

Report of Analysis

Client:	PSEG	Date Collected:	03/24/25
Project:	Central (Harrison) Gas Division Service MA00006789	Date Received:	03/24/25
Client Sample ID:	HR-01-032425	SDG No.:	Q1634
Lab Sample ID:	Q1634-01	Matrix:	SOIL
Analytical Method:	SW8270	% Solid:	91.1
Sample Wt/Vol:	50.03	Units:	g
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 :
Prep Method :	SW3541		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BF142147.D	5	03/25/25 08:50	03/28/25 02:06	PB167291

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
100-52-7	Benzaldehyde	0.51	U	0.51	1.10	mg/Kg
108-95-2	Phenol	0.073	U	0.073	0.56	mg/Kg
111-44-4	bis(2-Chloroethyl)ether	0.080	U	0.080	0.56	mg/Kg
95-57-8	2-Chlorophenol	0.080	U	0.080	0.56	mg/Kg
95-48-7	2-Methylphenol	0.098	U	0.098	0.56	mg/Kg
108-60-1	2,2-oxybis(1-Chloropropane)	0.12	U	0.12	0.56	mg/Kg
98-86-2	Acetophenone	0.097	U	0.097	0.56	mg/Kg
65794-96-9	3+4-Methylphenols	0.14	U	0.14	1.10	mg/Kg
621-64-7	n-Nitroso-di-n-propylamine	0.16	U	0.16	0.26	mg/Kg
67-72-1	Hexachloroethane	0.058	U	0.058	0.56	mg/Kg
98-95-3	Nitrobenzene	0.060	U	0.060	0.56	mg/Kg
78-59-1	Isophorone	0.11	U	0.11	0.56	mg/Kg
88-75-5	2-Nitrophenol	0.19	U	0.19	0.56	mg/Kg
105-67-9	2,4-Dimethylphenol	0.21	U	0.21	0.56	mg/Kg
111-91-1	bis(2-Chloroethoxy)methane	0.10	U	0.10	0.56	mg/Kg
120-83-2	2,4-Dichlorophenol	0.093	U	0.093	0.56	mg/Kg
91-20-3	Naphthalene	0.075	U	0.075	0.56	mg/Kg
106-47-8	4-Chloroaniline	0.12	U	0.12	0.56	mg/Kg
87-68-3	Hexachlorobutadiene	0.083	U	0.083	0.56	mg/Kg
105-60-2	Caprolactam	0.17	U	0.17	1.10	mg/Kg
59-50-7	4-Chloro-3-methylphenol	0.095	U	0.095	0.56	mg/Kg
91-57-6	2-Methylnaphthalene	0.084	U	0.084	0.56	mg/Kg
77-47-4	Hexachlorocyclopentadiene	0.38	U	0.38	1.10	mg/Kg
88-06-2	2,4,6-Trichlorophenol	0.065	U	0.065	0.56	mg/Kg
95-95-4	2,4,5-Trichlorophenol	0.096	U	0.096	0.56	mg/Kg
92-52-4	1,1-Biphenyl	0.072	U	0.072	0.56	mg/Kg
91-58-7	2-Chloronaphthalene	0.074	U	0.074	0.56	mg/Kg
88-74-4	2-Nitroaniline	0.16	U	0.16	0.56	mg/Kg
131-11-3	Dimethylphthalate	0.089	U	0.089	0.56	mg/Kg

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208-96-8	Acenaphthylene	0.095	U	0.095	0.56	mg/Kg
606-20-2	2,6-Dinitrotoluene	0.11	U	0.11	0.56	mg/Kg
99-09-2	3-Nitroaniline	0.15	U	0.15	0.56	mg/Kg
83-32-9	Acenaphthene	0.070	U	0.070	0.56	mg/Kg
51-28-5	2,4-Dinitrophenol	0.75	U	0.75	1.10	mg/Kg
100-02-7	4-Nitrophenol	0.35	U	0.35	1.10	mg/Kg
132-64-9	Dibenzofuran	0.075	U	0.075	0.56	mg/Kg
121-14-2	2,4-Dinitrotoluene	0.16	U	0.16	0.56	mg/Kg
84-66-2	Diethylphthalate	0.093	U	0.093	0.56	mg/Kg
7005-72-3	4-Chlorophenyl-phenylether	0.088	U	0.088	0.56	mg/Kg
86-73-7	Fluorene	0.083	U	0.083	0.56	mg/Kg
100-01-6	4-Nitroaniline	0.21	U	0.21	0.56	mg/Kg
534-52-1	4,6-Dinitro-2-methylphenol	0.34	U	0.34	1.10	mg/Kg
86-30-6	n-Nitrosodiphenylamine	0.11	U	0.11	0.56	mg/Kg
101-55-3	4-Bromophenyl-phenylether	0.092	U	0.092	0.56	mg/Kg
118-74-1	Hexachlorobenzene	0.083	U	0.083	0.56	mg/Kg
1912-24-9	Atrazine	0.11	U	0.11	0.56	mg/Kg
87-86-5	Pentachlorophenol	0.17	U	0.17	1.10	mg/Kg
85-01-8	Phenanthrene	0.069	U	0.069	0.56	mg/Kg
120-12-7	Anthracene	0.11	U	0.11	0.56	mg/Kg
86-74-8	Carbazole	0.10	U	0.10	0.56	mg/Kg
84-74-2	Di-n-butylphthalate	0.16	U	0.16	0.56	mg/Kg
206-44-0	Fluoranthene	0.099	U	0.099	0.56	mg/Kg
129-00-0	Pyrene	0.12	U	0.12	0.56	mg/Kg
85-68-7	Butylbenzylphthalate	0.23	U	0.23	0.56	mg/Kg
91-94-1	3,3-Dichlorobenzidine	0.12	U	0.12	1.10	mg/Kg
56-55-3	Benzo(a)anthracene	0.076	U	0.076	0.56	mg/Kg
218-01-9	Chrysene	0.066	U	0.066	0.56	mg/Kg
117-81-7	Bis(2-ethylhexyl)phthalate	0.19	U	0.19	0.56	mg/Kg
117-84-0	Di-n-octyl phthalate	0.29	U	0.29	1.10	mg/Kg
205-99-2	Benzo(b)fluoranthene	0.063	U	0.063	0.56	mg/Kg

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207-08-9	Benzo(k)fluoranthene	0.074	U	0.074	0.56	mg/Kg
50-32-8	Benzo(a)pyrene	0.097	U	0.097	0.56	mg/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	0.096	U	0.096	0.56	mg/Kg
53-70-3	Dibenzo(a,h)anthracene	0.090	U	0.090	0.56	mg/Kg
191-24-2	Benzo(g,h,i)perylene	0.085	U	0.085	0.56	mg/Kg
95-94-3	1,2,4,5-Tetrachlorobenzene	0.084	U	0.084	0.56	mg/Kg
123-91-1	1,4-Dioxane	0.15	U	0.15	0.56	mg/Kg
58-90-2	2,3,4,6-Tetrachlorophenol	0.090	U	0.090	0.56	mg/Kg

SURROGATES

367-12-4	2-Fluorophenol	100		18 - 112	67%	SPK: 150
13127-88-3	Phenol-d6	95.8		15 - 107	64%	SPK: 150
4165-60-0	Nitrobenzene-d5	78.1		18 - 107	78%	SPK: 100
321-60-8	2-Fluorobiphenyl	80.0		20 - 109	80%	SPK: 100
118-79-6	2,4,6-Tribromophenol	73.1		10 - 116	49%	SPK: 150
1718-51-0	Terphenyl-d14	51.4		10 - 105	51%	SPK: 100

INTERNAL STANDARDS

3855-82-1	1,4-Dichlorobenzene-d4	134000	6.869	
1146-65-2	Naphthalene-d8	468000	8.151	
15067-26-2	Acenaphthene-d10	220000	9.91	
1517-22-2	Phenanthrene-d10	312000	11.398	
1719-03-5	Chrysene-d12	284000	14.033	
1520-96-3	Perylene-d12	203000	15.509	

TENTATIVE IDENTIFIED COMPOUNDS

001074-43-7	Benzene, 1-methyl-3-propyl-	1000	J	7.14	ug/Kg
000105-05-5	Benzene, 1,4-diethyl-	910	J	7.19	ug/Kg
000629-50-5	Tridecane	1200	J	8.74	ug/Kg
	unknown9.104	1100	J	9.10	ug/Kg
003891-98-3	Dodecane, 2,6,10-trimethyl-	1600	J	9.17	ug/Kg
1000382-54-3	Carbonic acid, eicosyl vinyl ester	4500	J	9.31	ug/Kg
	unknown9.563	1500	J	9.56	ug/Kg

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1000342-70-4	1-Octadecanesulphonyl chloride	3000	J		9.63	ug/Kg
000629-62-9	Pentadecane	4000	J		9.84	ug/Kg
001560-78-7	2-Methyltetracosane	1100	J		10.2	ug/Kg
000544-76-3	Hexadecane	4300	J		10.3	ug/Kg
003892-00-0	Pentadecane, 2,6,10-trimethyl-	2300	J		10.6	ug/Kg
000112-40-3	Dodecane	8300	J		10.8	ug/Kg
000593-45-3	Octadecane	4500	J		11.3	ug/Kg
1000432-25-9	3-Ethyl-2,6,10-trimethylundecane	2100	J		11.3	ug/Kg
000646-31-1	Tetracosane	3400	J		11.7	ug/Kg
000629-78-7	Heptadecane	2800	J		12.1	ug/Kg
000629-94-7	Heneicosane	2200	J		12.5	ug/Kg
035599-77-0	Tridecane, 1-iodo-	1800	J		12.9	ug/Kg
000638-67-5	Tricosane	1100	J		13.2	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