

Report of Analysis

Client:	PSEG	Date Collected:	08/07/25
Project:	PSEG Metro Way	Date Received:	08/07/25
Client Sample ID:	TP-6	SDG No.:	Q2798
Lab Sample ID:	Q2798-01	Matrix:	SOIL
Analytical Method:	8270E	% Solid:	89.2
Sample Wt/Vol:	50.02	Units:	g
Soil Aliquot Vol:			uL
Extraction Type :		Decanted :	N
Injection Volume :		GPC Factor :	1.0
Prep Method :	SW3541	Level :	LOW
		Test:	SVOC-TCL BNA -20
		GPC Cleanup :	N
		PH :	

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BP025369.D	2	08/08/25 09:12	08/12/25 16:37	PB169176

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
100-52-7	Benzaldehyde	0.44	U	0.21	0.44	mg/Kg
108-95-2	Phenol	0.23	U	0.030	0.23	mg/Kg
111-44-4	bis(2-Chloroethyl)ether	0.23	U	0.033	0.23	mg/Kg
95-57-8	2-Chlorophenol	0.23	U	0.033	0.23	mg/Kg
95-48-7	2-Methylphenol	0.23	U	0.040	0.23	mg/Kg
108-60-1	2,2-oxybis(1-Chloropropane)	0.23	U	0.050	0.23	mg/Kg
98-86-2	Acetophenone	0.23	U	0.040	0.23	mg/Kg
65794-96-9	3+4-Methylphenols	0.44	U	0.055	0.44	mg/Kg
621-64-7	n-Nitroso-di-n-propylamine	0.11	U	0.064	0.11	mg/Kg
67-72-1	Hexachloroethane	0.23	U	0.024	0.23	mg/Kg
98-95-3	Nitrobenzene	0.23	U	0.025	0.23	mg/Kg
78-59-1	Isophorone	0.23	U	0.044	0.23	mg/Kg
88-75-5	2-Nitrophenol	0.23	U	0.078	0.23	mg/Kg
105-67-9	2,4-Dimethylphenol	0.23	U	0.087	0.23	mg/Kg
111-91-1	bis(2-Chloroethoxy)methane	0.23	U	0.041	0.23	mg/Kg
120-83-2	2,4-Dichlorophenol	0.23	U	0.038	0.23	mg/Kg
91-20-3	Naphthalene	0.23	U	0.031	0.23	mg/Kg
106-47-8	4-Chloroaniline	0.23	U	0.048	0.23	mg/Kg
87-68-3	Hexachlorobutadiene	0.23	U	0.034	0.23	mg/Kg
105-60-2	Caprolactam	0.44	U	0.070	0.44	mg/Kg
59-50-7	4-Chloro-3-methylphenol	0.23	U	0.039	0.23	mg/Kg
91-57-6	2-Methylnaphthalene	0.23	U	0.034	0.23	mg/Kg
77-47-4	Hexachlorocyclopentadiene	0.44	U	0.16	0.44	mg/Kg
88-06-2	2,4,6-Trichlorophenol	0.23	U	0.027	0.23	mg/Kg
95-95-4	2,4,5-Trichlorophenol	0.23	U	0.039	0.23	mg/Kg
92-52-4	1,1-Biphenyl	0.23	U	0.029	0.23	mg/Kg
91-58-7	2-Chloronaphthalene	0.23	U	0.030	0.23	mg/Kg
88-74-4	2-Nitroaniline	0.23	U	0.065	0.23	mg/Kg
131-11-3	Dimethylphthalate	0.23	U	0.036	0.23	mg/Kg

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Injection Volume :		Level :	LOW
		GPC Factor :	1.0
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208-96-8	Acenaphthylene	0.23	U	0.039	0.23	mg/Kg
606-20-2	2,6-Dinitrotoluene	0.23	U	0.045	0.23	mg/Kg
99-09-2	3-Nitroaniline	0.23	U	0.062	0.23	mg/Kg
83-32-9	Acenaphthene	0.23	U	0.029	0.23	mg/Kg
51-28-5	2,4-Dinitrophenol	0.44	U	0.31	0.44	mg/Kg
100-02-7	4-Nitrophenol	0.44	U	0.14	0.44	mg/Kg
132-64-9	Dibenzofuran	0.23	U	0.031	0.23	mg/Kg
121-14-2	2,4-Dinitrotoluene	0.23	U	0.067	0.23	mg/Kg
84-66-2	Diethylphthalate	0.23	U	0.038	0.23	mg/Kg
7005-72-3	4-Chlorophenyl-phenylether	0.23	U	0.036	0.23	mg/Kg
86-73-7	Fluorene	0.23	U	0.034	0.23	mg/Kg
100-01-6	4-Nitroaniline	0.23	U	0.086	0.23	mg/Kg
534-52-1	4,6-Dinitro-2-methylphenol	0.44	U	0.14	0.44	mg/Kg
86-30-6	n-Nitrosodiphenylamine	0.23	U	0.044	0.23	mg/Kg
101-55-3	4-Bromophenyl-phenylether	0.23	U	0.037	0.23	mg/Kg
118-74-1	Hexachlorobenzene	0.23	U	0.034	0.23	mg/Kg
1912-24-9	Atrazine	0.23	U	0.046	0.23	mg/Kg
87-86-5	Pentachlorophenol	0.44	U	0.069	0.44	mg/Kg
85-01-8	Phenanthrene	0.23	U	0.028	0.23	mg/Kg
120-12-7	Anthracene	0.23	U	0.045	0.23	mg/Kg
86-74-8	Carbazole	0.23	U	0.042	0.23	mg/Kg
84-74-2	Di-n-butylphthalate	0.23	U	0.064	0.23	mg/Kg
206-44-0	Fluoranthene	0.23	U	0.040	0.23	mg/Kg
129-00-0	Pyrene	0.23	U	0.048	0.23	mg/Kg
85-68-7	Butylbenzylphthalate	0.23	U	0.096	0.23	mg/Kg
91-94-1	3,3-Dichlorobenzidine	0.44	U	0.049	0.44	mg/Kg
56-55-3	Benzo(a)anthracene	0.23	U	0.031	0.23	mg/Kg
218-01-9	Chrysene	0.23	U	0.027	0.23	mg/Kg
117-81-7	Bis(2-ethylhexyl)phthalate	0.23	U	0.080	0.23	mg/Kg
117-84-0	Di-n-octyl phthalate	0.44	U	0.12	0.44	mg/Kg
205-99-2	Benzo(b)fluoranthene	0.23	U	0.026	0.23	mg/Kg

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Soil Aliquot Vol:		Final Vol:	1000 uL
Extraction Type :		Test:	SVOC-TCL BNA -20
	Decanted :	Level :	LOW
Injection Volume :	GPC Factor :	1.0	GPC Cleanup : N PH :
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207-08-9	Benzo(k)fluoranthene	0.23	U	0.030	0.23	mg/Kg
50-32-8	Benzo(a)pyrene	0.23	U	0.040	0.23	mg/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	0.23	U	0.039	0.23	mg/Kg
53-70-3	Dibenzo(a,h)anthracene	0.23	U	0.037	0.23	mg/Kg
191-24-2	Benzo(g,h,i)perylene	0.23	U	0.035	0.23	mg/Kg
95-94-3	1,2,4,5-Tetrachlorobenzene	0.23	U	0.034	0.23	mg/Kg
123-91-1	1,4-Dioxane	0.23	U	0.061	0.23	mg/Kg
58-90-2	2,3,4,6-Tetrachlorophenol	0.23	U	0.037	0.23	mg/Kg

SURROGATES

367-12-4	2-Fluorophenol	68.3		18 - 112	46%	SPK: 150
13127-88-3	Phenol-d6	64.9		15 - 107	43%	SPK: 150
4165-60-0	Nitrobenzene-d5	45.6		18 - 107	46%	SPK: 100
321-60-8	2-Fluorobiphenyl	40.3		20 - 109	40%	SPK: 100
118-79-6	2,4,6-Tribromophenol	78.5		10 - 116	52%	SPK: 150
1718-51-0	Terphenyl-d14	36.0		10 - 105	36%	SPK: 100

INTERNAL STANDARDS

3855-82-1	1,4-Dichlorobenzene-d4	200000	7.79	
1146-65-2	Naphthalene-d8	790000	10.566	
15067-26-2	Acenaphthene-d10	469000	14.413	
1517-22-2	Phenanthrene-d10	873000	17.225	
1719-03-5	Chrysene-d12	1040000	21.666	
1520-96-3	Perylene-d12	1220000	25.066	

TENTATIVE IDENTIFIED COMPOUNDS

000994-05-8	Butane, 2-methoxy-2-methyl-	600	J	3.04	ug/Kg
000112-79-8	9-Octadecenoic acid, (E)-	330	J	13.9	ug/Kg
000057-11-4	Octadecanoic acid	320	J	14.6	ug/Kg
000544-63-8	Tetradecanoic acid	340	J	14.7	ug/Kg
001002-84-2	Pentadecanoic acid	1700	J	14.7	ug/Kg
000638-53-9	Tridecanoic acid	6500	J	14.8	ug/Kg
002027-47-6	9-Octadecenoic acid	850	J	15.4	ug/Kg

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000057-10-3	n-Hexadecanoic acid	1500	J		15.8	ug/Kg
	unknown15.854	6500	J		15.9	ug/Kg
032811-40-8	(E)-4-(3-Hydroxyprop-1-en-1-yl)-2-	330	J		16.6	ug/Kg
082304-66-3	7,9-Di-tert-butyl-1-oxaspiro(4,5)d	570	J		17.9	ug/Kg
	unknown18.178	1100	J		18.2	ug/Kg
087345-53-7	3,5-Dimethoxy-4-hydroxycinnamaldehyde	740	J		18.4	ug/Kg
104330-63-4	cis-Sinapyl alcohol	570	J		18.5	ug/Kg
	unknown18.601	210	J		18.6	ug/Kg
000124-26-5	Octadecanamide	280	J		18.8	ug/Kg
000106-11-6	Octadecanoic acid, 2-(2-hydroxyethyl)	400	J		19.5	ug/Kg
000593-49-7	Heptacosane	220	J		21.4	ug/Kg
006422-86-2	1,4-Benzenedicarboxylic acid, bis(	2700	J		21.8	ug/Kg
1000356-27-6	Terephthalic acid, 3-methylbut-2-yl	1400	J		21.9	ug/Kg

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