

Report of Analysis

Client:	PSEG	Date Collected:	08/06/25
Project:	PSEG Metro Way	Date Received:	08/06/25
Client Sample ID:	TP-4-VOC	SDG No.:	Q2779
Lab Sample ID:	Q2779-03	Matrix:	SOIL
Analytical Method:	8260D	% Solid:	82.4
Sample Wt/Vol:	4.27	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
VW032031.D	1	08/06/25 16:12	VW080625

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
75-71-8	Dichlorodifluoromethane	0.0071	U	0.0016	0.0071	mg/Kg
74-87-3	Chloromethane	0.0071	U	0.0016	0.0071	mg/Kg
75-01-4	Vinyl Chloride	0.0071	U	0.0011	0.0071	mg/Kg
74-83-9	Bromomethane	0.0071	U	0.0015	0.0071	mg/Kg
75-00-3	Chloroethane	0.0071	U	0.0018	0.0071	mg/Kg
75-69-4	Trichlorofluoromethane	0.0071	U	0.0017	0.0071	mg/Kg
76-13-1	1,1,2-Trichlorotrifluoroethane	0.0071	U	0.0015	0.0071	mg/Kg
75-35-4	1,1-Dichloroethene	0.0071	U	0.0014	0.0071	mg/Kg
67-64-1	Acetone	0.080		0.0067	0.036	mg/Kg
75-15-0	Carbon Disulfide	0.0019	J	0.0015	0.0071	mg/Kg
1634-04-4	Methyl tert-butyl Ether	0.0071	U	0.0010	0.0071	mg/Kg
79-20-9	Methyl Acetate	0.0071	U	0.0022	0.0071	mg/Kg
75-09-2	Methylene Chloride	0.014	U	0.0050	0.014	mg/Kg
156-60-5	trans-1,2-Dichloroethene	0.0071	U	0.0012	0.0071	mg/Kg
75-34-3	1,1-Dichloroethane	0.0071	U	0.0011	0.0071	mg/Kg
110-82-7	Cyclohexane	0.0071	U	0.0011	0.0071	mg/Kg
78-93-3	2-Butanone	0.024	J	0.0093	0.036	mg/Kg
56-23-5	Carbon Tetrachloride	0.0071	U	0.0014	0.0071	mg/Kg
156-59-2	cis-1,2-Dichloroethene	0.0071	U	0.0011	0.0071	mg/Kg
74-97-5	Bromochloromethane	0.0071	U	0.0016	0.0071	mg/Kg
67-66-3	Chloroform	0.0071	U	0.0012	0.0071	mg/Kg
71-55-6	1,1,1-Trichloroethane	0.0071	U	0.0013	0.0071	mg/Kg
108-87-2	Methylcyclohexane	0.0071	U	0.0013	0.0071	mg/Kg
71-43-2	Benzene	0.0071	U	0.0011	0.0071	mg/Kg
107-06-2	1,2-Dichloroethane	0.0071	U	0.0011	0.0071	mg/Kg
79-01-6	Trichloroethene	0.0071	U	0.0012	0.0071	mg/Kg
78-87-5	1,2-Dichloropropane	0.0071	U	0.0013	0.0071	mg/Kg
75-27-4	Bromodichloromethane	0.0071	U	0.0011	0.0071	mg/Kg
108-10-1	4-Methyl-2-Pentanone	0.036	U	0.0051	0.036	mg/Kg
108-88-3	Toluene	0.0071	U	0.0011	0.0071	mg/Kg

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013152-05-1	Cyclooctene, 3-methyl-	9.70	J		12.3	ug/Kg
004057-42-5	unknown12.775	9.00	J		12.8	ug/Kg
004291-79-6	Cyclohexane, 1-methyl-2-propyl-	10.4	J		12.9	ug/Kg
135-98-8	sec-Butylbenzene	1.80	J		13.4	ug/Kg
000493-02-7	Naphthalene, decahydro-, trans-	17.4	J		13.9	ug/Kg
000535-77-3	Benzene, 1-methyl-3-(1-methylethyl	24.4	J		14.1	ug/Kg
000934-10-1	3-Phenylbut-1-ene	12.4	J		14.2	ug/Kg
1000152-47-3	unknown14.379	9.00	J		14.4	ug/Kg
000527-53-7	Benzene, 1,2,3,5-tetramethyl-	18.9	J		14.4	ug/Kg
000095-93-2	Benzene, 1,2,4,5-tetramethyl-	26.3	J		14.8	ug/Kg
017059-48-2	1H-Indene, 2,3-dihydro-1,6-dimethy	9.00	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