

Report of Analysis

Client:	PSEG	Date Collected:	07/24/25
Project:	Hackensack Sub HQ	Date Received:	07/24/25
Client Sample ID:	299	SDG No.:	Q2691
Lab Sample ID:	Q2691-07	Matrix:	SOIL
Analytical Method:	8260D	% Solid:	89.1
Sample Wt/Vol:	4.2 Units: g	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOC-TCLVOA-10
GC Column:	RXI-624 ID : 0.25	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Date Analyzed	Prep Batch ID
VW031925.D	1	07/24/25 14:35	VW072425

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
75-71-8	Dichlorodifluoromethane	0.0067	U	0.0015	0.0067	mg/Kg
74-87-3	Chloromethane	0.0067	U	0.0015	0.0067	mg/Kg
75-01-4	Vinyl Chloride	0.0067	U	0.0011	0.0067	mg/Kg
74-83-9	Bromomethane	0.0067	U	0.0014	0.0067	mg/Kg
75-00-3	Chloroethane	0.0067	U	0.0017	0.0067	mg/Kg
75-69-4	Trichlorofluoromethane	0.0067	U	0.0016	0.0067	mg/Kg
76-13-1	1,1,2-Trichlorotrifluoroethane	0.0067	U	0.0014	0.0067	mg/Kg
75-35-4	1,1-Dichloroethene	0.0067	U	0.0013	0.0067	mg/Kg
67-64-1	Acetone	0.018	J	0.0063	0.033	mg/Kg
75-15-0	Carbon Disulfide	0.0067	U	0.0014	0.0067	mg/Kg
1634-04-4	Methyl tert-butyl Ether	0.0067	U	0.00098	0.0067	mg/Kg
79-20-9	Methyl Acetate	0.0067	U	0.0021	0.0067	mg/Kg
75-09-2	Methylene Chloride	0.013	U	0.0047	0.013	mg/Kg
156-60-5	trans-1,2-Dichloroethene	0.0067	U	0.0011	0.0067	mg/Kg
75-34-3	1,1-Dichloroethane	0.0067	U	0.0011	0.0067	mg/Kg
110-82-7	Cyclohexane	0.0067	U	0.0011	0.0067	mg/Kg
78-93-3	2-Butanone	0.033	U	0.0087	0.033	mg/Kg
56-23-5	Carbon Tetrachloride	0.0067	U	0.0013	0.0067	mg/Kg
156-59-2	cis-1,2-Dichloroethene	0.0067	U	0.0010	0.0067	mg/Kg
74-97-5	Bromochloromethane	0.0067	U	0.0015	0.0067	mg/Kg
67-66-3	Chloroform	0.0067	U	0.0011	0.0067	mg/Kg
71-55-6	1,1,1-Trichloroethane	0.0067	U	0.0012	0.0067	mg/Kg
108-87-2	Methylcyclohexane	0.0022	J	0.0012	0.0067	mg/Kg
71-43-2	Benzene	0.0067	U	0.0011	0.0067	mg/Kg
107-06-2	1,2-Dichloroethane	0.0067	U	0.0011	0.0067	mg/Kg
79-01-6	Trichloroethene	0.0067	U	0.0011	0.0067	mg/Kg
78-87-5	1,2-Dichloropropane	0.0067	U	0.0012	0.0067	mg/Kg
75-27-4	Bromodichloromethane	0.0067	U	0.0010	0.0067	mg/Kg
108-10-1	4-Methyl-2-Pentanone	0.063		0.0048	0.033	mg/Kg
108-88-3	Toluene	0.015		0.0010	0.0067	mg/Kg

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75-65-0	Tert butyl alcohol	80.2	J		5.23	ug/Kg
000108-11-2	2-Pentanol, 4-methyl-	40.2	J		10.6	ug/Kg
103-65-1	n-propylbenzene	8.60	J		12.8	ug/Kg
000611-14-3	Benzene, 1-ethyl-2-methyl-	68.5	J		12.9	ug/Kg
108-67-8	1,3,5-Trimethylbenzene	24.3	J		12.9	ug/Kg
95-63-6	1,2,4-Trimethylbenzene	110	J		13.2	ug/Kg
000873-49-4	Benzene, cyclopropyl-	63.4	J		13.8	ug/Kg
104-51-8	n-Butylbenzene	4.90	J		13.8	ug/Kg
001758-88-9	Benzene, 2-ethyl-1,4-dimethyl-	48.3	J		14.1	ug/Kg
000488-23-3	Benzene, 1,2,3,4-tetramethyl-	37.5	J		14.4	ug/Kg
000934-74-7	Benzene, 1-ethyl-3,5-dimethyl-	45.1	J		14.5	ug/Kg
002039-89-6	Benzene, 2-ethenyl-1,4-dimethyl-	100	J		14.8	ug/Kg
017059-48-2	1H-Indene, 2,3-dihydro-1,6-dimethy	38.0	J		15.2	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