

Report of Analysis

Client:	G Environmental		Date Collected:	05/27/25	
Project:	DPW		Date Received:	05/28/25	
Client Sample ID:	MW10		SDG No.:	Q2134	
Lab Sample ID:	Q2134-01		Matrix:	Water	
Analytical Method:	8260D		% Solid:	0	
Sample Wt/Vol:	5	Units: mL	Final Vol:	5000	uL
Soil Aliquot Vol:		uL	Test:	VOCMS Group2	
GC Column:	DB-624UI	ID : 0.18	Level :	LOW	
Prep Method :					

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VX046395.D	1		05/29/25 13:30	VX052925

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
75-71-8	Dichlorodifluoromethane	0.22	U	0.22	1.00	ug/L
74-87-3	Chloromethane	0.32	U	0.32	1.00	ug/L
75-01-4	Vinyl Chloride	0.26	U	0.26	1.00	ug/L
74-83-9	Bromomethane	1.40	U	1.40	5.00	ug/L
75-00-3	Chloroethane	0.47	U	0.47	1.00	ug/L
75-69-4	Trichlorofluoromethane	0.33	U	0.33	1.00	ug/L
76-13-1	1,1,2-Trichlorotrifluoroethane	0.25	U	0.25	1.00	ug/L
75-35-4	1,1-Dichloroethene	0.23	U	0.23	1.00	ug/L
67-64-1	Acetone	1.50	U	1.50	5.00	ug/L
75-15-0	Carbon Disulfide	0.21	U	0.21	1.00	ug/L
1634-04-4	Methyl tert-butyl Ether	0.16	U	0.16	1.00	ug/L
79-20-9	Methyl Acetate	0.27	UQ	0.27	1.00	ug/L
75-09-2	Methylene Chloride	0.28	U	0.28	1.00	ug/L
156-60-5	trans-1,2-Dichloroethene	0.23	U	0.23	1.00	ug/L
75-34-3	1,1-Dichloroethane	0.23	U	0.23	1.00	ug/L
110-82-7	Cyclohexane	40.8		1.50	5.00	ug/L
78-93-3	2-Butanone	0.98	U	0.98	5.00	ug/L
56-23-5	Carbon Tetrachloride	0.25	U	0.25	1.00	ug/L
156-59-2	cis-1,2-Dichloroethene	0.19	U	0.19	1.00	ug/L
74-97-5	Bromochloromethane	0.22	U	0.22	1.00	ug/L
67-66-3	Chloroform	0.25	U	0.25	1.00	ug/L
71-55-6	1,1,1-Trichloroethane	0.20	U	0.20	1.00	ug/L
108-87-2	Methylcyclohexane	45.8		0.16	1.00	ug/L
71-43-2	Benzene	2.50		0.15	1.00	ug/L
107-06-2	1,2-Dichloroethane	0.22	U	0.22	1.00	ug/L
79-01-6	Trichloroethene	0.090	U	0.090	1.00	ug/L
78-87-5	1,2-Dichloropropane	0.20	U	0.20	1.00	ug/L
75-27-4	Bromodichloromethane	0.22	U	0.22	1.00	ug/L
108-10-1	4-Methyl-2-Pentanone	0.68	U	0.68	5.00	ug/L
108-88-3	Toluene	1.40		0.14	1.00	ug/L

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CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
000107-83-5	Pentane, 2-methyl-	140	J		2.82	ug/L
000096-14-0	Pentane, 3-methyl-	130	J		3.10	ug/L
000096-37-7	Cyclopentane, methyl-	93.8	J		4.30	ug/L
000565-59-3	Pentane, 2,3-dimethyl-	110	J		5.62	ug/L
000594-82-1	Butane, 2,2,3,3-tetramethyl-	130	J		6.25	ug/L
010574-36-4	2-Hexene, 3-methyl-, (Z)-	71.6	J		6.84	ug/L
103-65-1	n-propylbenzene	91.1	J		11.3	ug/L
000611-14-3	Benzene, 1-ethyl-2-methyl-	80.3	J		11.4	ug/L
108-67-8	1,3,5-Trimethylbenzene	8.70	J		11.5	ug/L
000622-96-8	Benzene, 1-ethyl-4-methyl-	130	J		11.6	ug/L
98-06-6	tert-Butylbenzene	3.60	J		11.7	ug/L
95-63-6	1,2,4-Trimethylbenzene	170	J		11.8	ug/L
135-98-8	sec-Butylbenzene	11.0	J		11.9	ug/L
99-87-6	p-Isopropyltoluene	2.90	J		12.0	ug/L
000526-73-8	Benzene, 1,2,3-trimethyl-	72.0	J		12.1	ug/L
000622-97-9	Benzene, 1-ethenyl-4-methyl-	200	J		12.2	ug/L
104-51-8	n-Butylbenzene	11.9	J		12.3	ug/L
000934-80-5	Benzene, 4-ethyl-1,2-dimethyl-	110	J		12.6	ug/L
007525-62-4	Benzene, 1-ethenyl-3-ethyl-	150	J		12.7	ug/L
000527-53-7	Benzene, 1,2,3,5-tetramethyl-	76.9	J		12.9	ug/L
000824-22-6	1H-Indene, 2,3-dihydro-4-methyl-	84.7	J		13.2	ug/L
000824-90-8	1-Phenyl-1-butene	190	J		13.3	ug/L

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