

Report of Analysis

Client:	PSEG	Date Collected:	04/11/25
Project:	Summit Gas and Appliance Service MA00006789	Date Received:	04/11/25
Client Sample ID:	SU-03-41125	SDG No.:	Q1783
Lab Sample ID:	Q1783-01	Matrix:	SOIL
Analytical Method:	SW8270	% Solid:	89.8
Sample Wt/Vol:	50.07 Units: g	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 :	SW3541		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
BM049924.D	5	04/14/25 09:40	04/14/25 21:43	PB167577

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
100-52-7	Benzaldehyde	0.52	U	0.52	1.10	mg/Kg
108-95-2	Phenol	0.074	U	0.074	0.57	mg/Kg
111-44-4	bis(2-Chloroethyl)ether	0.081	U	0.081	0.57	mg/Kg
95-57-8	2-Chlorophenol	0.081	U	0.081	0.57	mg/Kg
95-48-7	2-Methylphenol	0.100	U	0.100	0.57	mg/Kg
108-60-1	2,2-oxybis(1-Chloropropane)	0.13	U	0.13	0.57	mg/Kg
98-86-2	Acetophenone	0.098	U	0.098	0.57	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	0.27	mg/Kg
67-72-1	Hexachloroethane	0.059	U	0.059	0.57	mg/Kg
98-95-3	Nitrobenzene	0.061	U	0.061	0.57	mg/Kg
78-59-1	Isophorone	0.11	U	0.11	0.57	mg/Kg
88-75-5	2-Nitrophenol	0.19	U	0.19	0.57	mg/Kg
105-67-9	2,4-Dimethylphenol	0.22	U	0.22	0.57	mg/Kg
111-91-1	bis(2-Chloroethoxy)methane	0.10	U	0.10	0.57	mg/Kg
120-83-2	2,4-Dichlorophenol	0.094	U	0.094	0.57	mg/Kg
91-20-3	Naphthalene	0.076	U	0.076	0.57	mg/Kg
106-47-8	4-Chloroaniline	0.12	U	0.12	0.57	mg/Kg
87-68-3	Hexachlorobutadiene	0.084	U	0.084	0.57	mg/Kg
105-60-2	Caprolactam	0.17	U	0.17	1.10	mg/Kg
59-50-7	4-Chloro-3-methylphenol	0.096	U	0.096	0.57	mg/Kg
91-57-6	2-Methylnaphthalene	0.44	J	0.085	0.57	mg/Kg
77-47-4	Hexachlorocyclopentadiene	0.39	U	0.39	1.10	mg/Kg
88-06-2	2,4,6-Trichlorophenol	0.066	U	0.066	0.57	mg/Kg
95-95-4	2,4,5-Trichlorophenol	0.097	U	0.097	0.57	mg/Kg
92-52-4	1,1-Biphenyl	0.073	U	0.073	0.57	mg/Kg
91-58-7	2-Chloronaphthalene	0.075	U	0.075	0.57	mg/Kg
88-74-4	2-Nitroaniline	0.16	U	0.16	0.57	mg/Kg
131-11-3	Dimethylphthalate	0.090	U	0.090	0.57	mg/Kg

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208-96-8	Acenaphthylene	0.096	U	0.096	0.57	mg/Kg
606-20-2	2,6-Dinitrotoluene	0.11	U	0.11	0.57	mg/Kg
99-09-2	3-Nitroaniline	0.15	U	0.15	0.57	mg/Kg
83-32-9	Acenaphthene	0.071	U	0.071	0.57	mg/Kg
51-28-5	2,4-Dinitrophenol	0.76	U	0.76	1.10	mg/Kg
100-02-7	4-Nitrophenol	0.36	U	0.36	1.10	mg/Kg
132-64-9	Dibenzofuran	0.076	U	0.076	0.57	mg/Kg
121-14-2	2,4-Dinitrotoluene	0.17	U	0.17	0.57	mg/Kg
84-66-2	Diethylphthalate	0.094	U	0.094	0.57	mg/Kg
7005-72-3	4-Chlorophenyl-phenylether	0.089	U	0.089	0.57	mg/Kg
86-73-7	Fluorene	0.28	J	0.084	0.57	mg/Kg
100-01-6	4-Nitroaniline	0.21	U	0.21	0.57	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	0.57	mg/Kg
101-55-3	4-Bromophenyl-phenylether	0.093	U	0.093	0.57	mg/Kg
118-74-1	Hexachlorobenzene	0.084	U	0.084	0.57	mg/Kg
1912-24-9	Atrazine	0.11	U	0.11	0.57	mg/Kg
87-86-5	Pentachlorophenol	0.17	U	0.17	1.10	mg/Kg
85-01-8	Phenanthrene	1.20		0.070	0.57	mg/Kg
120-12-7	Anthracene	0.35	J	0.11	0.57	mg/Kg
86-74-8	Carbazole	0.10	U	0.10	0.57	mg/Kg
84-74-2	Di-n-butylphthalate	0.16	U	0.16	0.57	mg/Kg
206-44-0	Fluoranthene	1.20		0.10	0.57	mg/Kg
129-00-0	Pyrene	1.10		0.12	0.57	mg/Kg
85-68-7	Butylbenzylphthalate	0.24	U	0.24	0.57	mg/Kg
91-94-1	3,3-Dichlorobenzidine	0.12	U	0.12	1.10	mg/Kg
56-55-3	Benzo(a)anthracene	0.52	J	0.077	0.57	mg/Kg
218-01-9	Chrysene	0.50	J	0.066	0.57	mg/Kg
117-81-7	Bis(2-ethylhexyl)phthalate	0.20	U	0.20	0.57	mg/Kg
117-84-0	Di-n-octyl phthalate	0.29	U	0.29	1.10	mg/Kg
205-99-2	Benzo(b)fluoranthene	0.56	J	0.063	0.57	mg/Kg

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207-08-9	Benzo(k)fluoranthene	0.075	U	0.075	0.57	mg/Kg
50-32-8	Benzo(a)pyrene	0.49	J	0.098	0.57	mg/Kg
193-39-5	Indeno(1,2,3-cd)pyrene	0.23	J	0.097	0.57	mg/Kg
53-70-3	Dibenzo(a,h)anthracene	0.091	U	0.091	0.57	mg/Kg
191-24-2	Benzo(g,h,i)perylene	0.26	J	0.086	0.57	mg/Kg
95-94-3	1,2,4,5-Tetrachlorobenzene	0.085	U	0.085	0.57	mg/Kg
123-91-1	1,4-Dioxane	0.15	U	0.15	0.57	mg/Kg
58-90-2	2,3,4,6-Tetrachlorophenol	0.091	U	0.091	0.57	mg/Kg

SURROGATES

367-12-4	2-Fluorophenol	50.5		18 - 112	34%	SPK: 150
13127-88-3	Phenol-d6	47.1		15 - 107	31%	SPK: 150
4165-60-0	Nitrobenzene-d5	38.8		18 - 107	39%	SPK: 100
321-60-8	2-Fluorobiphenyl	35.4		20 - 109	35%	SPK: 100
118-79-6	2,4,6-Tribromophenol	47.6		10 - 116	32%	SPK: 150
1718-51-0	Terphenyl-d14	34.6		10 - 105	35%	SPK: 100

INTERNAL STANDARDS

3855-82-1	1,4-Dichlorobenzene-d4	442000	7.775
1146-65-2	Naphthalene-d8	1450000	10.569
15067-26-2	Acenaphthene-d10	901000	14.416
1517-22-2	Phenanthrene-d10	1750000	17.163
1719-03-5	Chrysene-d12	1790000	21.398
1520-96-3	Perylene-d12	2010000	24.397

TENTATIVE IDENTIFIED COMPOUNDS

000124-18-5	Decane	1700	J	7.40	ug/Kg
001074-43-7	Benzene, 1-methyl-3-propyl-	840	J	8.32	ug/Kg
1000382-90-9	Carbonic acid, prop-1-en-2-yl tetr	2800	J	9.00	ug/Kg
001595-16-0	Benzene, 1-methyl-4-(1-methylpropy	1400	J	9.13	ug/Kg
000629-50-5	Tridecane	1700	J	12.0	ug/Kg
90-12-0	1-Methylnaphthalene	330	J	12.5	ug/Kg
017302-32-8	Nonane, 3,7-dimethyl-	1200	J	13.0	ug/Kg

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000629-59-4	Tetradecane	3600	J		13.3	ug/Kg
002613-76-5	1H-Indene, 2,3-dihydro-1,1,3-trime	1100	J		13.4	ug/Kg
017312-62-8	Decane, 5-propyl-	1200	J		13.9	ug/Kg
062238-11-3	Decane, 2,3,5-trimethyl-	3900	J		14.3	ug/Kg
000112-89-0	Octadecane, 1-bromo-	1100	J		14.9	ug/Kg
000544-76-3	Hexadecane	3900	J		15.3	ug/Kg
003892-00-0	Pentadecane, 2,6,10-trimethyl-	1900	J		15.7	ug/Kg
000629-92-5	Nonadecane	5400	J		16.1	ug/Kg
002051-30-1	Octane, 2,6-dimethyl-	2800	J		16.2	ug/Kg
000629-94-7	Heneicosane	4400	J		16.9	ug/Kg
000646-31-1	Tetracosane	2400	J		17.0	ug/Kg
055045-08-4	Dodecane, 2-methyl-6-propyl-	4000	J		17.7	ug/Kg
000112-95-8	Eicosane	3300	J		18.4	ug/Kg
000629-78-7	Heptadecane	2600	J		19.0	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