

Report of Analysis

Client:	G Environmental		Date Collected:	02/27/25	
Project:	Fisal		Date Received:	02/27/25	
Client Sample ID:	T10		SDG No.:	Q1463	
Lab Sample ID:	Q1463-10		Matrix:	SOIL	
Analytical Method:	SW8260		% Solid:	92.8	
Sample Wt/Vol:	7.05	Units: g	Final Vol:	5000	uL
Soil Aliquot Vol:		uL	Test:	VOCMS Group2	
GC Column:	RXI-624	ID : 0.25	Level :	LOW	
Prep Method :					

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VY021458.D	1		03/07/25 14:05	VY030725

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
74-87-3	Chloromethane	0.89	U	0.89	3.80	ug/Kg
75-01-4	Vinyl Chloride	0.59	U	0.59	3.80	ug/Kg
74-83-9	Bromomethane	0.79	U	0.79	3.80	ug/Kg
75-00-3	Chloroethane	0.77	U	0.77	3.80	ug/Kg
75-65-0	Tert butyl alcohol	11.9	U	11.9	19.1	ug/Kg
75-35-4	1,1-Dichloroethene	0.60	U	0.60	3.80	ug/Kg
67-64-1	Acetone	4.80	U	4.80	19.1	ug/Kg
75-15-0	Carbon Disulfide	0.98	U	0.98	3.80	ug/Kg
1634-04-4	Methyl tert-butyl Ether	0.51	U	0.51	3.80	ug/Kg
75-09-2	Methylene Chloride	2.60	U	2.60	7.60	ug/Kg
156-60-5	trans-1,2-Dichloroethene	0.64	U	0.64	3.80	ug/Kg
75-34-3	1,1-Dichloroethane	0.48	U	0.48	3.80	ug/Kg
78-93-3	2-Butanone	4.30	U	4.30	19.1	ug/Kg
56-23-5	Carbon Tetrachloride	0.66	U	0.66	3.80	ug/Kg
156-59-2	cis-1,2-Dichloroethene	0.47	U	0.47	3.80	ug/Kg
67-66-3	Chloroform	0.51	U	0.51	3.80	ug/Kg
71-55-6	1,1,1-Trichloroethane	0.60	U	0.60	3.80	ug/Kg
71-43-2	Benzene	0.55	U	0.55	3.80	ug/Kg
107-06-2	1,2-Dichloroethane	0.47	U	0.47	3.80	ug/Kg
79-01-6	Trichloroethene	0.57	U	0.57	3.80	ug/Kg
78-87-5	1,2-Dichloropropane	0.50	U	0.50	3.80	ug/Kg
75-27-4	Bromodichloromethane	0.43	U	0.43	3.80	ug/Kg
108-10-1	4-Methyl-2-Pentanone	3.30	U	3.30	19.1	ug/Kg
108-88-3	Toluene	0.51	U	0.51	3.80	ug/Kg
10061-02-6	t-1,3-Dichloropropene	0.46	U	0.46	3.80	ug/Kg
10061-01-5	cis-1,3-Dichloropropene	0.44	U	0.44	3.80	ug/Kg
79-00-5	1,1,2-Trichloroethane	0.64	U	0.64	3.80	ug/Kg
591-78-6	2-Hexanone	3.70	U	3.70	19.1	ug/Kg
124-48-1	Dibromochloromethane	0.50	U	0.50	3.80	ug/Kg
127-18-4	Tetrachloroethene	0.68	U	0.68	3.80	ug/Kg

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CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
108-90-7	Chlorobenzene	0.57	U	0.57	3.80	ug/Kg
100-41-4	Ethyl Benzene	0.47	U	0.47	3.80	ug/Kg
179601-23-1	m/p-Xylenes	1.00	U	1.00	7.60	ug/Kg
95-47-6	o-Xylene	0.53	U	0.53	3.80	ug/Kg
100-42-5	Styrene	0.46	U	0.46	3.80	ug/Kg
75-25-2	Bromoform	0.62	U	0.62	3.80	ug/Kg
79-34-5	1,1,2,2-Tetrachloroethane	0.84	U	0.84	3.80	ug/Kg
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	55.5		70 (63) - 130 (155)	111%	SPK: 50
1868-53-7	Dibromofluoromethane	50.7		70 (70) - 130 (134)	101%	SPK: 50
2037-26-5	Toluene-d8	48.1		70 (74) - 130 (123)	96%	SPK: 50
460-00-4	4-Bromofluorobenzene	43.1		70 (38) - 130 (136)	86%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	290000	7.707			
540-36-3	1,4-Difluorobenzene	520000	8.616			
3114-55-4	Chlorobenzene-d5	457000	11.414			
3855-82-1	1,4-Dichlorobenzene-d4	176000	13.346			

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