

Report of Analysis

Client:	G Environmental		Date Collected:	08/26/25	
Project:	Lexington		Date Received:	08/26/25	
Client Sample ID:	MW1		SDG No.:	Q2963	
Lab Sample ID:	Q2963-01		Matrix:	Water	
Analytical Method:	8260D		% Solid:	0	
Sample Wt/Vol:	5	Units: mL	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
VN087673.D	1	08/26/25 16:24	VN082625

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	4.70	J	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	U	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	13.7		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	7.50		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	6.50		0.16	1.00	ug/L
71-43-2	Benzene	19.6		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	1.20		0.14	1.00	ug/L

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[illegible]

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001191-96-4	Cyclopropane, ethyl-	9.60	J		5.35	ug/L
003769-23-1	1-Hexene, 4-methyl-	6.30	J		5.80	ug/L
000096-37-7	Cyclopentane, methyl-	14.3	J		7.31	ug/L
000592-46-1	2,4-Hexadiene	7.70	J		8.71	ug/L
000591-49-1	Cyclohexene, 1-methyl-	6.30	J		10.4	ug/L
	unknown10.724	5.30	J		10.7	ug/L
000764-13-6	2,4-Hexadiene, 2,5-dimethyl-	8.80	J		11.0	ug/L
103-65-1	n-propylbenzene	7.50	J		13.0	ug/L
108-67-8	1,3,5-Trimethylbenzene	0.50	J		13.2	ug/L
95-63-6	1,2,4-Trimethylbenzene	4.10	J		13.5	ug/L
135-98-8	sec-Butylbenzene	1.10	J		13.6	ug/L
99-87-6	p-Isopropyltoluene	0.62	J		13.7	ug/L
026146-77-0	trans-Cinnamyl bromide	21.1	J		14.0	ug/L
001587-04-8	Benzene, 1-methyl-2-(2-propenyl)-	18.1	J		14.4	ug/L
000934-74-7	Benzene, 1-ethyl-3,5-dimethyl-	6.10	J		14.6	ug/L
002039-89-6	Benzene, 2-ethenyl-1,4-dimethyl-	7.60	J		14.9	ug/L
002234-20-0	2,4-Dimethylstyrene	6.30	J		15.0	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