

Report of Analysis

Client:	CDM Smith		Date Collected:	04/28/25	
Project:	Con Ed UTEN Mount Vernon, NY		Date Received:	04/29/25	
Client Sample ID:	SS-9		SDG No.:	Q1914	
Lab Sample ID:	Q1914-12		Matrix:	SOIL	
Analytical Method:	8260D		% Solid:	87.1	
Sample Wt/Vol:	7.92	Units: g	Final Vol:	5000	uL
Soil Aliquot Vol:		uL	Test:	VOCMS Group3	
GC Column:	RXI-624	ID : 0.25	Level :	LOW	
Prep Method :					

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VY022178.D	1		05/06/25 13:03	VY050625

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
71-43-2	Benzene	0.57	U	0.57	3.60	ug/Kg
108-88-3	Toluene	0.57	U	0.57	3.60	ug/Kg
100-41-4	Ethyl Benzene	0.49	U	0.49	3.60	ug/Kg
1330-20-7	Total Xylenes	1.49	U	1.49	10.8	ug/Kg
98-82-8	Isopropylbenzene	0.57	U	0.57	3.60	ug/Kg
103-65-1	n-propylbenzene	0.53	U	0.53	3.60	ug/Kg
108-67-8	1,3,5-Trimethylbenzene	0.59	U	0.59	3.60	ug/Kg
98-06-6	tert-Butylbenzene	0.49	U	0.49	3.60	ug/Kg
95-63-6	1,2,4-Trimethylbenzene	0.46	U	0.46	3.60	ug/Kg
135-98-8	sec-Butylbenzene	0.48	U	0.48	3.60	ug/Kg
99-87-6	p-Isopropyltoluene	0.45	U	0.45	3.60	ug/Kg
104-51-8	n-Butylbenzene	1.10	U	1.10	3.60	ug/Kg
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	56.4		63 - 155	113%	SPK: 50
1868-53-7	Dibromofluoromethane	52.0		70 - 134	104%	SPK: 50
2037-26-5	Toluene-d8	49.2		74 - 123	98%	SPK: 50
460-00-4	4-Bromofluorobenzene	38.8		38 - 136	78%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	197000	7.713			
540-36-3	1,4-Difluorobenzene	384000	8.616			
3114-55-4	Chlorobenzene-d5	340000	11.42			
3855-82-1	1,4-Dichlorobenzene-d4	121000	13.346			
TENTATIVE IDENTIFIED COMPOUNDS						
Total VOC TICs	Total VOC TICs	0	U		0	ug/Kg

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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