

Report of Analysis

Client:	PSEG	Date Collected:	08/05/25
Project:	Bridgewater Switch	Date Received:	08/05/25
Client Sample ID:	R5	SDG No.:	Q2774
Lab Sample ID:	Q2774-05	Matrix:	SOIL
Analytical Method:	8260D	% Solid:	74.5
Sample Wt/Vol:	5.09	Units:	g
Soil Aliquot Vol:			uL
GC Column:	RXI-624	ID :	0.25
Prep Method :		Level :	LOW
		Final Vol:	5000
		Test:	VOC-TCLVOA-10

File ID/Qc Batch:	Dilution:	Date Analyzed	Prep Batch ID
VW032012.D	1	08/05/25 14:59	VW080525

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
75-71-8	Dichlorodifluoromethane	0.0066	U	0.0015	0.0066	mg/Kg
74-87-3	Chloromethane	0.0066	U	0.0015	0.0066	mg/Kg
75-01-4	Vinyl Chloride	0.0066	U	0.0010	0.0066	mg/Kg
74-83-9	Bromomethane	0.0066	U	0.0014	0.0066	mg/Kg
75-00-3	Chloroethane	0.0066	U	0.0017	0.0066	mg/Kg
75-69-4	Trichlorofluoromethane	0.0066	U	0.0016	0.0066	mg/Kg
76-13-1	1,1,2-Trichlorotrifluoroethane	0.0066	U	0.0014	0.0066	mg/Kg
75-35-4	1,1-Dichloroethene	0.0066	U	0.0013	0.0066	mg/Kg
67-64-1	Acetone	0.033	U	0.0062	0.033	mg/Kg
75-15-0	Carbon Disulfide	0.0066	U	0.0014	0.0066	mg/Kg
1634-04-4	Methyl tert-butyl Ether	0.0066	U	0.00096	0.0066	mg/Kg
79-20-9	Methyl Acetate	0.0066	U	0.0020	0.0066	mg/Kg
75-09-2	Methylene Chloride	0.013	U	0.0047	0.013	mg/Kg
156-60-5	trans-1,2-Dichloroethene	0.0066	U	0.0011	0.0066	mg/Kg
75-34-3	1,1-Dichloroethane	0.0066	U	0.0011	0.0066	mg/Kg
110-82-7	Cyclohexane	0.0066	U	0.0010	0.0066	mg/Kg
78-93-3	2-Butanone	0.033	U	0.0086	0.033	mg/Kg
56-23-5	Carbon Tetrachloride	0.0066	U	0.0013	0.0066	mg/Kg
156-59-2	cis-1,2-Dichloroethene	0.0066	U	0.00099	0.0066	mg/Kg
74-97-5	Bromochloromethane	0.0066	U	0.0015	0.0066	mg/Kg
67-66-3	Chloroform	0.0066	U	0.0011	0.0066	mg/Kg
71-55-6	1,1,1-Trichloroethane	0.0066	U	0.0012	0.0066	mg/Kg
108-87-2	Methylcyclohexane	0.0066	U	0.0012	0.0066	mg/Kg
71-43-2	Benzene	0.0066	U	0.0010	0.0066	mg/Kg
107-06-2	1,2-Dichloroethane	0.0066	U	0.0010	0.0066	mg/Kg
79-01-6	Trichloroethene	0.0066	U	0.0011	0.0066	mg/Kg
78-87-5	1,2-Dichloropropane	0.0066	U	0.0012	0.0066	mg/Kg
75-27-4	Bromodichloromethane	0.0066	U	0.0010	0.0066	mg/Kg
108-10-1	4-Methyl-2-Pentanone	0.033	U	0.0047	0.033	mg/Kg
108-88-3	Toluene	0.0066	U	0.0010	0.0066	mg/Kg

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000080-56-8	.alpha.-Pinene	28.0	J		12.5	ug/Kg
000079-92-5	Camphene	20.1	J		12.7	ug/Kg
018172-67-3	Bicyclo[3.1.1]heptane, 6,6-dimethy	45.6	J		13.0	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