284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900, Fax : 908 789 8922

LAB CHRONICLE

OrderID:	Q2230	**OrderDate:**	6/4/2025 4:40:00 PM
Client:	CDM Smith	**Project:**	Con Ed UTEN Mount Vernon, NY
Contact:	Marcie Ann Encinas	**Location:**	N31,VOA Ref. #3 Water

LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
Q2230-01	**FB-060425**	**Water**			**06/04/25**			**06/04/25**
			VOCMS Group3	8260-Low			06/10/25	
Q2230-02	**GW-MW01-060425**	**Water**			**06/04/25**			**06/04/25**
			VOCMS Group3	8260-Low			06/10/25	
Q2230-05	**GW-MW901-060425**	**Water**			**06/04/25**			**06/04/25**
			VOCMS Group3	8260-Low			06/10/25	
Q2230-06	**TB060425**	**Water**			**06/04/25**			**06/04/25**
			VOCMS Group3	8260-Low			06/10/25	

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Hit Summary Sheet
SW-846

SDG No.: Q2230

Client: CDM Smith

Sample ID	Client ID	Matrix	Parameter	Concentration	C	MDL	RDL	Units
Client ID:	**GW-MW01-060425**							
Q2230-02	GW-MW01-060425	Water	Benzene	0.36	J	0.15	1.00	ug/L
Q2230-02	GW-MW01-060425	Water	Toluene	1.60		0.14	1.00	ug/L
Q2230-02	GW-MW01-060425	Water	Ethyl Benzene	3.80		0.13	1.00	ug/L
Q2230-02	GW-MW01-060425	Water	Total Xylenes	36.1		0.36	3.00	ug/L
Q2230-02	GW-MW01-060425	Water	Isopropylbenzene	1.30		0.12	1.00	ug/L
Q2230-02	GW-MW01-060425	Water	n-propylbenzene	2.50		0.13	1.00	ug/L
Q2230-02	GW-MW01-060425	Water	1,3,5-Trimethylbenzene	7.40		0.15	1.00	ug/L
Q2230-02	GW-MW01-060425	Water	1,2,4-Trimethylbenzene	26.0		0.14	1.00	ug/L
Q2230-02	GW-MW01-060425	Water	sec-Butylbenzene	1.00		0.13	1.00	ug/L
Q2230-02	GW-MW01-060425	Water	n-Butylbenzene	1.40		0.15	1.00	ug/L
			Total Voc :	81.5				
			Total Concentration:	81.5				
Client ID:	**GW-MW901-060425**							
Q2230-05	GW-MW901-06042	Water	Benzene	0.32	J	0.15	1.00	ug/L
Q2230-05	GW-MW901-06042	Water	Toluene	1.60		0.14	1.00	ug/L
Q2230-05	GW-MW901-06042	Water	Ethyl Benzene	3.70		0.13	1.00	ug/L
Q2230-05	GW-MW901-06042	Water	Total Xylenes	34.6		0.36	3.00	ug/L
Q2230-05	GW-MW901-06042	Water	Isopropylbenzene	1.20		0.12	1.00	ug/L
Q2230-05	GW-MW901-06042	Water	n-propylbenzene	2.30		0.13	1.00	ug/L
Q2230-05	GW-MW901-06042	Water	1,3,5-Trimethylbenzene	6.80		0.15	1.00	ug/L
Q2230-05	GW-MW901-06042	Water	1,2,4-Trimethylbenzene	24.8		0.14	1.00	ug/L
Q2230-05	GW-MW901-06042	Water	sec-Butylbenzene	0.90	J	0.13	1.00	ug/L
Q2230-05	GW-MW901-06042	Water	n-Butylbenzene	1.20		0.15	1.00	ug/L
			Total Voc :	77.4				
			Total Concentration:	77.4				

SAMPLE DATA

Alliance TECHNICAL GROUP
284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Report of Analysis

Client:	CDM Smith	Date Collected:	06/04/25
Project:	Con Ed UTEN Mount Vernon, NY	Date Received:	06/04/25
Client Sample ID:	FB-060425	SDG No.:	Q2230
Lab Sample ID:	Q2230-01	Matrix:	Water
Analytical Method:	8260D	% Solid:	0
Sample Wt/Vol:	5 Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOCMS Group3
GC Column:	DB-624UI ID : 0.18	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VX046594.D	1		06/10/25 12:09	VX061025

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
71-43-2	Benzene	1.00	U	0.15	1.00	ug/L
108-88-3	Toluene	1.00	U	0.14	1.00	ug/L
100-41-4	Ethyl Benzene	1.00	U	0.13	1.00	ug/L
1330-20-7	Total Xylenes	3.00	U	0.36	3.00	ug/L
98-82-8	Isopropylbenzene	1.00	U	0.12	1.00	ug/L
103-65-1	n-propylbenzene	1.00	U	0.13	1.00	ug/L
108-67-8	1,3,5-Trimethylbenzene	1.00	U	0.15	1.00	ug/L
98-06-6	tert-Butylbenzene	1.00	U	0.14	1.00	ug/L
95-63-6	1,2,4-Trimethylbenzene	1.00	U	0.14	1.00	ug/L
135-98-8	sec-Butylbenzene	1.00	U	0.13	1.00	ug/L
99-87-6	p-Isopropyltoluene	1.00	U	0.13	1.00	ug/L
104-51-8	n-Butylbenzene	1.00	U	0.15	1.00	ug/L
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	49.6		74 - 125	99%	SPK: 50
1868-53-7	Dibromofluoromethane	50.3		75 - 124	101%	SPK: 50
2037-26-5	Toluene-d8	54.7		86 - 113	109%	SPK: 50
460-00-4	4-Bromofluorobenzene	54.0		77 - 121	108%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	110000	5.562			
540-36-3	1,4-Difluorobenzene	215000	6.769			
3114-55-4	Chlorobenzene-d5	221000	10.055			
3855-82-1	1,4-Dichlorobenzene-d4	108000	12.018			

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

Alliance TECHNICAL GROUP
284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Report of Analysis

Client:	CDM Smith	Date Collected:	06/04/25
Project:	Con Ed UTEN Mount Vernon, NY	Date Received:	06/04/25
Client Sample ID:	GW-MW01-060425	SDG No.:	Q2230
Lab Sample ID:	Q2230-02	Matrix:	Water
Analytical Method:	8260D	% Solid:	0
Sample Wt/Vol:	5 Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOCMS Group3
GC Column:	DB-624UI ID : 0.18	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VX046602.D	1		06/10/25 14:59	VX061025

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
71-43-2	Benzene	0.36	J	0.15	1.00	ug/L
108-88-3	Toluene	1.60		0.14	1.00	ug/L
100-41-4	Ethyl Benzene	3.80		0.13	1.00	ug/L
1330-20-7	Total Xylenes	36.1		0.36	3.00	ug/L
98-82-8	Isopropylbenzene	1.30		0.12	1.00	ug/L
103-65-1	n-propylbenzene	2.50		0.13	1.00	ug/L
108-67-8	1,3,5-Trimethylbenzene	7.40		0.15	1.00	ug/L
98-06-6	tert-Butylbenzene	1.00	U	0.14	1.00	ug/L
95-63-6	1,2,4-Trimethylbenzene	26.0		0.14	1.00	ug/L
135-98-8	sec-Butylbenzene	1.00		0.13	1.00	ug/L
99-87-6	p-Isopropyltoluene	1.00	U	0.13	1.00	ug/L
104-51-8	n-Butylbenzene	1.40		0.15	1.00	ug/L
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	49.8		74 - 125	100%	SPK: 50
1868-53-7	Dibromofluoromethane	49.9		75 - 124	100%	SPK: 50
2037-26-5	Toluene-d8	54.2		86 - 113	108%	SPK: 50
460-00-4	4-Bromofluorobenzene	56.1		77 - 121	112%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	113000	5.562			
540-36-3	1,4-Difluorobenzene	225000	6.769			
3114-55-4	Chlorobenzene-d5	232000	10.055			
3855-82-1	1,4-Dichlorobenzene-d4	118000	12.018			

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

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Report of Analysis

Client:	CDM Smith	Date Collected:	06/04/25
Project:	Con Ed UTEN Mount Vernon, NY	Date Received:	06/04/25
Client Sample ID:	GW-MW901-060425	SDG No.:	Q2230
Lab Sample ID:	Q2230-05	Matrix:	Water
Analytical Method:	8260D	% Solid:	0
Sample Wt/Vol:	5 Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOCMS Group3
GC Column:	DB-624UI ID : 0.18	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VX046601.D	1		06/10/25 14:38	VX061025

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
71-43-2	Benzene	0.32	J	0.15	1.00	ug/L
108-88-3	Toluene	1.60		0.14	1.00	ug/L
100-41-4	Ethyl Benzene	3.70		0.13	1.00	ug/L
1330-20-7	Total Xylenes	34.6		0.36	3.00	ug/L
98-82-8	Isopropylbenzene	1.20		0.12	1.00	ug/L
103-65-1	n-propylbenzene	2.30		0.13	1.00	ug/L
108-67-8	1,3,5-Trimethylbenzene	6.80		0.15	1.00	ug/L
98-06-6	tert-Butylbenzene	1.00	U	0.14	1.00	ug/L
95-63-6	1,2,4-Trimethylbenzene	24.8		0.14	1.00	ug/L
135-98-8	sec-Butylbenzene	0.90	J	0.13	1.00	ug/L
99-87-6	p-Isopropyltoluene	1.00	U	0.13	1.00	ug/L
104-51-8	n-Butylbenzene	1.20		0.15	1.00	ug/L
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	50.2		74 - 125	100%	SPK: 50
1868-53-7	Dibromofluoromethane	49.9		75 - 124	100%	SPK: 50
2037-26-5	Toluene-d8	54.3		86 - 113	109%	SPK: 50
460-00-4	4-Bromofluorobenzene	55.6		77 - 121	111%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	112000	5.562			
540-36-3	1,4-Difluorobenzene	225000	6.769			
3114-55-4	Chlorobenzene-d5	232000	10.055			
3855-82-1	1,4-Dichlorobenzene-d4	117000	12.018			

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

Alliance TECHNICAL GROUP
284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Report of Analysis

Client:	CDM Smith	Date Collected:	06/04/25
Project:	Con Ed UTEN Mount Vernon, NY	Date Received:	06/04/25
Client Sample ID:	TB060425	SDG No.:	Q2230
Lab Sample ID:	Q2230-06	Matrix:	Water
Analytical Method:	8260D	% Solid:	0
Sample Wt/Vol:	5 Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOCMS Group3
GC Column:	DB-624UI ID : 0.18	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VX046595.D	1		06/10/25 12:30	VX061025

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
71-43-2	Benzene	1.00	U	0.15	1.00	ug/L
108-88-3	Toluene	1.00	U	0.14	1.00	ug/L
100-41-4	Ethyl Benzene	1.00	U	0.13	1.00	ug/L
1330-20-7	Total Xylenes	3.00	U	0.36	3.00	ug/L
98-82-8	Isopropylbenzene	1.00	U	0.12	1.00	ug/L
103-65-1	n-propylbenzene	1.00	U	0.13	1.00	ug/L
108-67-8	1,3,5-Trimethylbenzene	1.00	U	0.15	1.00	ug/L
98-06-6	tert-Butylbenzene	1.00	U	0.14	1.00	ug/L
95-63-6	1,2,4-Trimethylbenzene	1.00	U	0.14	1.00	ug/L
135-98-8	sec-Butylbenzene	1.00	U	0.13	1.00	ug/L
99-87-6	p-Isopropyltoluene	1.00	U	0.13	1.00	ug/L
104-51-8	n-Butylbenzene	1.00	U	0.15	1.00	ug/L
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	50.3		74 - 125	101%	SPK: 50
1868-53-7	Dibromofluoromethane	50.7		75 - 124	101%	SPK: 50
2037-26-5	Toluene-d8	54.5		86 - 113	109%	SPK: 50
460-00-4	4-Bromofluorobenzene	55.3		77 - 121	111%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	115000	5.568			
540-36-3	1,4-Difluorobenzene	224000	6.769			
3114-55-4	Chlorobenzene-d5	234000	10.055			
3855-82-1	1,4-Dichlorobenzene-d4	116000	12.018			

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

QC SUMMARY

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Surrogate Summary

SDG No.: Q2230

Client: CDM Smith

Analytical Method: SW8260-Low

Lab Sample ID	Client ID	Parameter	Spike	Result	RecoveryQual	Limits Low	Limits High
Q2230-01	FB-060425	1,2-Dichloroethane-d4	50	49.6	99	74	125
		Dibromofluoromethane	50	50.3	101	75	124
		Toluene-d8	50	54.7	109	86	113
		4-Bromofluorobenzene	50	54.0	108	77	121
Q2230-02	GW-MW01-060425	1,2-Dichloroethane-d4	50	49.8	100	74	125
		Dibromofluoromethane	50	49.9	100	75	124
		Toluene-d8	50	54.2	108	86	113
		4-Bromofluorobenzene	50	56.1	112	77	121
Q2230-03MS	GW-MW01-060425MS	1,2-Dichloroethane-d4	50	47.5	95	74	125
		Dibromofluoromethane	50	52.2	104	75	124
		Toluene-d8	50	51.1	102	86	113
		4-Bromofluorobenzene	50	50.9	102	77	121
Q2230-04MSD	GW-MW01-060425MSD	1,2-Dichloroethane-d4	50	45.9	92	74	125
		Dibromofluoromethane	50	50.4	101	75	124
		Toluene-d8	50	50.2	100	86	113
		4-Bromofluorobenzene	50	51.1	102	77	121
Q2230-05	GW-MW901-060425	1,2-Dichloroethane-d4	50	50.2	100	74	125
		Dibromofluoromethane	50	49.9	100	75	124
		Toluene-d8	50	54.3	109	86	113
		4-Bromofluorobenzene	50	55.6	111	77	121
Q2230-06	TB060425	1,2-Dichloroethane-d4	50	50.3	101	74	125
		Dibromofluoromethane	50	50.6	101	75	124
		Toluene-d8	50	54.5	109	86	113
		4-Bromofluorobenzene	50	55.3	111	77	121
VX0610WBL01	VX0610WBL01	1,2-Dichloroethane-d4	50	50.1	100	74	125
		Dibromofluoromethane	50	50.1	100	75	124
		Toluene-d8	50	54.3	109	86	113
		4-Bromofluorobenzene	50	54.8	110	77	121
VX0610WBS01	VX0610WBS01	1,2-Dichloroethane-d4	50	47.4	95	74	125
		Dibromofluoromethane	50	53.0	106	75	124
		Toluene-d8	50	51.5	103	86	113
		4-Bromofluorobenzene	50	53.1	106	77	121
VX0611WBL02	VX0611WBL02	1,2-Dichloroethane-d4	50	49.7	99	74	125
		Dibromofluoromethane	50	50.4	101	75	124
		Toluene-d8	50	54.3	109	86	113
		4-Bromofluorobenzene	50	55.3	111	77	121
VX0611WBS02	VX0611WBS02	1,2-Dichloroethane-d4	50	44.0	88	74	125
		Dibromofluoromethane	50	47.8	96	75	124
		Toluene-d8	50	47.2	94	86	113
		4-Bromofluorobenzene	50	48.5	97	77	121

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Matrix Spike/Matrix Spike Duplicate Summary

SW-846

SDG No.: Q2230

Client: CDM Smith

Analytical Method: SW8260-Low

Parameter	Spike	Sample Result	Result	Units	Rec	Rec Qual	RPD	RPD Qual	Limits Low	Limits High	Limits RPD
Lab Sample ID :	**Q2230-03MS**	**Client Sample ID :**	**GW-MW01-060425MS**					**Datafile :**	VX046641.D		
Benzene	50	0.36	58.1	ug/L	115				81	128	
Toluene	50	1.60	62.2	ug/L	121	*			85	115	
Ethyl Benzene	50	3.80	65.8	ug/L	124				81	128	
m/p-Xylenes	100	16.1	150	ug/L	134	*			69	129	
o-Xylene	50	20.0	94.6	ug/L	149	*			75	127	
Isopropylbenzene	50	1.30	62.9	ug/L	123	*			76	121	
N-propylbenzene	50	2.50	63.6	ug/L	122	*			79	116	
1,3,5-Trimethylbenzene	50	7.40	72.5	ug/L	130				40	169	
tert-Butylbenzene	50	0	61.8	ug/L	124	*			80	116	
1,2,4-Trimethylbenzene	50	26.0	100	ug/L	148	*			81	116	
Sec-butylbenzene	50	1.00	60.8	ug/L	120				70	122	
p-Isopropyltoluene	50	0	60.8	ug/L	122				64	124	
n-Butylbenzene	50	1.40	60.9	ug/L	119				53	140	

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Matrix Spike/Matrix Spike Duplicate Summary

SW-846

SDG No.: Q2230

Client: CDM Smith

Analytical Method: SW8260-Low

Parameter	Spike	Sample Result	Result	Units	Rec	Rec Qual	RPD	RPD Qual	Limits Low	Limits High	Limits RPD
Lab Sample ID :	**Q2230-04MSD**	**Client Sample ID :**	**GW-MW01-060425MSD**					**Datafile :**	VX046642.D		
Benzene	50	0.36	50.7	ug/L	101		13		81	128	20
Toluene	50	1.60	54.6	ug/L	106		13		85	115	20
Ethyl Benzene	50	3.80	58.5	ug/L	109		13		81	128	20
m/p-Xylenes	100	16.1	140	ug/L	124		8		69	129	20
o-Xylene	50	20.0	95.5	ug/L	151	*	1		75	127	20
Isopropylbenzene	50	1.30	53.5	ug/L	104		17		76	121	20
N-propylbenzene	50	2.50	55.2	ug/L	105		15		79	116	20
1,3,5-Trimethylbenzene	50	7.40	66.8	ug/L	119		9		40	169	20
tert-Butylbenzene	50	0	52.1	ug/L	104		17		80	116	20
1,2,4-Trimethylbenzene	50	26.0	110	ug/L	168	*	13		81	116	20
Sec-butylbenzene	50	1.00	52.0	ug/L	102		16		70	122	20
p-Isopropyltoluene	50	0	51.8	ug/L	104		16		64	124	20
n-Butylbenzene	50	1.40	53.5	ug/L	104		13		53	140	20

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q2230

Client: CDM Smith

Analytical Method: SW8260-Low **Datafile :** VX046589.D

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	Limits Low	Limits High	RPD
VX0610WBS01	Benzene	20	21.0	ug/L	105			82	109	
	Toluene	20	21.2	ug/L	106			82	110	
	Ethyl Benzene	20	20.3	ug/L	102			83	109	
	m/p-Xylenes	40	41.6	ug/L	104			82	110	
	o-Xylene	20	21.1	ug/L	106			83	109	
	Isopropylbenzene	20	20.2	ug/L	101			83	112	
	N-propylbenzene	20	20.0	ug/L	100			83	112	
	1,3,5-Trimethylbenzene	20	20.5	ug/L	103			85	112	
	tert-Butylbenzene	20	20.4	ug/L	102			83	112	
	1,2,4-Trimethylbenzene	20	20.8	ug/L	104			85	111	
	Sec-butylbenzene	20	20.5	ug/L	103			81	114	
	p-Isopropyltoluene	20	20.2	ug/L	101			78	116	
	n-Butylbenzene	20	19.9	ug/L	100			75	115	

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q2230

Client: CDM Smith

Analytical Method: SW8260-Low **Datafile :** VX046639.D

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	Limits Low	Limits High	RPD
VX0611WBS02	Benzene	20	19.6	ug/L	98			82	109	
	Toluene	20	20.2	ug/L	101			82	110	
	Ethyl Benzene	20	19.2	ug/L	96			83	109	
	m/p-Xylenes	40	40.1	ug/L	100			82	110	
	o-Xylene	20	20.1	ug/L	101			83	109	
	Isopropylbenzene	20	19.1	ug/L	96			83	112	
	N-propylbenzene	20	18.8	ug/L	94			83	112	
	1,3,5-Trimethylbenzene	20	18.9	ug/L	95			85	112	
	tert-Butylbenzene	20	19.2	ug/L	96			83	112	
	1,2,4-Trimethylbenzene	20	19.2	ug/L	96			85	111	
	Sec-butylbenzene	20	19.3	ug/L	97			81	114	
	p-Isopropyltoluene	20	18.9	ug/L	95			78	116	
	n-Butylbenzene	20	18.3	ug/L	92			75	115	

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

VOLATILE METHOD BLANK SUMMARY

EPA SAMPLE NO.

VX0610WBL01

Lab Name: CHEMTECH Contract: CAMP02

Lab Code: CHEM Case No.: Q2230 SAS No.: Q2230 SDG NO.: Q2230

Lab File ID: VX046588.D Lab Sample ID: VX0610WBL01

Date Analyzed: 06/10/2025 Time Analyzed: 09:57

GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: (Y/N) N

Instrument ID: MSVOA_X

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
VX0610WBS01	VX0610WBS01	VX046589.D	06/10/2025
FB-060425	Q2230-01	VX046594.D	06/10/2025
TB060425	Q2230-06	VX046595.D	06/10/2025
GW-MW901-060425	Q2230-05	VX046601.D	06/10/2025
GW-MW01-060425	Q2230-02	VX046602.D	06/10/2025

COMMENTS:

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

VOLATILE METHOD BLANK SUMMARY

EPA SAMPLE NO.

VX0611WBL02

Lab Name: CHEMTECH Contract: CAMP02

Lab Code: CHEM Case No.: Q2230 SAS No.: Q2230 SDG NO.: Q2230

Lab File ID: VX046637.D Lab Sample ID: VX0611WBL02

Date Analyzed: 06/11/2025 Time Analyzed: 15:14

GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: (Y/N) N

Instrument ID: MSVOA_X

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
VX0611WBS02	VX0611WBS02	VX046639.D	06/11/2025
GW-MW01-060425MS	Q2230-03MS	VX046641.D	06/11/2025
GW-MW01-060425MSD	Q2230-04MSD	VX046642.D	06/11/2025

COMMENTS:

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: CHEMTECH Contract: CAMP02

Lab Code: CHEM Case No.: Q2230 SAS No.: Q2230 SDG NO.: Q2230

Lab File ID: VX046516.D BFB Injection Date: 06/06/2025

Instrument ID: MSVOA_X BFB Injection Time: 08:47

GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: Y/N N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	21
75	30.0 - 60.0% of mass 95	56.5
95	Base Peak, 100% relative abundance	100
96	5.0 - 9.0% of mass 95	6.5
173	Less than 2.0% of mass 174	0.6 (1) 1
174	50.0 - 100.0% of mass 95	65.2
175	5.0 - 9.0% of mass 174	5 (7.7) 1
176	95.0 - 101.0% of mass 174	62.7 (96.1) 1
177	5.0 - 9.0% of mass 176	4.2 (6.6) 2

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
VSTDICC005	VSTDICC005	VX046518.D	06/06/2025	09:42
VSTDICC020	VSTDICC020	VX046519.D	06/06/2025	10:18
VSTDICCC050	VSTDICCC050	VX046520.D	06/06/2025	10:40
VSTDICC100	VSTDICC100	VX046521.D	06/06/2025	11:02
VSTDICC150	VSTDICC150	VX046522.D	06/06/2025	11:25
VSTDICC001	VSTDICC001	VX046524.D	06/06/2025	12:57

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: CHEMTECH Contract: CAMP02

Lab Code: CHEM Case No.: Q2230 SAS No.: Q2230 SDG NO.: Q2230

Lab File ID: VX046585.D BFB Injection Date: 06/10/2025

Instrument ID: MSVOA_X BFB Injection Time: 08:36

GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: Y/N N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	20.5
75	30.0 - 60.0% of mass 95	55
95	Base Peak, 100% relative abundance	100
96	5.0 - 9.0% of mass 95	6.7
173	Less than 2.0% of mass 174	0.3 (0.4) 1
174	50.0 - 100.0% of mass 95	70.1
175	5.0 - 9.0% of mass 174	5.3 (7.6) 1
176	95.0 - 101.0% of mass 174	68.4 (97.6) 1
177	5.0 - 9.0% of mass 176	4.6 (6.8) 2

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
VSTDCCC050	VSTDCCC050	VX046586.D	06/10/2025	09:07
VX0610WBL01	VX0610WBL01	VX046588.D	06/10/2025	09:57
VX0610WBS01	VX0610WBS01	VX046589.D	06/10/2025	10:18
FB-060425	Q2230-01	VX046594.D	06/10/2025	12:09
TB060425	Q2230-06	VX046595.D	06/10/2025	12:30
GW-MW901-060425	Q2230-05	VX046601.D	06/10/2025	14:38
GW-MW01-060425	Q2230-02	VX046602.D	06/10/2025	14:59

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: CHEMTECH Contract: CAMP02

Lab Code: CHEM Case No.: Q2230 SAS No.: Q2230 SDG NO.: Q2230

Lab File ID: VX046635.D BFB Injection Date: 06/11/2025

Instrument ID: MSVOA_X BFB Injection Time: 14:01

GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: Y/N N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	21
75	30.0 - 60.0% of mass 95	52.7
95	Base Peak, 100% relative abundance	100
96	5.0 - 9.0% of mass 95	6.8
173	Less than 2.0% of mass 174	1.1 (1.6) 1
174	50.0 - 100.0% of mass 95	66.1
175	5.0 - 9.0% of mass 174	5 (7.5) 1
176	95.0 - 101.0% of mass 174	63.6 (96.3) 1
177	5.0 - 9.0% of mass 176	4.5 (7) 2

1-Value is % mass 174 2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
VSTDCCC050	VSTDCCC050	VX046636.D	06/11/2025	14:46
VX0611WBL02	VX0611WBL02	VX046637.D	06/11/2025	15:14
VX0611WBS02	VX0611WBS02	VX046639.D	06/11/2025	15:58
GW-MW01-060425MS	Q2230-03MS	VX046641.D	06/11/2025	16:47
GW-MW01-060425MSD	Q2230-04MSD	VX046642.D	06/11/2025	17:09

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH Contract: CAMP02

Lab Code: CHEM Case No.: Q2230 SAS No.: Q2230 SDG NO.: Q2230

Lab File ID: VX046586.D Date Analyzed: 06/10/2025

Instrument ID: MSVOA_X Time Analyzed: 09:07

GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: (Y/N) N

	IS1 AREA #	RT #	IS2 AREA #	RT #	IS3 AREA #	RT #
12 HOUR STD	90758	5.56	150623	6.76	134767	10.06
UPPER LIMIT	181516	6.056	301246	7.263	269534	10.555
LOWER LIMIT	45379	5.056	75311.5	6.263	67383.5	9.555
EPA SAMPLE NO.						
FB-060425	109731	5.56	215111	6.77	220551	10.06
GW-MW01-060425	112673	5.56	225413	6.77	232357	10.06
GW-MW901-060425	112231	5.56	225068	6.77	232456	10.06
TB060425	114604	5.57	224360	6.77	234275	10.06
VX0610WBL01	106110	5.56	211553	6.77	217338	10.06
VX0610WBS01	84052	5.56	143506	6.77	131426	10.06

IS1 = Pentafluorobenzene
IS2 = 1,4-Difluorobenzene
IS3 = Chlorobenzene-d5

AREA UPPER LIMIT = +100% of internal standard area
AREA LOWER LIMIT = -50% of internal standard area
RT UPPER LIMIT = +0.50 minutes of internal standard RT
RT LOWER LIMIT = -0.50 minutes of internal standard RT

Column used to flag values outside QC limits with an asterisk.
* Values outside of QC limits.

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH Contract: CAMP02

Lab Code: CHEM Case No.: Q2230 SAS No.: Q2230 SDG NO.: Q2230

Lab File ID: VX046586.D Date Analyzed: 06/10/2025

Instrument ID: MSVOA_X Time Analyzed: 09:07

GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: (Y/N) N

	IS4 AREA #	RT #				
12 HOUR STD	66941	12.018				
UPPER LIMIT	133882	12.518				
LOWER LIMIT	33470.5	11.518				
EPA SAMPLE NO.						
FB-060425	108439	12.02				
GW-MW01-060425	117788	12.02				
GW-MW901-060425	116598	12.02				
TB060425	115739	12.02				
VX0610WBL01	104374	12.02				
VX0610WBS01	68222	12.02				

IS4 = 1,4-Dichlorobenzene-d4

AREA UPPER LIMIT = +100% of internal standard area
AREA LOWER LIMIT = -50% of internal standard area
RT UPPER LIMIT = +0.50 minutes of internal standard RT
RT LOWER LIMIT = -0.50 minutes of internal standard RT

Column used to flag values outside QC limits with an asterisk.
* Values outside of QC limits.

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH Contract: CAMP02

Lab Code: CHEM Case No.: Q2230 SAS No.: Q2230 SDG NO.: Q2230

Lab File ID: VX046636.D Date Analyzed: 06/11/2025

Instrument ID: MSVOA_X Time Analyzed: 14:46

GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: (Y/N) N

	IS1 AREA #	RT #	IS2 AREA #	RT #	IS3 AREA #	RT #
12 HOUR STD	77468	5.56	133266	6.76	119186	10.06
UPPER LIMIT	154936	6.062	266532	7.263	238372	10.555
LOWER LIMIT	38734	5.062	66633	6.263	59593	9.555
EPA SAMPLE NO.						
GW-MW01-060425MS	80829	5.57	141022	6.77	125449	10.06
GW-MW01-060425MSD	63422	5.57	110741	6.78	101042	10.06
VX0611WBL02	73470	5.56	147517	6.77	153084	10.06
VX0611WBS02	79398	5.57	136732	6.78	125565	10.06

IS1 = Pentafluorobenzene
IS2 = 1,4-Difluorobenzene
IS3 = Chlorobenzene-d5

AREA UPPER LIMIT = +100% of internal standard area
AREA LOWER LIMIT = -50% of internal standard area
RT UPPER LIMIT = +0.50 minutes of internal standard RT
RT LOWER LIMIT = -0.50 minutes of internal standard RT

Column used to flag values outside QC limits with an asterisk.
* Values outside of QC limits.

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH Contract: CAMP02

Lab Code: CHEM Case No.: Q2230 SAS No.: Q2230 SDG NO.: Q2230

Lab File ID: VX046636.D Date Analyzed: 06/11/2025

Instrument ID: MSVOA_X Time Analyzed: 14:46

GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: (Y/N) N

	IS4 AREA #	RT #				
12 HOUR STD	60295	12.018				
UPPER LIMIT	120590	12.518				
LOWER LIMIT	30147.5	11.518				
EPA SAMPLE NO.						
GW-MW01-060425MS	63001	12.02				
GW-MW01-060425MSD	50864	12.02				
VX0611WBL02	77047	12.02				
VX0611WBS02	66542	12.02				

IS4 = 1,4-Dichlorobenzene-d4

AREA UPPER LIMIT = +100% of internal standard area
AREA LOWER LIMIT = -50% of internal standard area
RT UPPER LIMIT = +0.50 minutes of internal standard RT
RT LOWER LIMIT = -0.50 minutes of internal standard RT

Column used to flag values outside QC limits with an asterisk.
* Values outside of QC limits.

QC SAMPLE DATA

Alliance TECHNICAL GROUP
284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Report of Analysis

Client:	CDM Smith	Date Collected:	
Project:	Con Ed UTEN Mount Vernon, NY	Date Received:	
Client Sample ID:	VX0610WBL01	SDG No.:	Q2230
Lab Sample ID:	VX0610WBL01	Matrix:	Water
Analytical Method:	8260D	% Solid:	0
Sample Wt/Vol:	5 Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOCMS Group3
GC Column:	DB-624UI ID : 0.18	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VX046588.D	1		06/10/25 09:57	VX061025

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
71-43-2	Benzene	1.00	U	0.15	1.00	ug/L
108-88-3	Toluene	1.00	U	0.14	1.00	ug/L
100-41-4	Ethyl Benzene	1.00	U	0.13	1.00	ug/L
1330-20-7	Total Xylenes	3.00	U	0.36	3.00	ug/L
98-82-8	Isopropylbenzene	1.00	U	0.12	1.00	ug/L
103-65-1	n-propylbenzene	1.00	U	0.13	1.00	ug/L
108-67-8	1,3,5-Trimethylbenzene	1.00	U	0.15	1.00	ug/L
98-06-6	tert-Butylbenzene	1.00	U	0.14	1.00	ug/L
95-63-6	1,2,4-Trimethylbenzene	1.00	U	0.14	1.00	ug/L
135-98-8	sec-Butylbenzene	1.00	U	0.13	1.00	ug/L
99-87-6	p-Isopropyltoluene	1.00	U	0.13	1.00	ug/L
104-51-8	n-Butylbenzene	1.00	U	0.15	1.00	ug/L
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	50.1		74 - 125	100%	SPK: 50
1868-53-7	Dibromofluoromethane	50.1		75 - 124	100%	SPK: 50
2037-26-5	Toluene-d8	54.3		86 - 113	109%	SPK: 50
460-00-4	4-Bromofluorobenzene	54.8		77 - 121	110%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	106000	5.562			
540-36-3	1,4-Difluorobenzene	212000	6.769			
3114-55-4	Chlorobenzene-d5	217000	10.055			
3855-82-1	1,4-Dichlorobenzene-d4	104000	12.018			

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

Alliance TECHNICAL GROUP
284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Report of Analysis

Client:	CDM Smith	Date Collected:	
Project:	Con Ed UTEN Mount Vernon, NY	Date Received:	
Client Sample ID:	VX0611WBL02	SDG No.:	Q2230
Lab Sample ID:	VX0611WBL02	Matrix:	Water
Analytical Method:	8260D	% Solid:	0
Sample Wt/Vol:	5 Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOCMS Group3
GC Column:	DB-624UI ID : 0.18	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VX046637.D	1		06/11/25 15:14	VX061125

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
71-43-2	Benzene	1.00	U	0.15	1.00	ug/L
108-88-3	Toluene	1.00	U	0.14	1.00	ug/L
100-41-4	Ethyl Benzene	1.00	U	0.13	1.00	ug/L
1330-20-7	Total Xylenes	3.00	U	0.36	3.00	ug/L
98-82-8	Isopropylbenzene	1.00	U	0.12	1.00	ug/L
103-65-1	n-propylbenzene	1.00	U	0.13	1.00	ug/L
108-67-8	1,3,5-Trimethylbenzene	1.00	U	0.15	1.00	ug/L
98-06-6	tert-Butylbenzene	1.00	U	0.14	1.00	ug/L
95-63-6	1,2,4-Trimethylbenzene	1.00	U	0.14	1.00	ug/L
135-98-8	sec-Butylbenzene	1.00	U	0.13	1.00	ug/L
99-87-6	p-Isopropyltoluene	1.00	U	0.13	1.00	ug/L
104-51-8	n-Butylbenzene	1.00	U	0.15	1.00	ug/L
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	49.7		74 - 125	99%	SPK: 50
1868-53-7	Dibromofluoromethane	50.4		75 - 124	101%	SPK: 50
2037-26-5	Toluene-d8	54.3		86 - 113	109%	SPK: 50
460-00-4	4-Bromofluorobenzene	55.3		77 - 121	111%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	73500	5.556			
540-36-3	1,4-Difluorobenzene	148000	6.769			
3114-55-4	Chlorobenzene-d5	153000	10.055			
3855-82-1	1,4-Dichlorobenzene-d4	77000	12.018			

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

Alliance TECHNICAL GROUP
284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Report of Analysis

Client:	CDM Smith	Date Collected:	
Project:	Con Ed UTEN Mount Vernon, NY	Date Received:	
Client Sample ID:	VX0610WBS01	SDG No.:	Q2230
Lab Sample ID:	VX0610WBS01	Matrix:	Water
Analytical Method:	8260D	% Solid:	0
Sample Wt/Vol:	5 Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOCMS Group3
GC Column:	DB-624UI ID : 0.18	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VX046589.D	1		06/10/25 10:18	VX061025

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
71-43-2	Benzene	21.0		0.15	1.00	ug/L
108-88-3	Toluene	21.2		0.14	1.00	ug/L
100-41-4	Ethyl Benzene	20.3		0.13	1.00	ug/L
1330-20-7	Total Xylenes	62.7		0.36	3.00	ug/L
98-82-8	Isopropylbenzene	20.2		0.12	1.00	ug/L
103-65-1	n-propylbenzene	20.0		0.13	1.00	ug/L
108-67-8	1,3,5-Trimethylbenzene	20.5		0.15	1.00	ug/L
98-06-6	tert-Butylbenzene	20.4		0.14	1.00	ug/L
95-63-6	1,2,4-Trimethylbenzene	20.8		0.14	1.00	ug/L
135-98-8	sec-Butylbenzene	20.5		0.13	1.00	ug/L
99-87-6	p-Isopropyltoluene	20.2		0.13	1.00	ug/L
104-51-8	n-Butylbenzene	19.9		0.15	1.00	ug/L
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	47.4		74 - 125	95%	SPK: 50
1868-53-7	Dibromofluoromethane	53.0		75 - 124	106%	SPK: 50
2037-26-5	Toluene-d8	51.5		86 - 113	103%	SPK: 50
460-00-4	4-Bromofluorobenzene	53.1		77 - 121	106%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	84100	5.562			
540-36-3	1,4-Difluorobenzene	144000	6.769			
3114-55-4	Chlorobenzene-d5	131000	10.055			
3855-82-1	1,4-Dichlorobenzene-d4	68200	12.018			

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

Alliance TECHNICAL GROUP
284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Report of Analysis

Client:	CDM Smith	Date Collected:	
Project:	Con Ed UTEN Mount Vernon, NY	Date Received:	
Client Sample ID:	VX0611WBS02	SDG No.:	Q2230
Lab Sample ID:	VX0611WBS02	Matrix:	Water
Analytical Method:	8260D	% Solid:	0
Sample Wt/Vol:	5 Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOCMS Group3
GC Column:	DB-624UI ID : 0.18	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VX046639.D	1		06/11/25 15:58	VX061125

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
71-43-2	Benzene	19.6		0.15	1.00	ug/L
108-88-3	Toluene	20.2		0.14	1.00	ug/L
100-41-4	Ethyl Benzene	19.2		0.13	1.00	ug/L
1330-20-7	Total Xylenes	60.2		0.36	3.00	ug/L
98-82-8	Isopropylbenzene	19.1		0.12	1.00	ug/L
103-65-1	n-propylbenzene	18.8		0.13	1.00	ug/L
108-67-8	1,3,5-Trimethylbenzene	18.9		0.15	1.00	ug/L
98-06-6	tert-Butylbenzene	19.2		0.14	1.00	ug/L
95-63-6	1,2,4-Trimethylbenzene	19.2		0.14	1.00	ug/L
135-98-8	sec-Butylbenzene	19.3		0.13	1.00	ug/L
99-87-6	p-Isopropyltoluene	18.9		0.13	1.00	ug/L
104-51-8	n-Butylbenzene	18.3		0.15	1.00	ug/L
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	44.0		74 - 125	88%	SPK: 50
1868-53-7	Dibromofluoromethane	47.8		75 - 124	96%	SPK: 50
2037-26-5	Toluene-d8	47.2		86 - 113	94%	SPK: 50
460-00-4	4-Bromofluorobenzene	48.5		77 - 121	97%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	79400	5.568			
540-36-3	1,4-Difluorobenzene	137000	6.775			
3114-55-4	Chlorobenzene-d5	126000	10.055			
3855-82-1	1,4-Dichlorobenzene-d4	66500	12.018			

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

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Report of Analysis

Client:	CDM Smith	Date Collected:	06/04/25
Project:	Con Ed UTEN Mount Vernon, NY	Date Received:	06/04/25
Client Sample ID:	GW-MW01-060425MS	SDG No.:	Q2230
Lab Sample ID:	Q2230-03MS	Matrix:	Water
Analytical Method:	8260D	% Solid:	0
Sample Wt/Vol:	5 Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOCMS Group3
GC Column:	DB-624UI ID : 0.18	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VX046641.D	1		06/11/25 16:47	VX061125

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
71-43-2	Benzene	58.1		0.15	1.00	ug/L
108-88-3	Toluene	62.2		0.14	1.00	ug/L
100-41-4	Ethyl Benzene	65.8		0.13	1.00	ug/L
1330-20-7	Total Xylenes	245		0.36	3.00	ug/L
98-82-8	Isopropylbenzene	62.9		0.12	1.00	ug/L
103-65-1	n-propylbenzene	63.6		0.13	1.00	ug/L
108-67-8	1,3,5-Trimethylbenzene	72.5		0.15	1.00	ug/L
98-06-6	tert-Butylbenzene	61.8		0.14	1.00	ug/L
95-63-6	1,2,4-Trimethylbenzene	100		0.14	1.00	ug/L
135-98-8	sec-Butylbenzene	60.8		0.13	1.00	ug/L
99-87-6	p-Isopropyltoluene	60.8		0.13	1.00	ug/L
104-51-8	n-Butylbenzene	60.9		0.15	1.00	ug/L
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	47.5		74 - 125	95%	SPK: 50
1868-53-7	Dibromofluoromethane	52.2		75 - 124	104%	SPK: 50
2037-26-5	Toluene-d8	51.1		86 - 113	102%	SPK: 50
460-00-4	4-Bromofluorobenzene	50.9		77 - 121	102%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	80800	5.568			
540-36-3	1,4-Difluorobenzene	141000	6.769			
3114-55-4	Chlorobenzene-d5	125000	10.055			
3855-82-1	1,4-Dichlorobenzene-d4	63000	12.018			

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

Alliance TECHNICAL GROUP
284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

Report of Analysis

Client:	CDM Smith	Date Collected:	06/04/25
Project:	Con Ed UTEN Mount Vernon, NY	Date Received:	06/04/25
Client Sample ID:	GW-MW01-060425MSD	SDG No.:	Q2230
Lab Sample ID:	Q2230-04MSD	Matrix:	Water
Analytical Method:	8260D	% Solid:	0
Sample Wt/Vol:	5 Units: mL	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOCMS Group3
GC Column:	DB-624UI ID : 0.18	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VX046642.D	1		06/11/25 17:09	VX061125

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
71-43-2	Benzene	50.7		0.15	1.00	ug/L
108-88-3	Toluene	54.6		0.14	1.00	ug/L
100-41-4	Ethyl Benzene	58.5		0.13	1.00	ug/L
1330-20-7	Total Xylenes	236		0.36	3.00	ug/L
98-82-8	Isopropylbenzene	53.5		0.12	1.00	ug/L
103-65-1	n-propylbenzene	55.2		0.13	1.00	ug/L
108-67-8	1,3,5-Trimethylbenzene	66.8		0.15	1.00	ug/L
98-06-6	tert-Butylbenzene	52.1		0.14	1.00	ug/L
95-63-6	1,2,4-Trimethylbenzene	110		0.14	1.00	ug/L
135-98-8	sec-Butylbenzene	52.0		0.13	1.00	ug/L
99-87-6	p-Isopropyltoluene	51.8		0.13	1.00	ug/L
104-51-8	n-Butylbenzene	53.5		0.15	1.00	ug/L
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	45.9		74 - 125	92%	SPK: 50
1868-53-7	Dibromofluoromethane	50.4		75 - 124	101%	SPK: 50
2037-26-5	Toluene-d8	50.2		86 - 113	100%	SPK: 50
460-00-4	4-Bromofluorobenzene	51.1		77 - 121	102%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	63400	5.568			
540-36-3	1,4-Difluorobenzene	111000	6.775			
3114-55-4	Chlorobenzene-d5	101000	10.055			
3855-82-1	1,4-Dichlorobenzene-d4	50900	12.018			

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

CALIBRATION SUMMARY

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: CAMP02

Lab Code: CHEM Case No.: Q2230 SAS No.: Q2230 SDG No.: Q2230

Instrument ID: MSVOA_X Calibration Date(s): 06/06/2025 06/06/2025

Heated Purge: (Y/N) N Calibration Time(s): 09:42 12:57

GC Column: DB-624UI ID: 0.18 (mm)

LAB FILE ID: RRF005 = VX046518.D RRF020 = VX046519.D RRF050 = VX046520.D
RRF100 = VX046521.D RRF150 = VX046522.D RRF001 = VX046524.D

COMPOUND	RRF005	RRF020	RRF050	RRF100	RRF150	RRF001	$\overline{RRF}$	% RSD
Benzene	1.597	1.503	1.427	1.380	1.442	1.522	1.479	5.3
Toluene	0.980	0.911	0.884	0.849	0.888	0.938	0.908	5
Ethyl Benzene	2.093	2.029	2.018	1.971	2.041	2.166	2.053	3.3
m/p-Xylenes	0.756	0.750	0.737	0.714	0.738	0.773	0.745	2.7
o-Xylene	0.728	0.731	0.716	0.703	0.724	0.719	0.720	1.4
Isopropylbenzene	3.936	3.949	3.909	3.828	4.064	3.896	3.930	2
n-propylbenzene	4.659	4.775	4.693	4.553	4.813	5.051	4.757	3.6
1,3,5-Trimethylbenzene	3.329	3.297	3.243	3.147	3.314	3.322	3.275	2.1
tert-Butylbenzene	3.341	3.364	3.287	3.248	3.482	3.284	3.334	2.5
1,2,4-Trimethylbenzene	3.309	3.360	3.256	3.194	3.356	3.406	3.313	2.3
sec-Butylbenzene	4.350	4.296	4.172	4.096	4.334	4.458	4.284	3
p-Isopropyltoluene	3.509	3.554	3.484	3.397	3.575	3.910	3.572	5
n-Butylbenzene	3.319	3.378	3.354	3.310	3.487	4.197	3.507	9.8
1,2-Dichloroethane-d4	0.997	0.848	0.873	0.828	0.900		0.890	7.4
Dibromofluoromethane	0.392	0.353	0.368	0.347	0.379		0.368	5
Toluene-d8	1.362	1.159	1.188	1.132	1.220		1.212	7.4
4-Bromofluorobenzene	0.564	0.482	0.493	0.468	0.501		0.502	7.4

* Compounds with required minimum RRF and maximum %RSD values.
All other compounds must meet a minimum RRF of 0.010.
RRF of 1,4-Dioxane = Value should be divide by 1000.

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: CAMP02

Lab Code: CHEM Case No.: Q2230 SAS No.: Q2230 SDG No.: Q2230

Instrument ID: MSVOA_X Calibration Date/Time: 06/10/2025 09:07

Lab File ID: VX046586.D Init. Calib. Date(s): 06/06/2025 06/06/2025

Heated Purge: (Y/N) N Init. Calib. Time(s): 09:42 12:57

GC Column: DB-624UI ID: 0.18 (mm)

COMPOUND	RRF	RRF050	MIN RRF	%D	MAX%D
Benzene	1.479	1.615		9.19	20
Toluene	0.908	0.992		9.25	20
Ethyl Benzene	2.053	2.179		6.14	20
m/p-Xylenes	0.745	0.802		7.65	20
o-Xylene	0.720	0.775		7.64	20
Isopropylbenzene	3.930	4.253		8.22	20
n-propylbenzene	4.757	5.071		6.6	20
1,3,5-Trimethylbenzene	3.275	3.526		7.66	20
tert-Butylbenzene	3.334	3.617		8.49	20
1,2,4-Trimethylbenzene	3.313	3.499		5.61	20
sec-Butylbenzene	4.284	4.548		6.16	20
p-Isopropyltoluene	3.572	3.774		5.66	20
n-Butylbenzene	3.507	3.673		4.73	20
1,2-Dichloroethane-d4	0.890	0.757		-14.94	20
Dibromofluoromethane	0.368	0.360		-2.17	20
Toluene-d8	1.212	1.152		-4.95	20
4-Bromofluorobenzene	0.502	0.468		-6.77	20

All other compounds must meet a minimum RRF of 0.010.
RRF of 1,4-Dioxane = Value should be divide by 1000.

284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: CAMP02

Lab Code: CHEM Case No.: Q2230 SAS No.: Q2230 SDG No.: Q2230

Instrument ID: MSVOA_X Calibration Date/Time: 06/11/2025 14:46

Lab File ID: VX046636.D Init. Calib. Date(s): 06/06/2025 06/06/2025

Heated Purge: (Y/N) N Init. Calib. Time(s): 09:42 12:57

GC Column: DB-624UI ID: 0.18 (mm)

COMPOUND	RRF	RRF050	MIN RRF	%D	MAX%D
Benzene	1.479	1.534		3.72	20
Toluene	0.908	0.958		5.51	20
Ethyl Benzene	2.053	2.124		3.46	20
m/p-Xylenes	0.745	0.792		6.31	20
o-Xylene	0.720	0.786		9.17	20
Isopropylbenzene	3.930	4.163		5.93	20
n-propylbenzene	4.757	4.937		3.78	20
1,3,5-Trimethylbenzene	3.275	3.443		5.13	20
tert-Butylbenzene	3.334	3.554		6.6	20
1,2,4-Trimethylbenzene	3.313	3.476		4.92	20
sec-Butylbenzene	4.284	4.505		5.16	20
p-Isopropyltoluene	3.572	3.722		4.2	20
n-Butylbenzene	3.507	3.559		1.48	20
1,2-Dichloroethane-d4	0.890	0.830		-6.74	20
Dibromofluoromethane	0.368	0.382		3.8	20
Toluene-d8	1.212	1.219		0.58	20
4-Bromofluorobenzene	0.502	0.503		0.2	20

All other compounds must meet a minimum RRF of 0.010.
RRF of 1,4-Dioxane = Value should be divide by 1000.