

Report of Analysis

Client:	CHA Companies, Inc.		Date Collected:	
Project:	OMH Tank Pull-011 - 078673-01 C-03		Date Received:	
Client Sample ID:	VX1127MBS01		SDG No.:	P4960
Lab Sample ID:	VX1127MBS01		Matrix:	SOIL
Analytical Method:	SW8260		% Solid:	100
Sample Wt/Vol:	5	Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	100	uL	Test:	VOCMS Group1
GC Column:	DB-624UI	ID : 0.18	Level :	MED
Prep Method :				

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VX044027.D	1		11/27/24 10:01	VX112724

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
1634-04-4	Methyl tert-butyl Ether	2200		67.0	500	ug/Kg
71-43-2	Benzene	2100		72.0	500	ug/Kg
108-88-3	Toluene	2100		67.0	500	ug/Kg
100-41-4	Ethyl Benzene	2100		62.0	500	ug/Kg
179601-23-1	m/p-Xylenes	4100		140	1000	ug/Kg
95-47-6	o-Xylene	2100		70.0	500	ug/Kg
98-82-8	Isopropylbenzene	2200		67.0	500	ug/Kg
103-65-1	n-propylbenzene	2200		64.0	500	ug/Kg
108-67-8	1,3,5-Trimethylbenzene	2200		64.0	500	ug/Kg
98-06-6	tert-Butylbenzene	2200		67.0	500	ug/Kg
95-63-6	1,2,4-Trimethylbenzene	2200		140	500	ug/Kg
135-98-8	sec-Butylbenzene	2100		67.0	500	ug/Kg
99-87-6	p-Isopropyltoluene	2200		58.0	500	ug/Kg
104-51-8	n-Butylbenzene	2100		63.0	500	ug/Kg
91-20-3	Naphthalene	2000		150	500	ug/Kg
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	56.2		50 - 163	112%	SPK: 50
1868-53-7	Dibromofluoromethane	54.0		54 - 147	108%	SPK: 50
2037-26-5	Toluene-d8	52.4		58 - 134	105%	SPK: 50
460-00-4	4-Bromofluorobenzene	52.9		29 - 146	106%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	117000	5.544			
540-36-3	1,4-Difluorobenzene	213000	6.757			
3114-55-4	Chlorobenzene-d5	181000	10.049			
3855-82-1	1,4-Dichlorobenzene-d4	80900	12.018			

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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