

Report of Analysis

Client:	G Environmental		Date Collected:	07/02/25	
Project:	Capra		Date Received:	07/03/25	
Client Sample ID:	GCAP3W		SDG No.:	Q2503	
Lab Sample ID:	Q2503-02		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:	DB-624UI	ID : 0.18	Level :	LOW	
Prep Method :					

File ID/Qc Batch:	Dilution:	Date Analyzed	Prep Batch ID
VX046887.D	1	07/03/25 14:58	VX070325

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
75-71-8	Dichlorodifluoromethane	0.22	U	0.22	5.00	ug/L
74-87-3	Chloromethane	0.50	J	0.32	5.00	ug/L
75-01-4	Vinyl Chloride	0.26	U	0.26	5.00	ug/L
74-83-9	Bromomethane	1.40	U	1.40	5.00	ug/L
75-00-3	Chloroethane	0.47	U	0.47	5.00	ug/L
75-69-4	Trichlorofluoromethane	0.33	U	0.33	5.00	ug/L
76-13-1	1,1,2-Trichlorotrifluoroethane	0.25	U	0.25	5.00	ug/L
75-35-4	1,1-Dichloroethene	0.23	U	0.23	5.00	ug/L
67-64-1	Acetone	1.50	U	1.50	25.0	ug/L
75-15-0	Carbon Disulfide	0.21	U	0.21	5.00	ug/L
1634-04-4	Methyl tert-butyl Ether	0.16	U	0.16	5.00	ug/L
79-20-9	Methyl Acetate	0.27	U	0.27	5.00	ug/L
75-09-2	Methylene Chloride	0.28	U	0.28	5.00	ug/L
156-60-5	trans-1,2-Dichloroethene	0.23	U	0.23	5.00	ug/L
75-34-3	1,1-Dichloroethane	0.23	U	0.23	5.00	ug/L
110-82-7	Cyclohexane	14.1		1.50	5.00	ug/L
78-93-3	2-Butanone	0.98	U	0.98	25.0	ug/L
56-23-5	Carbon Tetrachloride	0.25	U	0.25	5.00	ug/L
156-59-2	cis-1,2-Dichloroethene	0.19	U	0.19	5.00	ug/L
74-97-5	Bromochloromethane	0.22	U	0.22	5.00	ug/L
67-66-3	Chloroform	0.25	U	0.25	5.00	ug/L
71-55-6	1,1,1-Trichloroethane	0.20	U	0.20	5.00	ug/L
108-87-2	Methylcyclohexane	48.9		0.16	5.00	ug/L
71-43-2	Benzene	0.15	U	0.15	5.00	ug/L
107-06-2	1,2-Dichloroethane	0.22	U	0.22	5.00	ug/L
79-01-6	Trichloroethene	0.090	U	0.090	5.00	ug/L
78-87-5	1,2-Dichloropropane	0.20	U	0.20	5.00	ug/L
75-27-4	Bromodichloromethane	0.22	U	0.22	5.00	ug/L
108-10-1	4-Methyl-2-Pentanone	0.68	U	0.68	25.0	ug/L
108-88-3	Toluene	0.38	J	0.14	5.00	ug/L

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CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
000079-29-8	Butane, 2,3-dimethyl-	8.10	J		2.80	ug/L
000107-83-5	Pentane, 2-methyl-	29.4	J		2.84	ug/L
000096-14-0	Pentane, 3-methyl-	25.4	J		3.12	ug/L
000096-37-7	Cyclopentane, methyl-	45.1	J		4.32	ug/L
000589-34-4	Hexane, 3-methyl-	33.9	J		5.84	ug/L
001759-58-6	Cyclopentane, 1,3-dimethyl-, trans	40.0	J		6.17	ug/L
002453-00-1	Cyclopentane, 1,3-dimethyl-	20.7	J		6.27	ug/L
000872-56-0	Isopropylcyclobutane	26.0	J		6.37	ug/L
001640-89-7	Cyclopentane, ethyl-	9.00	J		7.66	ug/L
016883-48-0	Cyclopentane, 1,2,4-trimethyl-, (1	6.20	J		7.78	ug/L
015890-40-1	Cyclopentane, 1,2,3-trimethyl-, (1	10.6	J		7.96	ug/L
000583-57-3	Cyclohexane, 1,2-dimethyl- (cis/tr	9.00	J		8.98	ug/L
002207-03-6	Cyclohexane, 1,3-dimethyl-, trans-	6.20	J		9.10	ug/L
001678-91-7	Cyclohexane, ethyl-	6.50	J		9.56	ug/L
103-65-1	n-propylbenzene	2.00	J		11.3	ug/L
135-98-8	sec-Butylbenzene	0.61	J		11.9	ug/L
104-51-8	n-Butylbenzene	0.95	J		12.3	ug/L
001758-88-9	Benzene, 2-ethyl-1,4-dimethyl-	6.40	J		12.6	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