

Report of Analysis

Client: Ambric Technology Corporation
Project: 6506 Haverford Road
Client Sample ID: VW1119SBL01
Lab Sample ID: VW1119SBL01
Analytical Method: 8260D
Sample Wt/Vol: 5 g

Level: LOW
Final Vol: 5000 uL

Date Collected:
Date Received:
SDG No.: Q3619
Matrix: SOIL
% Solid: 100
Test: VOCMS Group10

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
1634-04-4	Methyl tert-butyl Ether	0.00073	U	1	0.00073	0.0050	mg/Kg	11/19/25 10:36	VW111925
71-43-2	Benzene	0.00079	U	1	0.00079	0.0050	mg/Kg	11/19/25 10:36	VW111925
108-88-3	Toluene	0.00078	U	1	0.00078	0.0050	mg/Kg	11/19/25 10:36	VW111925
106-93-4	1,2-Dibromoethane	0.00088	U	1	0.00088	0.0050	mg/Kg	11/19/25 10:36	VW111925
100-41-4	Ethyl Benzene	0.00067	U	1	0.00067	0.0050	mg/Kg	11/19/25 10:36	VW111925
179601-23-1	m/p-Xylenes	0.0012	U	1	0.0012	0.010	mg/Kg	11/19/25 10:36	VW111925
1330-20-7	Total Xylenes	0.0020	U	1	0.0020	0.015	mg/Kg	11/19/25 10:36	VW111925
95-47-6	o-Xylene	0.00082	U	1	0.00082	0.0050	mg/Kg	11/19/25 10:36	VW111925
98-82-8	Isopropylbenzene	0.00078	U	1	0.00078	0.0050	mg/Kg	11/19/25 10:36	VW111925
108-67-8	1,3,5-Trimethylbenzene	0.00082	U	1	0.00082	0.0050	mg/Kg	11/19/25 10:36	VW111925
95-63-6	1,2,4-Trimethylbenzene	0.00064	U	1	0.00064	0.0050	mg/Kg	11/19/25 10:36	VW111925
91-20-3	Naphthalene	0.0026	U	1	0.0026	0.0050	mg/Kg	11/19/25 10:36	VW111925
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	60.5			63 - 155	121%	SPK: 50		
1868-53-7	Dibromofluoromethane	51.6			70 - 134	103%	SPK: 50		
2037-26-5	Toluene-d8	45.5			74 - 123	91%	SPK: 50		
460-00-4	4-Bromofluorobenzene	54.4			52 - 138	109%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	218000							
540-36-3	1,4-Difluorobenzene	509000							
3114-55-4	Chlorobenzene-d5	542000							
3855-82-1	1,4-Dichlorobenzene-d4	261000							

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