5	Indeno(1,2,3-cd)pyrene	45.3		1.00	5.00	ug/L
53-70-3	Dibenzo(a,h)anthracene	45.0		1.20	5.00	ug/L
191-24-2	Benzo(g,h,i)perylene	41.8		1.20	5.00	ug/L
95-94-3	1,2,4,5-Tetrachlorobenzene	49.2		1.10	5.00	ug/L
123-91-1	1,4-Dioxane	39.6		1.30	5.00	ug/L
58-90-2	2,3,4,6-Tetrachlorophenol	48.7		0.79	5.00	ug/L
SURROGATES						
367-12-4	2-Fluorophenol	128		15 (10) - 110 (139)	85%	SPK: 150
13127-88-3	Phenol-d6	122		15 (10) - 110 (134)	81%	SPK: 150
4165-60-0	Nitrobenzene-d5	109		30 (49) - 130 (133)	109%	SPK: 100
321-60-8	2-Fluorobiphenyl	94.0		30 (52) - 130 (132)	94%	SPK: 100
118-79-6	2,4,6-Tribromophenol	174	*	15 (44) - 110 (137)	116%	SPK: 150
1718-51-0	Terphenyl-d14	95.1		30 (48) - 130 (125)	95%	SPK: 100
INTERNAL STANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	219000	6.887			
1146-65-2	Naphthalene-d8	851000	8.169			
15067-26-2	Acenaphthene-d10	458000	9.922			
1517-22-2	Phenanthrene-d10	790000	11.404			
1719-03-5	Chrysene-d12	569000	14.045			
1520-96-3	Perylene-d12	469000	15.516			

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