

Report of Analysis

Client:	PSEG	Date Collected:	04/11/25
Project:	Central (Harrison) Gas Division Service MA00006789	Date Received:	04/11/25
Client Sample ID:	HR-03-041125	SDG No.:	Q1785
Lab Sample ID:	Q1785-03	Matrix:	SOIL
Analytical Method:	SW8270	% Solid:	94.7
Sample Wt/Vol:	50.06	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
BM049940.D	2	04/14/25 09:40	04/15/25 18:32	PB167577

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
100-52-7	Benzaldehyde	0.20	U	0.20	0.42	mg/Kg
108-95-2	Phenol	0.028	U	0.028	0.22	mg/Kg
111-44-4	bis(2-Chloroethyl)ether	0.031	U	0.031	0.22	mg/Kg
95-57-8	2-Chlorophenol	0.031	U	0.031	0.22	mg/Kg
95-48-7	2-Methylphenol	0.038	U	0.038	0.22	mg/Kg
108-60-1	2,2-oxybis(1-Chloropropane)	0.048	U	0.048	0.22	mg/Kg
98-86-2	Acetophenone	0.037	U	0.037	0.22	mg/Kg
65794-96-9	3+4-Methylphenols	0.052	U	0.052	0.42	mg/Kg
621-64-7	n-Nitroso-di-n-propylamine	0.060	U	0.060	0.10	mg/Kg
67-72-1	Hexachloroethane	0.022	U	0.022	0.22	mg/Kg
98-95-3	Nitrobenzene	0.023	U	0.023	0.22	mg/Kg
78-59-1	Isophorone	0.042	U	0.042	0.22	mg/Kg
88-75-5	2-Nitrophenol	0.074	U	0.074	0.22	mg/Kg
105-67-9	2,4-Dimethylphenol	0.082	U	0.082	0.22	mg/Kg
111-91-1	bis(2-Chloroethoxy)methane	0.039	U	0.039	0.22	mg/Kg
120-83-2	2,4-Dichlorophenol	0.036	U	0.036	0.22	mg/Kg
91-20-3	Naphthalene	0.029	U	0.029	0.22	mg/Kg
106-47-8	4-Chloroaniline	0.045	U	0.045	0.22	mg/Kg
87-68-3	Hexachlorobutadiene	0.032	U	0.032	0.22	mg/Kg
105-60-2	Caprolactam	0.066	U	0.066	0.42	mg/Kg
59-50-7	4-Chloro-3-methylphenol	0.036	U	0.036	0.22	mg/Kg
91-57-6	2-Methylnaphthalene	0.032	U	0.032	0.22	mg/Kg
77-47-4	Hexachlorocyclopentadiene	0.15	U	0.15	0.42	mg/Kg
88-06-2	2,4,6-Trichlorophenol	0.025	U	0.025	0.22	mg/Kg
95-95-4	2,4,5-Trichlorophenol	0.037	U	0.037	0.22	mg/Kg
92-52-4	1,1-Biphenyl	0.028	U	0.028	0.22	mg/Kg
91-58-7	2-Chloronaphthalene	0.029	U	0.029	0.22	mg/Kg
88-74-4	2-Nitroaniline	0.061	U	0.061	0.22	mg/Kg
131-11-3	Dimethylphthalate	0.034	U	0.034	0.22	mg/Kg

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208-96-8	Acenaphthylene	0.037	U	0.037	0.22	mg/Kg
606-20-2	2,6-Dinitrotoluene	0.043	U	0.043	0.22	mg/Kg
99-09-2	3-Nitroaniline	0.058	U	0.058	0.22	mg/Kg
83-32-9	Acenaphthene	0.027	U	0.027	0.22	mg/Kg
51-28-5	2,4-Dinitrophenol	0.29	U	0.29	0.42	mg/Kg
100-02-7	4-Nitrophenol	0.14	U	0.14	0.42	mg/Kg
132-64-9	Dibenzofuran	0.029	U	0.029	0.22	mg/Kg
121-14-2	2,4-Dinitrotoluene	0.063	U	0.063	0.22	mg/Kg
84-66-2	Diethylphthalate	0.036	U	0.036	0.22	mg/Kg
7005-72-3	4-Chlorophenyl-phenylether	0.034	U	0.034	0.22	mg/Kg
86-73-7	Fluorene	0.032	U	0.032	0.22	mg/Kg
100-01-6	4-Nitroaniline	0.081	U	0.081	0.22	mg/Kg
534-52-1	4,6-Dinitro-2-methylphenol	0.13	U	0.13	0.42	mg/Kg
86-30-6	n-Nitrosodiphenylamine	0.042	U	0.042	0.22	mg/Kg
101-55-3	4-Bromophenyl-phenylether	0.035	U	0.035	0.22	mg/Kg
118-74-1	Hexachlorobenzene	0.032	U	0.032	0.22	mg/Kg
1912-24-9	Atrazine	0.043	U	0.043	0.22	mg/Kg
87-86-5	Pentachlorophenol	0.065	U	0.065	0.42	mg/Kg
85-01-8	Phenanthrene	0.027	U	0.027	0.22	mg/Kg
120-12-7	Anthracene	0.042	U	0.042	0.22	mg/Kg
86-74-8	Carbazole	0.040	U	0.040	0.22	mg/Kg
84-74-2	Di-n-butylphthalate	0.061	U	0.061	0.22	mg/Kg
206-44-0	Fluoranthene	0.038	U	0.038	0.22	mg/Kg
129-00-0	Pyrene	0.046	U	0.046	0.22	mg/Kg
85-68-7	Butylbenzylphthalate	0.090	U	0.090	0.22	mg/Kg
91-94-1	3,3-Dichlorobenzidine	0.046	U	0.046	0.42	mg/Kg
56-55-3	Benzo(a)anthracene	0.029	U	0.029	0.22	mg/Kg
218-01-9	Chrysene	0.025	U	0.025	0.22	mg/Kg
117-81-7	Bis(2-ethylhexyl)phthalate	0.075	U	0.075	0.22	mg/Kg
117-84-0	Di-n-octyl phthalate	0.11	U	0.11	0.42	mg/Kg
205-99-2	Benzo(b)fluoranthene	0.024	U	0.024	0.22	mg/Kg

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207-08-9	Benzo(k)fluoranthene	0.028	U	0.028	0.22	mg/Kg
50-32-8	Benzo(a)pyrene	0.037	U	0.037	0.22	mg/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	0.037	U	0.037	0.22	mg/Kg
53-70-3	Dibenzo(a,h)anthracene	0.035	U	0.035	0.22	mg/Kg
191-24-2	Benzo(g,h,i)perylene	0.033	U	0.033	0.22	mg/Kg
95-94-3	1,2,4,5-Tetrachlorobenzene	0.032	U	0.032	0.22	mg/Kg
123-91-1	1,4-Dioxane	0.057	U	0.057	0.22	mg/Kg
58-90-2	2,3,4,6-Tetrachlorophenol	0.035	U	0.035	0.22	mg/Kg

SURROGATES

367-12-4	2-Fluorophenol	76.1		18 - 112	51%	SPK: 150
13127-88-3	Phenol-d6	72.7		15 - 107	48%	SPK: 150
4165-60-0	Nitrobenzene-d5	52.6		18 - 107	53%	SPK: 100
321-60-8	2-Fluorobiphenyl	53.7		20 - 109	54%	SPK: 100
118-79-6	2,4,6-Tribromophenol	81.7		10 - 116	54%	SPK: 150
1718-51-0	Terphenyl-d14	55.2		10 - 105	55%	SPK: 100

INTERNAL STANDARDS

3855-82-1	1,4-Dichlorobenzene-d4	419000	7.769
1146-65-2	Naphthalene-d8	1390000	10.569
15067-26-2	Acenaphthene-d10	848000	14.416
1517-22-2	Phenanthrene-d10	1650000	17.162
1719-03-5	Chrysene-d12	1620000	21.397
1520-96-3	Perylene-d12	1770000	24.409

TENTATIVE IDENTIFIED COMPOUNDS

001120-21-4	Undecane	740	J	9.00	ug/Kg
000629-50-5	Tridecane	970	J	12.0	ug/Kg
000629-59-4	Tetradecane	1500	J	13.3	ug/Kg
020027-77-4	Naphthalene, 1,2,3,4-tetrahydro-5,	530	J	13.4	ug/Kg
055045-10-8	Tridecane, 6-propyl-	570	J	13.9	ug/Kg
000629-62-9	Pentadecane	1800	J	14.3	ug/Kg
007098-21-7	Tritetracontane	580	J	14.9	ug/Kg

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000544-76-3	Hexadecane	1800	J		15.3	ug/Kg
003892-00-0	Pentadecane, 2,6,10-trimethyl-	1000	J		15.7	ug/Kg
017312-55-9	Decane, 3,8-dimethyl-	550	J		15.8	ug/Kg
000629-78-7	Heptadecane	3400	J		16.1	ug/Kg
000593-45-3	Octadecane	2600	J		16.9	ug/Kg
000629-93-6	Octadecane, 1-iodo-	1500	J		17.0	ug/Kg
000629-94-7	Heneicosane	2400	J		17.7	ug/Kg
000629-92-5	Nonadecane	2000	J		18.4	ug/Kg
000112-95-8	Eicosane	1600	J		19.0	ug/Kg
000057-11-4	Octadecanoic acid	720	J		19.3	ug/Kg
035599-77-0	Tridecane, 1-iodo-	1200	J		19.6	ug/Kg
000630-01-3	Hexacosane	850	J		20.1	ug/Kg
000630-02-4	Octacosane	530	J		20.6	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