

Hit Summary Sheet SW-846

SDG No.: Q1761
Client: G Environmental

Sample ID	Client ID	Matrix	Parameter	Concentration	C	MDL	RDL	Units
Client ID:	GST3							
Q1761-01	GST3	SOIL	Benzene, 4-ethyl-1,2-dimethyl- *	4.90	J	0	0	ug/Kg
			Total Tics :	4.90				
			Total Concentration:	4.90				



SAMPLE DATA

Report of Analysis

Client:	G Environmental	Date Collected:	04/08/25
Project:	Stockton	Date Received:	04/09/25
Client Sample ID:	GST3	SDG No.:	Q1761
Lab Sample ID:	Q1761-01	Matrix:	SOIL
Analytical Method:	SW8260	% Solid:	78.6
Sample Wt/Vol:	7.82 Units: g	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOC-TCLVOA-10
GC Column:	RXI-624 ID : 0.25	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VY021855.D	1		04/11/25 13:33	VY041125

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
75-71-8	Dichlorodifluoromethane	0.93	U	0.93	4.10	ug/Kg
74-87-3	Chloromethane	0.93	U	0.93	4.10	ug/Kg
75-01-4	Vinyl Chloride	0.64	U	0.64	4.10	ug/Kg
74-83-9	Bromomethane	0.87	U	0.87	4.10	ug/Kg
75-00-3	Chloroethane	1.00	U	1.00	4.10	ug/Kg
75-69-4	Trichlorofluoromethane	0.98	U	0.98	4.10	ug/Kg
76-13-1	1,1,2-Trichlorotrifluoroethane	0.86	U	0.86	4.10	ug/Kg
75-35-4	1,1-Dichloroethene	0.81	U	0.81	4.10	ug/Kg
67-64-1	Acetone	3.90	U	3.90	20.3	ug/Kg
75-15-0	Carbon Disulfide	0.86	U	0.86	4.10	ug/Kg
1634-04-4	Methyl tert-butyl Ether	0.59	U	0.59	4.10	ug/Kg
79-20-9	Methyl Acetate	1.30	U	1.30	4.10	ug/Kg
75-09-2	Methylene Chloride	2.90	U	2.90	8.10	ug/Kg
156-60-5	trans-1,2-Dichloroethene	0.70	U	0.70	4.10	ug/Kg
75-34-3	1,1-Dichloroethane	0.65	U	0.65	4.10	ug/Kg
110-82-7	Cyclohexane	0.64	U	0.64	4.10	ug/Kg
78-93-3	2-Butanone	5.30	U	5.30	20.3	ug/Kg
56-23-5	Carbon Tetrachloride	0.79	U	0.79	4.10	ug/Kg
156-59-2	cis-1,2-Dichloroethene	0.61	U	0.61	4.10	ug/Kg
74-97-5	Bromochloromethane	0.94	U	0.94	4.10	ug/Kg
67-66-3	Chloroform	0.68	U	0.68	4.10	ug/Kg
71-55-6	1,1,1-Trichloroethane	0.76	U	0.76	4.10	ug/Kg
108-87-2	Methylcyclohexane	0.74	U	0.74	4.10	ug/Kg
71-43-2	Benzene	0.64	U	0.64	4.10	ug/Kg
107-06-2	1,2-Dichloroethane	0.64	U	0.64	4.10	ug/Kg
79-01-6	Trichloroethene	0.66	U	0.66	4.10	ug/Kg
78-87-5	1,2-Dichloropropane	0.74	U	0.74	4.10	ug/Kg
75-27-4	Bromodichloromethane	0.63	U	0.63	4.10	ug/Kg
108-10-1	4-Methyl-2-Pentanone	2.90	U	2.90	20.3	ug/Kg
108-88-3	Toluene	0.63	U	0.63	4.10	ug/Kg

Report of Analysis

Client:	G Environmental		Date Collected:	04/08/25	
Project:	Stockton		Date Received:	04/09/25	
Client Sample ID:	GST3		SDG No.:	Q1761	
Lab Sample ID:	Q1761-01		Matrix:	SOIL	
Analytical Method:	SW8260		% Solid:	78.6	
Sample Wt/Vol:	7.82	Units: g	Final Vol:	5000	uL
Soil Aliquot Vol:		uL	Test:	VOC-TCLVOA-10	
GC Column:	RXI-624	ID : 0.25	Level :	LOW	
Prep Method :					

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VY021855.D	1		04/11/25 13:33	VY041125

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
10061-02-6	t-1,3-Dichloropropene	0.53	U	0.53	4.10	ug/Kg
10061-01-5	cis-1,3-Dichloropropene	0.50	U	0.50	4.10	ug/Kg
79-00-5	1,1,2-Trichloroethane	0.75	U	0.75	4.10	ug/Kg
591-78-6	2-Hexanone	3.00	U	3.00	20.3	ug/Kg
124-48-1	Dibromochloromethane	0.71	U	0.71	4.10	ug/Kg
106-93-4	1,2-Dibromoethane	0.72	U	0.72	4.10	ug/Kg
127-18-4	Tetrachloroethene	0.85	U	0.85	4.10	ug/Kg
108-90-7	Chlorobenzene	0.74	U	0.74	4.10	ug/Kg
100-41-4	Ethyl Benzene	0.55	U	0.55	4.10	ug/Kg
179601-23-1	m/p-Xylenes	1.00	U	1.00	8.10	ug/Kg
95-47-6	o-Xylene	0.67	U	0.67	4.10	ug/Kg
100-42-5	Styrene	0.58	U	0.58	4.10	ug/Kg
75-25-2	Bromoform	0.70	U	0.70	4.10	ug/Kg
98-82-8	Isopropylbenzene	0.63	U	0.63	4.10	ug/Kg
79-34-5	1,1,2,2-Tetrachloroethane	0.98	U	0.98	4.10	ug/Kg
541-73-1	1,3-Dichlorobenzene	1.40	U	1.40	4.10	ug/Kg
106-46-7	1,4-Dichlorobenzene	1.30	U	1.30	4.10	ug/Kg
95-50-1	1,2-Dichlorobenzene	1.20	U	1.20	4.10	ug/Kg
96-12-8	1,2-Dibromo-3-Chloropropane	1.50	U	1.50	4.10	ug/Kg
120-82-1	1,2,4-Trichlorobenzene	2.40	U	2.40	4.10	ug/Kg
87-61-6	1,2,3-Trichlorobenzene	2.60	U	2.60	4.10	ug/Kg
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	46.6		70 (63) - 130 (155)	93%	SPK: 50
1868-53-7	Dibromofluoromethane	49.6		70 (70) - 130 (134)	99%	SPK: 50
2037-26-5	Toluene-d8	47.9		70 (74) - 130 (123)	96%	SPK: 50
460-00-4	4-Bromofluorobenzene	40.5		70 (38) - 130 (136)	81%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	317000	7.713			
540-36-3	1,4-Difluorobenzene	560000	8.616			
3114-55-4	Chlorobenzene-d5	469000	11.42			
3855-82-1	1,4-Dichlorobenzene-d4	180000	13.346			

TENTATIVE IDENTIFIED COMPOUNDS

Report of Analysis

Client:	G Environmental		Date Collected:	04/08/25	
Project:	Stockton		Date Received:	04/09/25	
Client Sample ID:	GST3		SDG No.:	Q1761	
Lab Sample ID:	Q1761-01		Matrix:	SOIL	
Analytical Method:	SW8260		% Solid:	78.6	
Sample Wt/Vol:	7.82	Units: g	Final Vol:	5000	uL
Soil Aliquot Vol:		uL	Test:	VOC-TCLVOA-10	
GC Column:	RXI-624	ID : 0.25	Level :	LOW	
Prep Method :					

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VY021855.D	1		04/11/25 13:33	VY041125

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
000934-80-5	Benzene, 4-ethyl-1,2-dimethyl-	4.90	J		13.9	ug/Kg

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

Surrogate Summary

SDG No.: Q1761

Client: G Environmental

Analytical Method: SW8260D

Lab Sample ID	Client ID	Parameter	Spike	Result	RecoveryQual	Limits	
						Low	High
Q1761-01	GST3	1,2-Dichloroethane-d4	50	46.6	93	70 (63)	130 (155)
		Dibromofluoromethane	50	49.6	99	70 (70)	130 (134)
		Toluene-d8	50	48.0	96	70 (74)	130 (123)
		4-Bromofluorobenzene	50	40.5	81	70 (38)	130 (136)
VY0411SBL01	VY0411SBL01	1,2-Dichloroethane-d4	50	47.0	94	70 (63)	130 (155)
		Dibromofluoromethane	50	49.2	98	70 (70)	130 (134)
		Toluene-d8	50	47.4	95	70 (74)	130 (123)
		4-Bromofluorobenzene	50	39.1	78	70 (38)	130 (136)
VY0411SBS01	VY0411SBS01	1,2-Dichloroethane-d4	50	52.3	105	70 (63)	130 (155)
		Dibromofluoromethane	50	53.2	106	70 (70)	130 (134)
		Toluene-d8	50	53.8	108	70 (74)	130 (123)
		4-Bromofluorobenzene	50	55.3	111	70 (38)	130 (136)
VY0411SBSD01	VY0411SBSD01	1,2-Dichloroethane-d4	50	50.6	101	70 (63)	130 (155)
		Dibromofluoromethane	50	53.9	108	70 (70)	130 (134)
		Toluene-d8	50	53.1	106	70 (74)	130 (123)
		4-Bromofluorobenzene	50	54.5	109	70 (38)	130 (136)

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q1761

Client: G Environmental

Analytical Method: SW8260D

Datafile : VY021852.D

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	Low	Limits High	RPD
VY0411SBS01	Dichlorodifluoromethane	20	17.7	ug/Kg	89			40 (64)	160 (136)	
	Chloromethane	20	17.5	ug/Kg	88			40 (70)	160 (130)	
	Vinyl chloride	20	17.2	ug/Kg	86			70 (72)	130 (129)	
	Bromomethane	20	18.9	ug/Kg	95			40 (58)	160 (141)	
	Chloroethane	20	18.0	ug/Kg	90			40 (69)	160 (130)	
	Trichlorofluoromethane	20	18.6	ug/Kg	93			40 (69)	160 (134)	
	1,1,2-Trichlorotrifluoroethane	20	19.6	ug/Kg	98			70 (81)	130 (123)	
	1,1-Dichloroethene	20	19.1	ug/Kg	96			70 (79)	130 (121)	
	Acetone	100	100	ug/Kg	100			40 (60)	160 (131)	
	Carbon disulfide	20	17.4	ug/Kg	87			40 (45)	160 (154)	
	Methyl tert-butyl Ether	20	19.8	ug/Kg	99			70 (77)	130 (129)	
	Methyl Acetate	20	20.7	ug/Kg	104			70 (69)	130 (149)	
	Methylene Chloride	20	20.1	ug/Kg	101			70 (56)	130 (174)	
	trans-1,2-Dichloroethene	20	18.9	ug/Kg	95			70 (80)	130 (123)	
	1,1-Dichloroethane	20	19.7	ug/Kg	99			70 (82)	130 (123)	
	Cyclohexane	20	17.9	ug/Kg	90			70 (76)	130 (122)	
	2-Butanone	100	95.3	ug/Kg	95			40 (69)	160 (131)	
	Carbon Tetrachloride	20	19.7	ug/Kg	99			70 (76)	130 (129)	
	cis-1,2-Dichloroethene	20	19.9	ug/Kg	100			70 (82)	130 (123)	
	Bromochloromethane	20	19.7	ug/Kg	99			70 (80)	130 (127)	
	Chloroform	20	19.7	ug/Kg	99			70 (82)	130 (125)	
	1,1,1-Trichloroethane	20	19.6	ug/Kg	98			70 (80)	130 (126)	
	Methylcyclohexane	20	18.8	ug/Kg	94			70 (77)	130 (123)	
	Benzene	20	19.7	ug/Kg	99			70 (84)	130 (121)	
	1,2-Dichloroethane	20	19.8	ug/Kg	99			70 (81)	130 (126)	
	Trichloroethene	20	19.8	ug/Kg	99			70 (83)	130 (122)	
	1,2-Dichloropropane	20	19.5	ug/Kg	98			70 (83)	130 (122)	
	Bromodichloromethane	20	20.0	ug/Kg	100			70 (82)	130 (123)	
	4-Methyl-2-Pentanone	100	97.2	ug/Kg	97			40 (70)	160 (135)	
	Toluene	20	20.1	ug/Kg	101			70 (83)	130 (122)	
	t-1,3-Dichloropropene	20	19.4	ug/Kg	97			70 (78)	130 (124)	
	cis-1,3-Dichloropropene	20	19.9	ug/Kg	100			70 (81)	130 (122)	
	1,1,2-Trichloroethane	20	20.1	ug/Kg	101			70 (82)	130 (125)	
	2-Hexanone	100	97.2	ug/Kg	97			40 (66)	160 (138)	
	Dibromochloromethane	20	20.3	ug/Kg	102			70 (79)	130 (125)	
	1,2-Dibromoethane	20	20.4	ug/Kg	102			70 (80)	130 (125)	
	Tetrachloroethene	20	22.1	ug/Kg	111			70 (83)	130 (125)	
	Chlorobenzene	20	19.8	ug/Kg	99			70 (84)	130 (122)	
	Ethyl Benzene	20	19.5	ug/Kg	98			70 (82)	130 (124)	
	m/p-Xylenes	40	39.5	ug/Kg	99			70 (83)	130 (124)	
	o-Xylene	20	20.0	ug/Kg	100			70 (83)	130 (123)	
	Styrene	20	19.8	ug/Kg	99			70 (82)	130 (124)	
	Bromoform	20	20.4	ug/Kg	102			70 (75)	130 (127)	
	Isopropylbenzene	20	19.0	ug/Kg	95			70 (82)	130 (124)	
	1,1,2,2-Tetrachloroethane	20	17.9	ug/Kg	90			70 (77)	130 (127)	
	1,3-Dichlorobenzene	20	19.4	ug/Kg	97			70 (83)	130 (122)	
	1,4-Dichlorobenzene	20	19.5	ug/Kg	98			70 (84)	130 (121)	
	1,2-Dichlorobenzene	20	20.1	ug/Kg	101			70 (83)	130 (124)	

() = LABORATORY INHOUSE LIMIT

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q1761

Client: G Environmental

Analytical Method: SW8260D

Datafile : VY021852.D

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	Low	Limits High	RPD
VY0411SBS01	1,2-Dibromo-3-Chloropropane	20	20.9	ug/Kg	104			40 (66)	160 (134)	
	1,2,4-Trichlorobenzene	20	21.1	ug/Kg	106			70 (78)	130 (127)	
	1,2,3-Trichlorobenzene	20	21.4	ug/Kg	107			70 (70)	130 (137)	

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q1761

Client: G Environmental

Analytical Method: SW8260D

Datafile : VY021853.D

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	Low	Limits High	RPD
VY0411SBSD01	Dichlorodifluoromethane	20	18.3	ug/Kg	92	3		40 (64)	160 (136)	30 (20)
	Chloromethane	20	16.4	ug/Kg	82	7		40 (70)	160 (130)	30 (20)
	Vinyl chloride	20	17.0	ug/Kg	85	1		70 (72)	130 (129)	30 (20)
	Bromomethane	20	17.5	ug/Kg	88	8		40 (58)	160 (141)	30 (20)
	Chloroethane	20	17.1	ug/Kg	86	5		40 (69)	160 (130)	30 (20)
	Trichlorofluoromethane	20	19.0	ug/Kg	95	2		40 (69)	160 (134)	30 (20)
	1,1,2-Trichlorotrifluoroethane	20	19.5	ug/Kg	98	0		70 (81)	130 (123)	30 (20)
	1,1-Dichloroethene	20	19.8	ug/Kg	99	3		70 (79)	130 (121)	30 (20)
	Acetone	100	84.8	ug/Kg	85	16		40 (60)	160 (131)	30 (20)
	Carbon disulfide	20	17.9	ug/Kg	90	3		40 (45)	160 (154)	30 (20)
	Methyl tert-butyl Ether	20	19.6	ug/Kg	98	1		70 (77)	130 (129)	30 (20)
	Methyl Acetate	20	21.8	ug/Kg	109	5		70 (69)	130 (149)	30 (20)
	Methylene Chloride	20	20.4	ug/Kg	102	1		70 (56)	130 (174)	30 (20)
	trans-1,2-Dichloroethene	20	19.5	ug/Kg	98	3		70 (80)	130 (123)	30 (20)
	1,1-Dichloroethane	20	19.9	ug/Kg	100	1		70 (82)	130 (123)	30 (20)
	Cyclohexane	20	18.2	ug/Kg	91	1		70 (76)	130 (122)	30 (20)
	2-Butanone	100	90.6	ug/Kg	91	4		40 (69)	160 (131)	30 (20)
	Carbon Tetrachloride	20	20.2	ug/Kg	101	2		70 (76)	130 (129)	30 (20)
	cis-1,2-Dichloroethene	20	20.2	ug/Kg	101	1		70 (82)	130 (123)	30 (20)
	Bromochloromethane	20	18.0	ug/Kg	90	10		70 (80)	130 (127)	30 (20)
	Chloroform	20	20.0	ug/Kg	100	1		70 (82)	130 (125)	30 (20)
	1,1,1-Trichloroethane	20	20.1	ug/Kg	101	3		70 (80)	130 (126)	30 (20)
	Methylcyclohexane	20	19.1	ug/Kg	96	2		70 (77)	130 (123)	30 (20)
	Benzene	20	20.3	ug/Kg	102	3		70 (84)	130 (121)	30 (20)
	1,2-Dichloroethane	20	19.6	ug/Kg	98	1		70 (81)	130 (126)	30 (20)
	Trichloroethene	20	20.9	ug/Kg	104	5		70 (83)	130 (122)	30 (20)
	1,2-Dichloropropane	20	19.9	ug/Kg	100	2		70 (83)	130 (122)	30 (20)
	Bromodichloromethane	20	20.0	ug/Kg	100	0		70 (82)	130 (123)	30 (20)
	4-Methyl-2-Pentanone	100	94.4	ug/Kg	94	3		40 (70)	160 (135)	30 (20)
	Toluene	20	20.5	ug/Kg	103	2		70 (83)	130 (122)	30 (20)
	t-1,3-Dichloropropene	20	19.9	ug/Kg	100	3		70 (78)	130 (124)	30 (20)
	cis-1,3-Dichloropropene	20	19.6	ug/Kg	98	2		70 (81)	130 (122)	30 (20)
	1,1,2-Trichloroethane	20	20.6	ug/Kg	103	2		70 (82)	130 (125)	30 (20)
	2-Hexanone	100	93.4	ug/Kg	93	4		40 (66)	160 (138)	30 (20)
	Dibromochloromethane	20	20.5	ug/Kg	103	1		70 (79)	130 (125)	30 (20)
	1,2-Dibromoethane	20	19.9	ug/Kg	100	2		70 (80)	130 (125)	30 (20)
	Tetrachloroethene	20	22.4	ug/Kg	112	1		70 (83)	130 (125)	30 (20)
	Chlorobenzene	20	20.4	ug/Kg	102	3		70 (84)	130 (122)	30 (20)
	Ethyl Benzene	20	20.1	ug/Kg	101	3		70 (82)	130 (124)	30 (20)
	m/p-Xylenes	40	40.8	ug/Kg	102	3		70 (83)	130 (124)	30 (20)
	o-Xylene	20	20.5	ug/Kg	103	3		70 (83)	130 (123)	30 (20)
	Styrene	20	20.6	ug/Kg	103	4		70 (82)	130 (124)	30 (20)
	Bromoform	20	20.4	ug/Kg	102	0		70 (75)	130 (127)	30 (20)
	Isopropylbenzene	20	19.7	ug/Kg	99	4		70 (82)	130 (124)	30 (20)
	1,1,2,2-Tetrachloroethane	20	18.0	ug/Kg	90	0		70 (77)	130 (127)	30 (20)
	1,3-Dichlorobenzene	20	20.3	ug/Kg	102	5		70 (83)	130 (122)	30 (20)
	1,4-Dichlorobenzene	20	20.2	ug/Kg	101	3		70 (84)	130 (121)	30 (20)
	1,2-Dichlorobenzene	20	20.1	ug/Kg	101	0		70 (83)	130 (124)	30 (20)

() = LABORATORY INHOUSE LIMIT

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q1761

Client: G Environmental

Analytical Method: SW8260D

Datafile : VY021853.D

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	Low	Limits High	RPD
VY0411SBSD01	1,2-Dibromo-3-Chloropropane	20	19.9	ug/Kg	100	4		40 (66)	160 (134)	30 (20)
	1,2,4-Trichlorobenzene	20	21.9	ug/Kg	110	4		70 (78)	130 (127)	30 (20)
	1,2,3-Trichlorobenzene	20	22.5	ug/Kg	113	5		70 (70)	130 (137)	30 (20)

VOLATILE METHOD BLANK SUMMARY

EPA SAMPLE NO.

VY0411SBL01

Lab Name: CHEMTECH

Contract: GENV01

Lab Code: CHEM Case No.: Q1761

SAS No.: Q1761 SDG NO.: Q1761

Lab File ID: VY021851.D

Lab Sample ID: VY0411SBL01

Date Analyzed: 04/11/2025

Time Analyzed: 11:35

GC Column: RXI-624 ID: 0.25 (mm)

Heated Purge: (Y/N) Y

Instrument ID: MSVOA_Y

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

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
VY0411SBS01	VY0411SBS01	VY021852.D	04/11/2025
VY0411SBSD01	VY0411SBSD01	VY021853.D	04/11/2025
GST3	Q1761-01	VY021855.D	04/11/2025

COMMENTS: _____

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: CHEMTECH Contract: GENV01
Lab Code: CHEM Case No.: Q1761 SAS No.: Q1761 SDG NO.: Q1761
Lab File ID: VY021655.D BFB Injection Date: 03/27/2025
Instrument ID: MSVOA_Y BFB Injection Time: 09:23
GC Column: RXI-624 ID: 0.25 (mm) Heated Purge: Y/N Y

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	20.4
75	30.0 - 60.0% of mass 95	53
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	1.7 (2) 1
174	50.0 - 100.0% of mass 95	86.7
175	5.0 - 9.0% of mass 174	6.6 (7.6) 1
176	95.0 - 101.0% of mass 174	82.9 (95.6) 1
177	5.0 - 9.0% of mass 176	5.6 (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
VSTDIC005	VSTDIC005	VY021656.D	03/27/2025	10:08
VSTDIC010	VSTDIC010	VY021657.D	03/27/2025	10:31
VSTDIC020	VSTDIC020	VY021658.D	03/27/2025	10:54
VSTDIC050	VSTDIC050	VY021659.D	03/27/2025	11:16
VSTDIC100	VSTDIC100	VY021660.D	03/27/2025	11:39
VSTDIC150	VSTDIC150	VY021661.D	03/27/2025	12:02

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: CHEMTECH Contract: GENV01
Lab Code: CHEM Case No.: Q1761 SAS No.: Q1761 SDG NO.: Q1761
Lab File ID: VY021849.D BFB Injection Date: 04/11/2025
Instrument ID: MSVOA_Y BFB Injection Time: 08:46
GC Column: RXI-624 ID: 0.25 (mm) Heated Purge: Y/N Y

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	19.3
75	30.0 - 60.0% of mass 95	52
95	Base Peak, 100% relative abundance	100
96	5.0 - 9.0% of mass 95	6.6
173	Less than 2.0% of mass 174	1.2 (1.4) 1
174	50.0 - 100.0% of mass 95	87
175	5.0 - 9.0% of mass 174	6.6 (7.6) 1
176	95.0 - 101.0% of mass 174	82.8 (95.1) 1
177	5.0 - 9.0% of mass 176	5.6 (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	VY021850.D	04/11/2025	09:18
VY0411SBL01	VY0411SBL01	VY021851.D	04/11/2025	11:35
VY0411SBS01	VY0411SBS01	VY021852.D	04/11/2025	12:06
VY0411SBSD01	VY0411SBSD01	VY021853.D	04/11/2025	12:28
GST3	Q1761-01	VY021855.D	04/11/2025	13:33

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH Contract: GENV01
Lab Code: CHEM Case No.: Q1761 SAS No.: Q1761 SDG NO.: Q1761
Lab File ID: VY021850.D Date Analyzed: 04/11/2025
Instrument ID: MSVOA_Y Time Analyzed: 09:18
GC Column: RXI-624 ID: 0.25 (mm) Heated Purge: (Y/N) Y

	IS1 AREA #	RT #	IS2 AREA #	RT #	IS3 AREA #	RT #
12 HOUR STD	297119	7.71	443480	8.62	386894	11.42
UPPER LIMIT	594238	8.207	886960	9.115	773788	11.92
LOWER LIMIT	148560	7.207	221740	8.115	193447	10.92
EPA SAMPLE NO.						
GST3	317324	7.71	560087	8.62	469329	11.42
VY0411SBL01	321354	7.71	579974	8.62	477999	11.42
VY0411SBS01	261639	7.71	412441	8.62	367570	11.42
VY0411SBSD01	261813	7.71	409914	8.62	362626	11.41

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.

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH Contract: GENV01
Lab Code: CHEM Case No.: Q1761 SAS No.: Q1761 SDG NO.: Q1761
Lab File ID: VY021850.D Date Analyzed: 04/11/2025
Instrument ID: MSVOA_Y Time Analyzed: 09:18
GC Column: RXI-624 ID: 0.25 (mm) Heated Purge: (Y/N) Y

	IS4 AREA #	RT #				
12 HOUR STD	196263	13.346				
UPPER LIMIT	392526	13.846				
LOWER LIMIT	98131.5	12.846				
EPA SAMPLE NO.						
GST3	180059	13.35				
VY0411SBL01	174550	13.35				
VY0411SBS01	190311	13.35				
VY0411SBSD01	188335	13.35				

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

Report of Analysis

Client:	G Environmental	Date Collected:	
Project:	Stockton	Date Received:	
Client Sample ID:	VY0411SBL01	SDG No.:	Q1761
Lab Sample ID:	VY0411SBL01	Matrix:	SOIL
Analytical Method:	SW8260	% Solid:	100
Sample Wt/Vol:	5 Units: g	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOC-TCLVOA-10
GC Column:	RXI-624 ID : 0.25	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VY021851.D	1		04/11/25 11:35	VY041125

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
75-71-8	Dichlorodifluoromethane	1.10	U	1.10	5.00	ug/Kg
74-87-3	Chloromethane	1.10	U	1.10	5.00	ug/Kg
75-01-4	Vinyl Chloride	0.79	U	0.79	5.00	ug/Kg
74-83-9	Bromomethane	1.10	U	1.10	5.00	ug/Kg
75-00-3	Chloroethane	1.30	U	1.30	5.00	ug/Kg
75-69-4	Trichlorofluoromethane	1.20	U	1.20	5.00	ug/Kg
76-13-1	1,1,2-Trichlorotrifluoroethane	1.10	U	1.10	5.00	ug/Kg
75-35-4	1,1-Dichloroethene	1.00	U	1.00	5.00	ug/Kg
67-64-1	Acetone	4.70	U	4.70	25.0	ug/Kg
75-15-0	Carbon Disulfide	1.10	U	1.10	5.00	ug/Kg
1634-04-4	Methyl tert-butyl Ether	0.73	U	0.73	5.00	ug/Kg
79-20-9	Methyl Acetate	1.50	U	1.50	5.00	ug/Kg
75-09-2	Methylene Chloride	3.50	U	3.50	10.0	ug/Kg
156-60-5	trans-1,2-Dichloroethene	0.86	U	0.86	5.00	ug/Kg
75-34-3	1,1-Dichloroethane	0.80	U	0.80	5.00	ug/Kg
110-82-7	Cyclohexane	0.79	U	0.79	5.00	ug/Kg
78-93-3	2-Butanone	6.50	U	6.50	25.0	ug/Kg
56-23-5	Carbon Tetrachloride	0.97	U	0.97	5.00	ug/Kg
156-59-2	cis-1,2-Dichloroethene	0.75	U	0.75	5.00	ug/Kg
74-97-5	Bromochloromethane	1.20	U	1.20	5.00	ug/Kg
67-66-3	Chloroform	0.84	U	0.84	5.00	ug/Kg
71-55-6	1,1,1-Trichloroethane	0.93	U	0.93	5.00	ug/Kg
108-87-2	Methylcyclohexane	0.91	U	0.91	5.00	ug/Kg
71-43-2	Benzene	0.79	U	0.79	5.00	ug/Kg
107-06-2	1,2-Dichloroethane	0.79	U	0.79	5.00	ug/Kg
79-01-6	Trichloroethene	0.81	U	0.81	5.00	ug/Kg
78-87-5	1,2-Dichloropropane	0.91	U	0.91	5.00	ug/Kg
75-27-4	Bromodichloromethane	0.78	U	0.78	5.00	ug/Kg
108-10-1	4-Methyl-2-Pentanone	3.60	U	3.60	25.0	ug/Kg
108-88-3	Toluene	0.78	U	0.78	5.00	ug/Kg

Report of Analysis

Client:	G Environmental	Date Collected:	
Project:	Stockton	Date Received:	
Client Sample ID:	VY0411SBL01	SDG No.:	Q1761
Lab Sample ID:	VY0411SBL01	Matrix:	SOIL
Analytical Method:	SW8260	% Solid:	100
Sample Wt/Vol:	5	Units:	g
Soil Aliquot Vol:			uL
GC Column:	RXI-624	ID :	0.25
Prep Method :		Level :	LOW

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VY021851.D	1		04/11/25 11:35	VY041125

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
10061-02-6	t-1,3-Dichloropropene	0.65	U	0.65	5.00	ug/Kg
10061-01-5	cis-1,3-Dichloropropene	0.62	U	0.62	5.00	ug/Kg
79-00-5	1,1,2-Trichloroethane	0.92	U	0.92	5.00	ug/Kg
591-78-6	2-Hexanone	3.70	U	3.70	25.0	ug/Kg
124-48-1	Dibromochloromethane	0.87	U	0.87	5.00	ug/Kg
106-93-4	1,2-Dibromoethane	0.88	U	0.88	5.00	ug/Kg
127-18-4	Tetrachloroethene	1.10	U	1.10	5.00	ug/Kg
108-90-7	Chlorobenzene	0.91	U	0.91	5.00	ug/Kg
100-41-4	Ethyl Benzene	0.67	U	0.67	5.00	ug/Kg
179601-23-1	m/p-Xylenes	1.20	U	1.20	10.0	ug/Kg
95-47-6	o-Xylene	0.82	U	0.82	5.00	ug/Kg
100-42-5	Styrene	0.71	U	0.71	5.00	ug/Kg
75-25-2	Bromoform	0.86	U	0.86	5.00	ug/Kg
98-82-8	Isopropylbenzene	0.78	U	0.78	5.00	ug/Kg
79-34-5	1,1,2,2-Tetrachloroethane	1.20	U	1.20	5.00	ug/Kg
541-73-1	1,3-Dichlorobenzene	1.70	U	1.70	5.00	ug/Kg
106-46-7	1,4-Dichlorobenzene	1.60	U	1.60	5.00	ug/Kg
95-50-1	1,2-Dichlorobenzene	1.50	U	1.50	5.00	ug/Kg
96-12-8	1,2-Dibromo-3-Chloropropane	1.80	U	1.80	5.00	ug/Kg
120-82-1	1,2,4-Trichlorobenzene	3.00	U	3.00	5.00	ug/Kg
87-61-6	1,2,3-Trichlorobenzene	3.20	U	3.20	5.00	ug/Kg
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	47.0		70 (63) - 130 (155)	94%	SPK: 50
1868-53-7	Dibromofluoromethane	49.2		70 (70) - 130 (134)	98%	SPK: 50
2037-26-5	Toluene-d8	47.4		70 (74) - 130 (123)	95%	SPK: 50
460-00-4	4-Bromofluorobenzene	39.1		70 (38) - 130 (136)	78%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	321000	7.713			
540-36-3	1,4-Difluorobenzene	580000	8.615			
3114-55-4	Chlorobenzene-d5	478000	11.42			
3855-82-1	1,4-Dichlorobenzene-d4	175000	13.352			

Report of Analysis

Client:	G Environmental		Date Collected:	
Project:	Stockton		Date Received:	
Client Sample ID:	VY0411SBL01		SDG No.:	Q1761
Lab Sample ID:	VY0411SBL01		Matrix:	SOIL
Analytical Method:	SW8260		% Solid:	100
Sample Wt/Vol:	5	Units: g	Final Vol:	5000 uL
Soil Aliquot Vol:		uL	Test:	VOC-TCLVOA-10
GC Column:	RXI-624	ID : 0.25	Level :	LOW
Prep Method :				

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VY021851.D	1		04/11/25 11:35	VY041125

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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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

Report of Analysis

Client:	G Environmental	Date Collected:	
Project:	Stockton	Date Received:	
Client Sample ID:	VY0411SBS01	SDG No.:	Q1761
Lab Sample ID:	VY0411SBS01	Matrix:	SOIL
Analytical Method:	SW8260	% Solid:	100
Sample Wt/Vol:	5	Units:	g
Soil Aliquot Vol:			uL
GC Column:	RXI-624	ID :	0.25
Prep Method :		Level :	LOW

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VY021852.D	1		04/11/25 12:06	VY041125

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
75-71-8	Dichlorodifluoromethane	17.7		1.10	5.00	ug/Kg
74-87-3	Chloromethane	17.5		1.10	5.00	ug/Kg
75-01-4	Vinyl Chloride	17.2		0.79	5.00	ug/Kg
74-83-9	Bromomethane	18.9		1.10	5.00	ug/Kg
75-00-3	Chloroethane	18.0		1.30	5.00	ug/Kg
75-69-4	Trichlorofluoromethane	18.6		1.20	5.00	ug/Kg
76-13-1	1,1,2-Trichlorotrifluoroethane	19.6		1.10	5.00	ug/Kg
75-35-4	1,1-Dichloroethene	19.1		1.00	5.00	ug/Kg
67-64-1	Acetone	100		4.70	25.0	ug/Kg
75-15-0	Carbon Disulfide	17.4		1.10	5.00	ug/Kg
1634-04-4	Methyl tert-butyl Ether	19.8		0.73	5.00	ug/Kg
79-20-9	Methyl Acetate	20.7		1.50	5.00	ug/Kg
75-09-2	Methylene Chloride	20.1		3.50	10.0	ug/Kg
156-60-5	trans-1,2-Dichloroethene	18.9		0.86	5.00	ug/Kg
75-34-3	1,1-Dichloroethane	19.7		0.80	5.00	ug/Kg
110-82-7	Cyclohexane	17.9		0.79	5.00	ug/Kg
78-93-3	2-Butanone	95.3		6.50	25.0	ug/Kg
56-23-5	Carbon Tetrachloride	19.7		0.97	5.00	ug/Kg
156-59-2	cis-1,2-Dichloroethene	19.9		0.75	5.00	ug/Kg
74-97-5	Bromochloromethane	19.7		1.20	5.00	ug/Kg
67-66-3	Chloroform	19.7		0.84	5.00	ug/Kg
71-55-6	1,1,1-Trichloroethane	19.6		0.93	5.00	ug/Kg
108-87-2	Methylcyclohexane	18.8		0.91	5.00	ug/Kg
71-43-2	Benzene	19.7		0.79	5.00	ug/Kg
107-06-2	1,2-Dichloroethane	19.8		0.79	5.00	ug/Kg
79-01-6	Trichloroethene	19.8		0.81	5.00	ug/Kg
78-87-5	1,2-Dichloropropane	19.5		0.91	5.00	ug/Kg
75-27-4	Bromodichloromethane	20.0		0.78	5.00	ug/Kg
108-10-1	4-Methyl-2-Pentanone	97.2		3.60	25.0	ug/Kg
108-88-3	Toluene	20.1		0.78	5.00	ug/Kg

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
10061-02-6	t-1,3-Dichloropropene	19.4		0.65	5.00	ug/Kg
10061-01-5	cis-1,3-Dichloropropene	19.9		0.62	5.00	ug/Kg
79-00-5	1,1,2-Trichloroethane	20.1		0.92	5.00	ug/Kg
591-78-6	2-Hexanone	97.2		3.70	25.0	ug/Kg
124-48-1	Dibromochloromethane	20.3		0.87	5.00	ug/Kg
106-93-4	1,2-Dibromoethane	20.4		0.88	5.00	ug/Kg
127-18-4	Tetrachloroethene	22.1		1.10	5.00	ug/Kg
108-90-7	Chlorobenzene	19.8		0.91	5.00	ug/Kg
100-41-4	Ethyl Benzene	19.5		0.67	5.00	ug/Kg
179601-23-1	m/p-Xylenes	39.5		1.20	10.0	ug/Kg
95-47-6	o-Xylene	20.0		0.82	5.00	ug/Kg
100-42-5	Styrene	19.8		0.71	5.00	ug/Kg
75-25-2	Bromoform	20.4		0.86	5.00	ug/Kg
98-82-8	Isopropylbenzene	19.0		0.78	5.00	ug/Kg
79-34-5	1,1,2,2-Tetrachloroethane	17.9		1.20	5.00	ug/Kg
541-73-1	1,3-Dichlorobenzene	19.4		1.70	5.00	ug/Kg
106-46-7	1,4-Dichlorobenzene	19.5		1.60	5.00	ug/Kg
95-50-1	1,2-Dichlorobenzene	20.1		1.50	5.00	ug/Kg
96-12-8	1,2-Dibromo-3-Chloropropane	20.9		1.80	5.00	ug/Kg
120-82-1	1,2,4-Trichlorobenzene	21.1		3.00	5.00	ug/Kg
87-61-6	1,2,3-Trichlorobenzene	21.4		3.20	5.00	ug/Kg
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	52.3		70 (63) - 130 (155)	105%	SPK: 50
1868-53-7	Dibromofluoromethane	53.2		70 (70) - 130 (134)	106%	SPK: 50
2037-26-5	Toluene-d8	53.8		70 (74) - 130 (123)	108%	SPK: 50
460-00-4	4-Bromofluorobenzene	55.3		70 (38) - 130 (136)	111%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	262000	7.707			
540-36-3	1,4-Difluorobenzene	412000	8.616			
3114-55-4	Chlorobenzene-d5	368000	11.42			
3855-82-1	1,4-Dichlorobenzene-d4	190000	13.347			

Report of Analysis

Client:	G Environmental		Date Collected:	
Project:	Stockton		Date Received:	
Client Sample ID:	VY0411SBS01		SDG No.:	Q1761
Lab Sample ID:	VY0411SBS01		Matrix:	SOIL
Analytical Method:	SW8260		% Solid:	100
Sample Wt/Vol:	5	Units: g	Final Vol:	5000 uL
Soil Aliquot Vol:		uL	Test:	VOC-TCLVOA-10
GC Column:	RXI-624	ID : 0.25	Level :	LOW
Prep Method :				

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VY021852.D	1		04/11/25 12:06	VY041125

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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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

Report of Analysis

Client:	G Environmental	Date Collected:	
Project:	Stockton	Date Received:	
Client Sample ID:	VY0411SBSD01	SDG No.:	Q1761
Lab Sample ID:	VY0411SBSD01	Matrix:	SOIL
Analytical Method:	SW8260	% Solid:	100
Sample Wt/Vol:	5 Units: g	Final Vol:	5000 uL
Soil Aliquot Vol:	uL	Test:	VOC-TCLVOA-10
GC Column:	RXI-624 ID : 0.25	Level :	LOW
Prep Method :			

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VY021853.D	1		04/11/25 12:28	VY041125

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
75-71-8	Dichlorodifluoromethane	18.3		1.10	5.00	ug/Kg
74-87-3	Chloromethane	16.4		1.10	5.00	ug/Kg
75-01-4	Vinyl Chloride	17.0		0.79	5.00	ug/Kg
74-83-9	Bromomethane	17.5		1.10	5.00	ug/Kg
75-00-3	Chloroethane	17.1		1.30	5.00	ug/Kg
75-69-4	Trichlorofluoromethane	19.0		1.20	5.00	ug/Kg
76-13-1	1,1,2-Trichlorotrifluoroethane	19.5		1.10	5.00	ug/Kg
75-35-4	1,1-Dichloroethene	19.8		1.00	5.00	ug/Kg
67-64-1	Acetone	84.8		4.70	25.0	ug/Kg
75-15-0	Carbon Disulfide	17.9		1.10	5.00	ug/Kg
1634-04-4	Methyl tert-butyl Ether	19.6		0.73	5.00	ug/Kg
79-20-9	Methyl Acetate	21.8		1.50	5.00	ug/Kg
75-09-2	Methylene Chloride	20.4		3.50	10.0	ug/Kg
156-60-5	trans-1,2-Dichloroethene	19.5		0.86	5.00	ug/Kg
75-34-3	1,1-Dichloroethane	19.9		0.80	5.00	ug/Kg
110-82-7	Cyclohexane	18.2		0.79	5.00	ug/Kg
78-93-3	2-Butanone	90.6		6.50	25.0	ug/Kg
56-23-5	Carbon Tetrachloride	20.2		0.97	5.00	ug/Kg
156-59-2	cis-1,2-Dichloroethene	20.2		0.75	5.00	ug/Kg
74-97-5	Bromochloromethane	18.0		1.20	5.00	ug/Kg
67-66-3	Chloroform	20.0		0.84	5.00	ug/Kg
71-55-6	1,1,1-Trichloroethane	20.1		0.93	5.00	ug/Kg
108-87-2	Methylcyclohexane	19.1		0.91	5.00	ug/Kg
71-43-2	Benzene	20.3		0.79	5.00	ug/Kg
107-06-2	1,2-Dichloroethane	19.6		0.79	5.00	ug/Kg
79-01-6	Trichloroethene	20.9		0.81	5.00	ug/Kg
78-87-5	1,2-Dichloropropane	19.9		0.91	5.00	ug/Kg
75-27-4	Bromodichloromethane	20.0		0.78	5.00	ug/Kg
108-10-1	4-Methyl-2-Pentanone	94.4		3.60	25.0	ug/Kg
108-88-3	Toluene	20.5		0.78	5.00	ug/Kg

Report of Analysis

Client:	G Environmental	Date Collected:	
Project:	Stockton	Date Received:	
Client Sample ID:	VY0411SBSD01	SDG No.:	Q1761
Lab Sample ID:	VY0411SBSD01	Matrix:	SOIL
Analytical Method:	SW8260	% Solid:	100
Sample Wt/Vol:	5	Units:	g
Soil Aliquot Vol:			uL
GC Column:	RXI-624	ID :	0.25
Prep Method :		Level :	LOW

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VY021853.D	1		04/11/25 12:28	VY041125

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
10061-02-6	t-1,3-Dichloropropene	19.9		0.65	5.00	ug/Kg
10061-01-5	cis-1,3-Dichloropropene	19.6		0.62	5.00	ug/Kg
79-00-5	1,1,2-Trichloroethane	20.6		0.92	5.00	ug/Kg
591-78-6	2-Hexanone	93.4		3.70	25.0	ug/Kg
124-48-1	Dibromochloromethane	20.5		0.87	5.00	ug/Kg
106-93-4	1,2-Dibromoethane	19.9		0.88	5.00	ug/Kg
127-18-4	Tetrachloroethene	22.4		1.10	5.00	ug/Kg
108-90-7	Chlorobenzene	20.4		0.91	5.00	ug/Kg
100-41-4	Ethyl Benzene	20.1		0.67	5.00	ug/Kg
179601-23-1	m/p-Xylenes	40.8		1.20	10.0	ug/Kg
95-47-6	o-Xylene	20.5		0.82	5.00	ug/Kg
100-42-5	Styrene	20.6		0.71	5.00	ug/Kg
75-25-2	Bromoform	20.4		0.86	5.00	ug/Kg
98-82-8	Isopropylbenzene	19.7		0.78	5.00	ug/Kg
79-34-5	1,1,2,2-Tetrachloroethane	18.0		1.20	5.00	ug/Kg
541-73-1	1,3-Dichlorobenzene	20.3		1.70	5.00	ug/Kg
106-46-7	1,4-Dichlorobenzene	20.2		1.60	5.00	ug/Kg
95-50-1	1,2-Dichlorobenzene	20.1		1.50	5.00	ug/Kg
96-12-8	1,2-Dibromo-3-Chloropropane	19.9		1.80	5.00	ug/Kg
120-82-1	1,2,4-Trichlorobenzene	21.9		3.00	5.00	ug/Kg
87-61-6	1,2,3-Trichlorobenzene	22.5		3.20	5.00	ug/Kg
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	50.6		70 (63) - 130 (155)	101%	SPK: 50
1868-53-7	Dibromofluoromethane	53.9		70 (70) - 130 (134)	108%	SPK: 50
2037-26-5	Toluene-d8	53.1		70 (74) - 130 (123)	106%	SPK: 50
460-00-4	4-Bromofluorobenzene	54.5		70 (38) - 130 (136)	109%	SPK: 50
INTERNAL STANDARDS						
363-72-4	Pentafluorobenzene	262000	7.707			
540-36-3	1,4-Difluorobenzene	410000	8.616			
3114-55-4	Chlorobenzene-d5	363000	11.414			
3855-82-1	1,4-Dichlorobenzene-d4	188000	13.347			

Report of Analysis

Client:	G Environmental		Date Collected:	
Project:	Stockton		Date Received:	
Client Sample ID:	VY0411SBSD01		SDG No.:	Q1761
Lab Sample ID:	VY0411SBSD01		Matrix:	SOIL
Analytical Method:	SW8260		% Solid:	100
Sample Wt/Vol:	5	Units: g	Final Vol:	5000 uL
Soil Aliquot Vol:		uL	Test:	VOC-TCLVOA-10
GC Column:	RXI-624	ID : 0.25	Level :	LOW
Prep Method :				

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VY021853.D	1		04/11/25 12:28	VY041125

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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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



CALIBRATION SUMMARY

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: GENV01

Lab Code: CHEM Case No.: Q1761 SAS No.: Q1761 SDG No.: Q1761

Instrument ID: MSVOA_Y Calibration Date(s): 03/27/2025 03/27/2025

Heated Purge: (Y/N) Y Calibration Time(s): 10:08 12:02

GC Column: RXI-624 ID: 0.25 (mm)

LAB FILE ID:		RRF005 = VY021656.D		RRF010 = VY021657.D		RRF020 = VY021658.D		RRF050 = VY021659.D		RRF100 = VY021660.D		RRF150 = VY021661.D	
COMPOUND		RRF005	RRF010	RRF020	RRF050	RRF100	RRF150	RRF	% RSD				
Dichlorodifluoromethane		0.602	0.577	0.493	0.506	0.485	0.447	0.518	11.4				
Chloromethane		0.876	0.775	0.713	0.692	0.672	0.644	0.729	11.6				
Vinyl Chloride		0.977	0.886	0.810	0.800	0.758	0.714	0.824	11.5				
Bromomethane		0.709	0.613	0.537	0.513	0.494	0.506	0.562	14.9				
Chloroethane		0.644	0.570	0.544	0.512	0.490	0.476	0.539	11.5				
Trichlorofluoromethane		1.212	1.124	1.104	1.013	0.967	0.914	1.056	10.5				
1,1,2-Trichlorotrifluoroethane		0.668	0.601	0.553	0.536	0.510	0.473	0.557	12.4				
1,1-Dichloroethene		0.589	0.564	0.508	0.514	0.496	0.477	0.524	8.2				
Acetone		0.145	0.123	0.102	0.107	0.087	0.102	0.111	18.3				
Carbon Disulfide		1.956	1.828	1.709	1.696	1.623	1.552	1.727	8.4				
Methyl tert-butyl Ether		1.380	1.335	1.247	1.325	1.256	1.292	1.306	3.9				
Methyl Acetate		0.329	0.286	0.269	0.265	0.243	0.257	0.275	11				
Methylene Chloride		0.870	0.694	0.594	0.564	0.518	0.514	0.626	21.8				
trans-1,2-Dichloroethene		0.634	0.608	0.568	0.567	0.551	0.536	0.577	6.3				
1,1-Dichloroethane		1.178	1.091	1.024	1.040	0.992	0.968	1.049	7.3				
Cyclohexane		1.170	1.032	0.912	0.918	0.886	0.818	0.956	13.1				
2-Butanone		0.181	0.166	0.149	0.159	0.140	0.155	0.158	8.9				
Carbon Tetrachloride		0.637	0.594	0.546	0.564	0.547	0.507	0.566	7.9				
cis-1,2-Dichloroethene		0.700	0.658	0.626	0.652	0.631	0.629	0.649	4.3				
Bromochloromethane		0.513	0.470	0.424	0.426	0.420	0.402	0.442	9.3				
Chloroform		1.235	1.141	1.062	1.081	1.026	1.000	1.091	7.9				
1,1,1-Trichloroethane		1.155	1.013	0.959	0.967	0.929	0.893	0.986	9.3				
Methylcyclohexane		0.593	0.599	0.569	0.619	0.612	0.550	0.590	4.5				
Benzene		1.606	1.527	1.429	1.499	1.445	1.371	1.479	5.6				
1,2-Dichloroethane		0.467	0.445	0.399	0.420	0.392	0.379	0.417	8.1				
Trichloroethene		0.412	0.398	0.362	0.374	0.366	0.348	0.377	6.4				
1,2-Dichloropropane		0.383	0.366	0.334	0.353	0.340	0.325	0.350	6.2				
Bromodichloromethane		0.571	0.546	0.504	0.525	0.508	0.484	0.523	6				
4-Methyl-2-Pentanone		0.232	0.240	0.219	0.241	0.222	0.226	0.230	3.9				
Toluene		0.940	0.943	0.893	0.951	0.928	0.885	0.923	3				

* 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.

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name:	<u>CHEMTECH</u>	Contract:	<u>GENV01</u>
Lab Code:	<u>CHEM</u>	Case No.:	<u>Q1761</u>
Instrument ID:	<u>MSVOA_Y</u>	SAS No.:	<u>Q1761</u>
Heated Purge: (Y/N)	<u>Y</u>	SDG No.:	<u>Q1761</u>
GC Column:	<u>RXI-624</u>	Calibration Date(s):	<u>03/27/2025</u>
ID:	<u>0.25 (mm)</u>	Calibration Time(s):	<u>10:08</u>
			<u>12:02</u>

LAB FILE ID:		RRF005 = VY021656.D		RRF010 = VY021657.D		RRF020 = VY021658.D		
		RRF050 = VY021659.D		RRF100 = VY021660.D		RRF150 = VY021661.D		
COMPOUND	RRF005	RRF010	RRF020	RRF050	RRF100	RRF150	RRF	% RSD
t-1,3-Dichloropropene	0.456	0.470	0.439	0.486	0.475	0.461	0.465	3.5
cis-1,3-Dichloropropene	0.553	0.558	0.528	0.565	0.549	0.534	0.548	2.6
1,1,2-Trichloroethane	0.280	0.283	0.249	0.261	0.247	0.240	0.260	6.9
2-Hexanone	0.154	0.154	0.144	0.164	0.148	0.154	0.153	4.2
Dibromochloromethane	0.371	0.367	0.340	0.364	0.348	0.337	0.355	4.1
1,2-Dibromoethane	0.267	0.250	0.230	0.254	0.233	0.229	0.244	6.3
Tetrachloroethene	0.534	0.503	0.453	0.449	0.412	0.382	0.455	12.3
Chlorobenzene	1.244	1.178	1.110	1.155	1.116	1.073	1.146	5.3
Ethyl Benzene	1.949	1.916	1.918	2.051	2.023	1.935	1.965	2.9
m/p-Xylenes	0.726	0.753	0.744	0.787	0.774	0.731	0.753	3.2
o-Xylene	0.665	0.670	0.676	0.738	0.725	0.694	0.695	4.4
Styrene	1.064	1.110	1.134	1.232	1.229	1.181	1.158	5.8
Bromoform	0.241	0.239	0.229	0.237	0.225	0.221	0.232	3.4
Isopropylbenzene	3.628	3.701	3.582	3.826	3.843	3.702	3.714	2.8
1,1,2,2-Tetrachloroethane	0.764	0.720	0.616	0.657	0.619	0.631	0.668	9.1
1,3-Dichlorobenzene	1.895	1.774	1.707	1.748	1.728	1.689	1.757	4.2
1,4-Dichlorobenzene	1.876	1.794	1.704	1.741	1.673	1.626	1.736	5.2
1,2-Dichlorobenzene	1.598	1.607	1.496	1.524	1.486	1.465	1.529	3.9
1,2-Dibromo-3-Chloropropane	0.112	0.111	0.097	0.102	0.096	0.102	0.103	6.6
1,2,4-Trichlorobenzene	0.771	0.781	0.787	0.878	0.870	0.889	0.829	6.6
1,2,3-Trichlorobenzene	0.653	0.681	0.673	0.746	0.737	0.753	0.707	6.1
1,2-Dichloroethane-d4	0.612	0.559	0.502	0.548	0.510	0.520	0.542	7.5
Dibromofluoromethane	0.357	0.321	0.290	0.339	0.325	0.313	0.324	7
Toluene-d8	1.277	1.240	1.127	1.306	1.252	1.201	1.234	5.1
4-Bromofluorobenzene	0.447	0.410	0.378	0.439	0.424	0.407	0.417	6

* 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.

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: GENV01

Lab Code: CHEM Case No.: Q1761 SAS No.: Q1761 SDG No.: Q1761

Instrument ID: MSVOA_Y Calibration Date/Time: 04/11/2025 09:18

Lab File ID: VY021850.D Init. Calib. Date(s): 03/27/2025 03/27/2025

Heated Purge: (Y/N) Y Init. Calib. Time(s): 10:08 12:02

GC Column: RXI-624 ID: 0.25 (mm)

COMPOUND	RRF	RRF050	MIN RRF	%D	MAX%D
Dichlorodifluoromethane	0.518	0.500		-3.47	20
Chloromethane	0.729	0.591	0.1	-18.93	20
Vinyl Chloride	0.824	0.724		-12.14	20
Bromomethane	0.562	0.463		-17.62	20
Chloroethane	0.539	0.483		-10.39	20
Trichlorofluoromethane	1.056	1.051		-0.47	20
1,1,2-Trichlorotrifluoroethane	0.557	0.581		4.31	20
1,1-Dichloroethene	0.524	0.541		3.24	20
Acetone	0.111	0.094		-15.31	20
Carbon Disulfide	1.727	1.635		-5.33	20
Methyl tert-butyl Ether	1.306	1.320		1.07	20
Methyl Acetate	0.275	0.290		5.45	20
Methylene Chloride	0.626	0.581		-7.19	20
trans-1,2-Dichloroethene	0.577	0.594		2.95	20
1,1-Dichloroethane	1.049	1.065	0.1	1.52	20
Cyclohexane	0.956	0.919		-3.87	20
2-Butanone	0.158	0.139		-12.02	20
Carbon Tetrachloride	0.566	0.635		12.19	20
cis-1,2-Dichloroethene	0.649	0.680		4.78	20
Bromochloromethane	0.442	0.405		-8.37	20
Chloroform	1.091	1.112		1.92	20
1,1,1-Trichloroethane	0.986	1.021		3.55	20
Methylcyclohexane	0.590	0.661		12.03	20
Benzene	1.479	1.585		7.17	20
1,2-Dichloroethane	0.417	0.424		1.68	20
Trichloroethene	0.377	0.432		14.59	20
1,2-Dichloropropane	0.350	0.373		6.57	20
Bromodichloromethane	0.523	0.552		5.55	20
4-Methyl-2-Pentanone	0.230	0.221		-3.91	20
Toluene	0.923	1.019		10.4	20
t-1,3-Dichloropropene	0.465	0.489		5.16	20
cis-1,3-Dichloropropene	0.548	0.585		6.75	20
1,1,2-Trichloroethane	0.260	0.270		3.85	20
2-Hexanone	0.153	0.146		-4.57	20
Dibromochloromethane	0.355	0.384		8.17	20
1,2-Dibromoethane	0.244	0.260		6.56	20
Tetrachloroethene	0.455	0.571		25.5	20
Chlorobenzene	1.146	1.268	0.3	10.65	20

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

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: GENV01

Lab Code: CHEM Case No.: Q1761 SAS No.: Q1761 SDG No.: Q1761

Instrument ID: MSVOA_Y Calibration Date/Time: 04/11/2025 09:18

Lab File ID: VY021850.D Init. Calib. Date(s): 03/27/2025 03/27/2025

Heated Purge: (Y/N) Y Init. Calib. Time(s): 10:08 12:02

GC Column: RXI-624 ID: 0.25 (mm)

COMPOUND	RRF	RRF050	MIN RRF	%D	MAX%D
Ethyl Benzene	1.965	2.263		15.16	20
m/p-Xylenes	0.753	0.863		14.61	20
o-Xylene	0.695	0.813		16.98	20
Styrene	1.158	1.347		16.32	20
Bromoform	0.232	0.254	0.1	9.48	20
Isopropylbenzene	3.714	4.282		15.29	20
1,1,2,2-Tetrachloroethane	0.668	0.611	0.3	-8.53	20
1,3-Dichlorobenzene	1.757	1.945		10.7	20
1,4-Dichlorobenzene	1.736	1.895		9.16	20
1,2-Dichlorobenzene	1.529	1.699		11.12	20
1,2-Dibromo-3-Chloropropane	0.103	0.103		0	20
1,2,4-Trichlorobenzene	0.829	1.048		26.42	20
1,2,3-Trichlorobenzene	0.707	0.882		24.75	20
1,2-Dichloroethane-d4	0.542	0.511		-5.72	20
Dibromofluoromethane	0.324	0.349		7.72	20
Toluene-d8	1.234	1.339		8.51	20
4-Bromofluorobenzene	0.417	0.461		10.55	20

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



SAMPLE RAW DATA

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
 Data File : VY021855.D
 Acq On : 11 Apr 2025 13:33
 Operator : SY/MD
 Sample : Q1761-01
 Misc : 7.82g/5.0mL/MSVOA_Y/SOIL/A
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 GST3

A
B
C
D
E
F
G
H
I
J

Quant Time: Apr 12 02:48:30 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M
 Quant Title : SW846 8260
 QLast Update : Fri Mar 28 02:30:29 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	7.713	168	317324	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.616	114	560087	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.420	117	469329	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.346	152	180059	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.061	65	160307	46.626	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	=	93.260%
35) Dibromofluoromethane	7.634	113	180049	49.557	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	=	99.120%
50) Toluene-d8	10.109	98	662731	47.948	ug/l	0.00
Spiked Amount	50.000	Range	58 - 134	Recovery	=	95.900%
62) 4-Bromofluorobenzene	12.408	95	189531	40.540	ug/l	0.00
Spiked Amount	50.000	Range	30 - 143	Recovery	=	81.080%

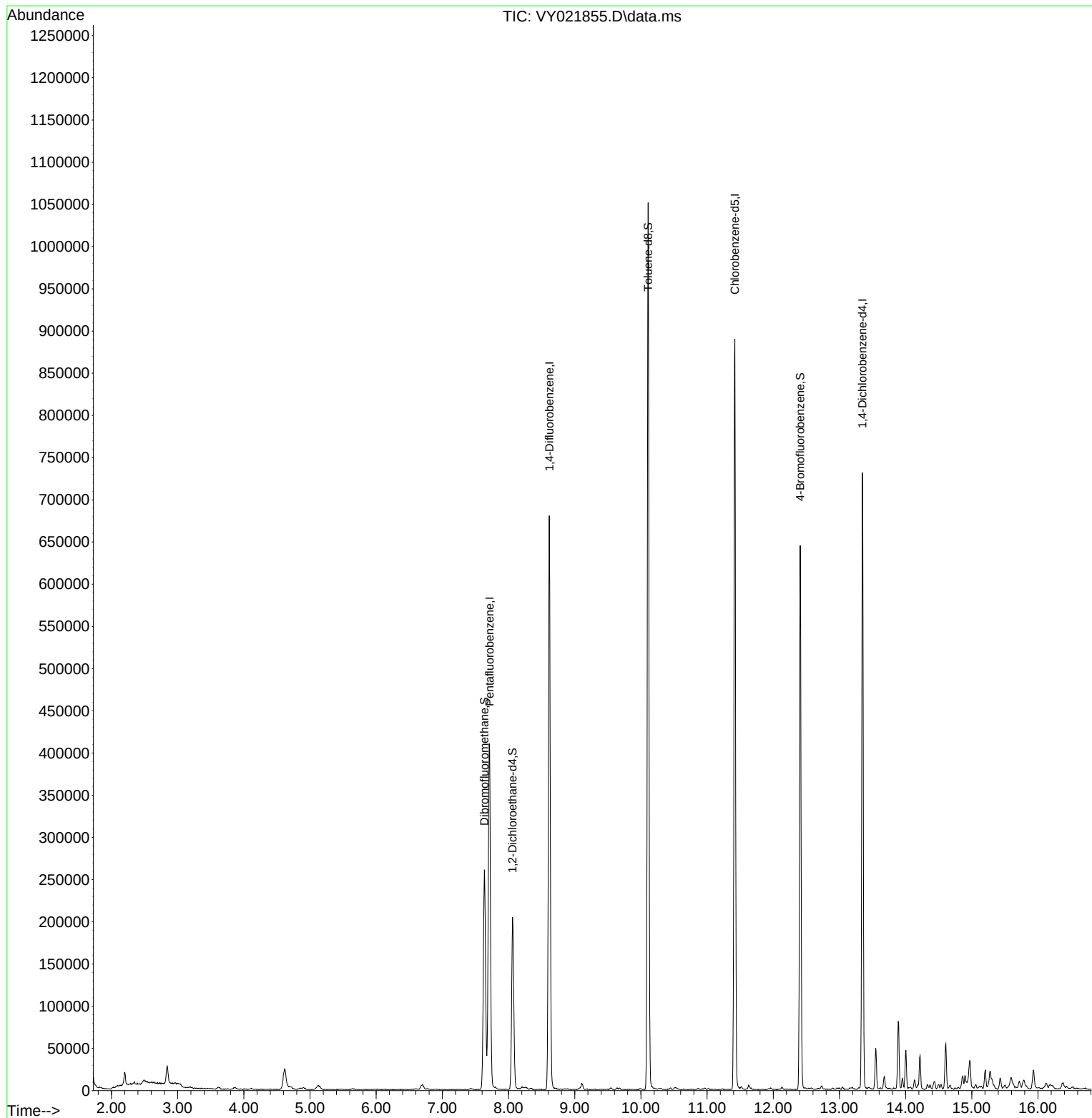
Target Compounds	Qvalue

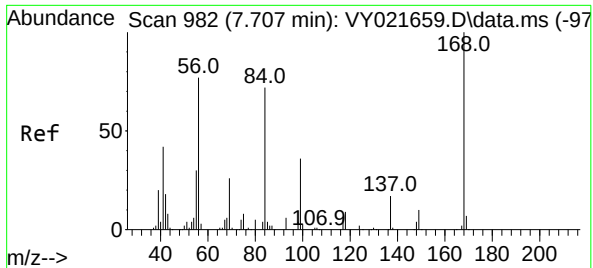
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
Data File : VY021855.D
Acq On : 11 Apr 2025 13:33
Operator : SY/MD
Sample : Q1761-01
Misc : 7.82g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
GST3

Quant Time: Apr 12 02:48:30 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M
Quant Title : SW846 8260
QLast Update : Fri Mar 28 02:30:29 2025
Response via : Initial Calibration

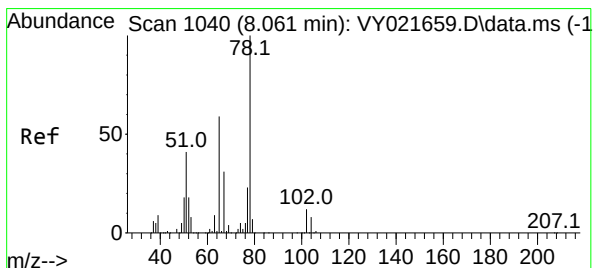
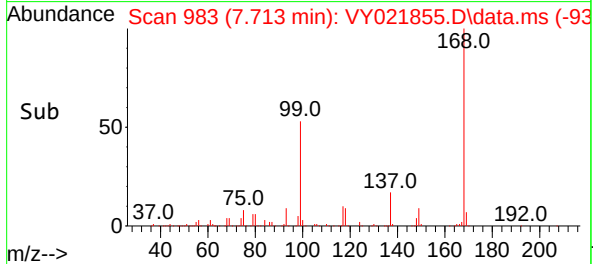
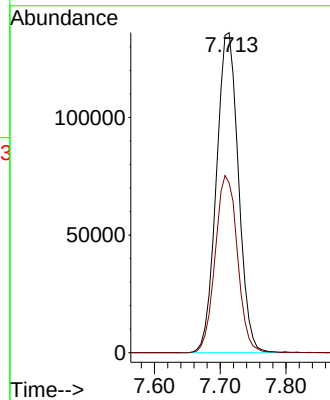
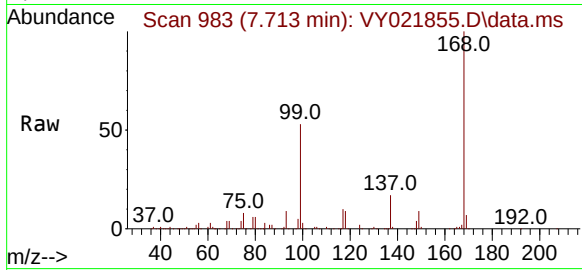




#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 7.713 min Scan# 99
Delta R.T. 0.006 min
Lab File: VY021855.D
Acq: 11 Apr 2025 13:33

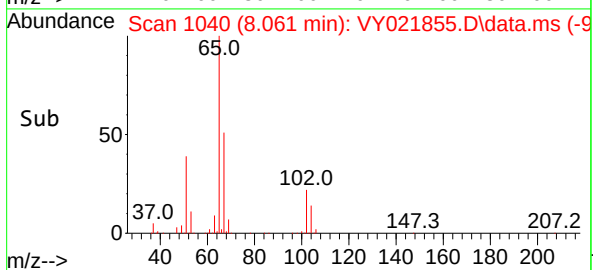
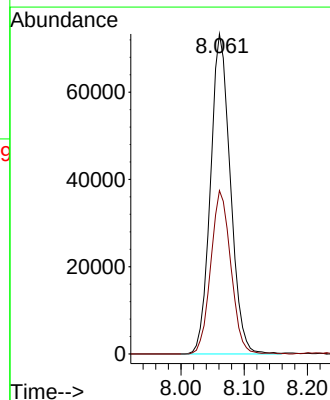
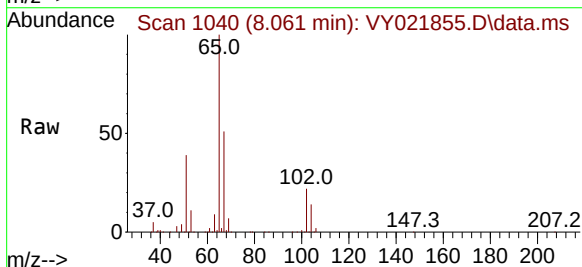
Instrument :
MSVOA_Y
ClientSampleId :
GST3

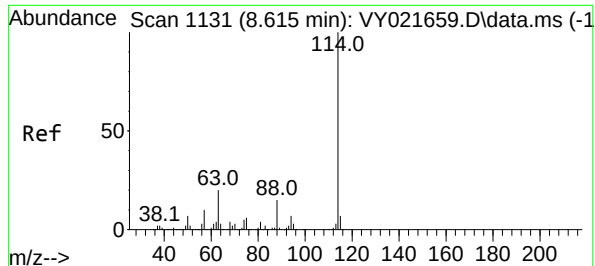
Tgt Ion:168 Resp: 317324
Ion Ratio Lower Upper
168 100
99 53.1 46.7 70.1



#33
1,2-Dichloroethane-d4
Concen: 46.626 ug/l
RT: 8.061 min Scan# 1040
Delta R.T. 0.000 min
Lab File: VY021855.D
Acq: 11 Apr 2025 13:33

Tgt Ion: 65 Resp: 160307
Ion Ratio Lower Upper
65 100
67 51.4 0.0 104.6





#34

1,4-Difluorobenzene

Concen: 50.000 ug/l

RT: 8.616 min Scan# 1131

Delta R.T. 0.000 min

Lab File: VY021855.D

Acq: 11 Apr 2025 13:33

Instrument :

MSVOA_Y

ClientSampleId :

GST3

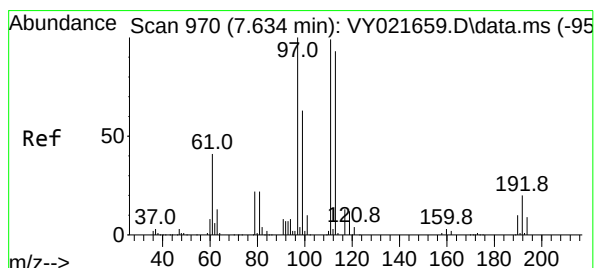
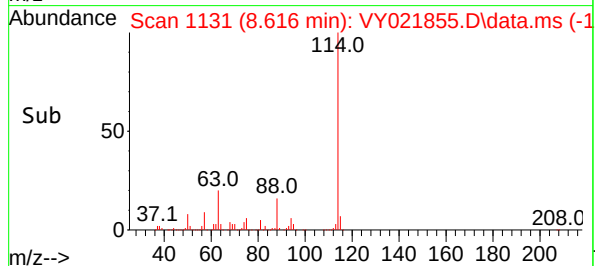
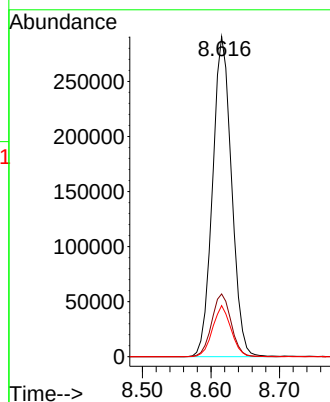
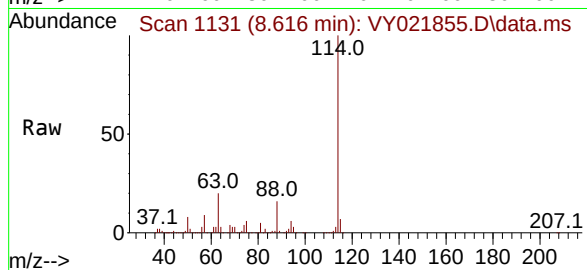
Tgt Ion:114 Resp: 560087

Ion Ratio Lower Upper

114 100

63 19.6 0.0 40.2

88 16.0 0.0 29.8



#35

Dibromofluoromethane

Concen: 49.557 ug/l

RT: 7.634 min Scan# 970

Delta R.T. 0.000 min

Lab File: VY021855.D

Acq: 11 Apr 2025 13:33

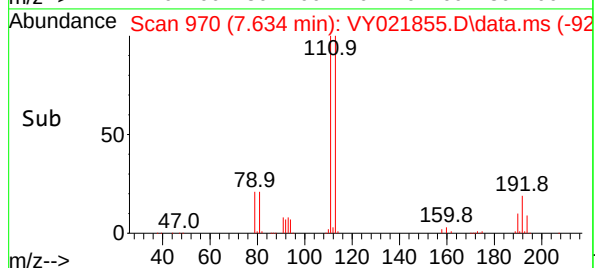
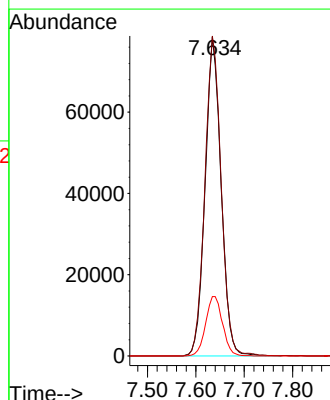
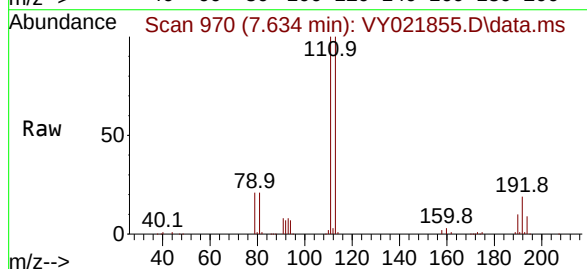
Tgt Ion:113 Resp: 180049

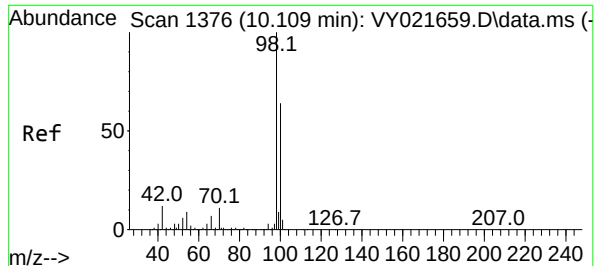
Ion Ratio Lower Upper

113 100

111 102.4 82.6 123.8

192 19.7 16.2 24.4





#50

Toluene-d8

Concen: 47.948 ug/l

RT: 10.109 min Scan# 11

Delta R.T. 0.000 min

Lab File: VY021855.D

Acq: 11 Apr 2025 13:33

Instrument :

MSVOA_Y

ClientSampleId :

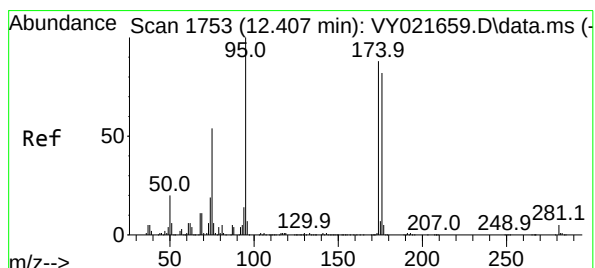
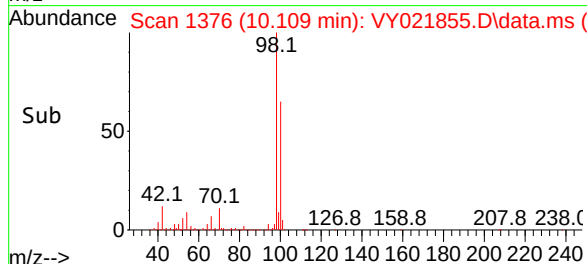
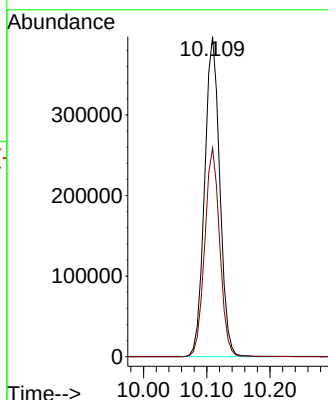
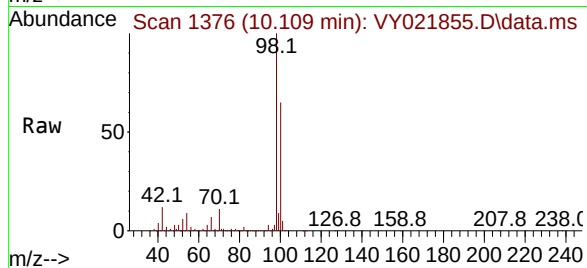
GST3

Tgt Ion: 98 Resp: 662731

Ion Ratio Lower Upper

98 100

100 64.5 52.1 78.1



#62

4-Bromofluorobenzene

Concen: 40.540 ug/l

RT: 12.408 min Scan# 1753

Delta R.T. 0.000 min

Lab File: VY021855.D

Acq: 11 Apr 2025 13:33

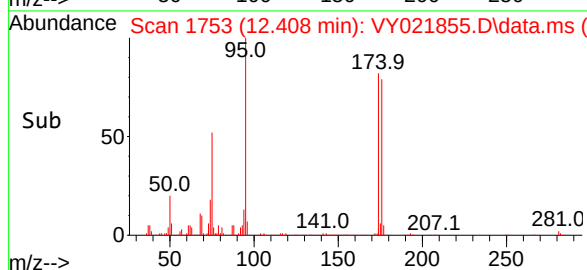
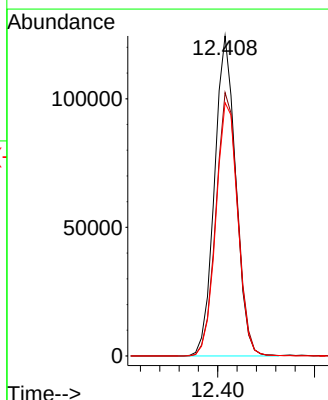
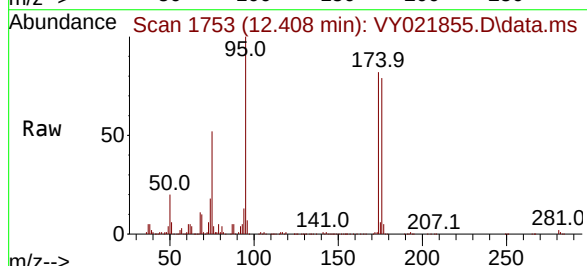
Tgt Ion: 95 Resp: 189531

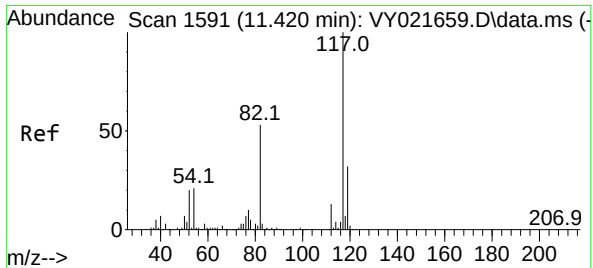
Ion Ratio Lower Upper

95 100

174 84.6 0.0 170.8

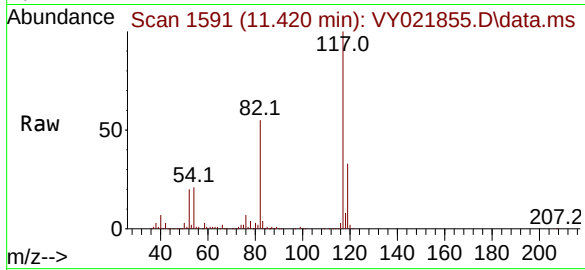
176 81.5 0.0 162.0



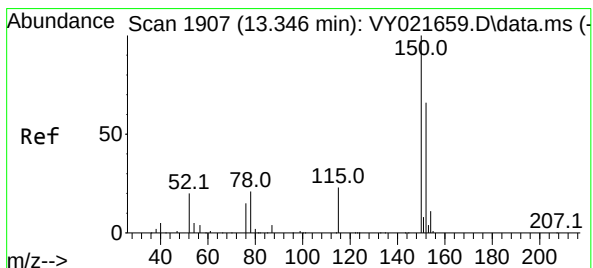
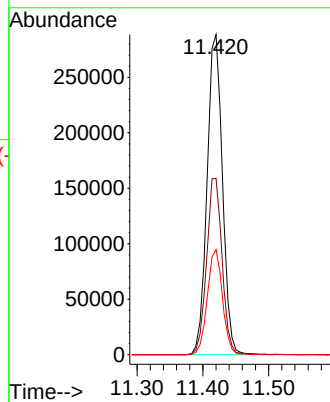
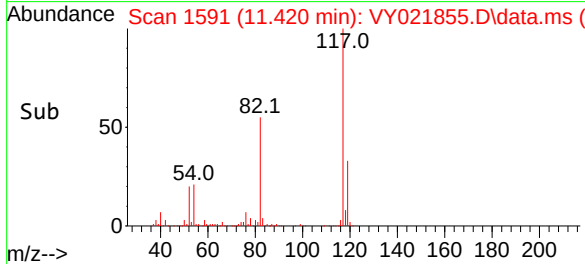


#63
Chlorobenzene-d5
Concen: 50.000 ug/l
RT: 11.420 min Scan# 117
Delta R.T. 0.000 min
Lab File: VY021855.D
Acq: 11 Apr 2025 13:33

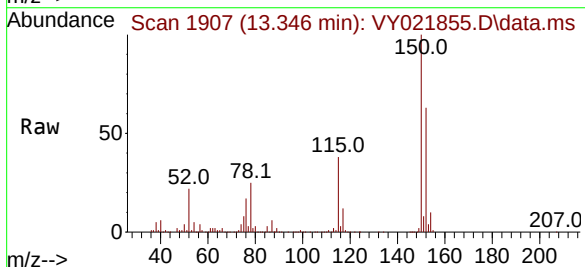
Instrument :
MSVOA_Y
ClientSampleId :
GST3



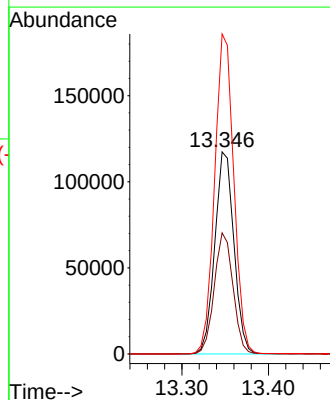
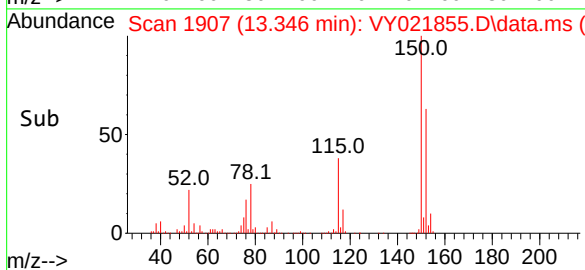
Tgt Ion:117 Resp: 469329
Ion Ratio Lower Upper
117 100
82 55.1 42.8 64.2
119 32.9 25.7 38.5



#72
1,4-Dichlorobenzene-d4
Concen: 50.000 ug/l
RT: 13.346 min Scan# 1907
Delta R.T. 0.000 min
Lab File: VY021855.D
Acq: 11 Apr 2025 13:33



Tgt Ion:152 Resp: 180059
Ion Ratio Lower Upper
152 100
115 59.0 28.8 86.5
150 158.4 0.0 348.4



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
Data File : VY021855.D
Acq On : 11 Apr 2025 13:33
Operator : SY/MD
Sample : Q1761-01
Misc : 7.82g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
GST3

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M

Title : SW846 8260

Signal : TIC: VY021855.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.196	70	78	84	rBV	15916	30434	1.72%	0.312%
2	2.842	178	184	191	rBV2	20153	40648	2.29%	0.417%
3	4.616	464	475	488	rBV3	23977	82478	4.65%	0.846%
4	5.122	546	558	572	rVB8	5206	20613	1.16%	0.212%
5	7.634	959	970	976	rBV	260364	606024	34.18%	6.218%
6	7.707	976	982	996	rVB	408015	951933	53.69%	9.768%
7	8.061	1030	1040	1053	rBV	203880	448622	25.30%	4.603%
8	8.616	1120	1131	1145	rBV	680272	1322317	74.59%	13.568%
9	9.103	1201	1211	1219	rVB4	7375	17868	1.01%	0.183%
10	10.109	1368	1376	1390	rBV	1050917	1772888	100.00%	18.191%
11	11.420	1583	1591	1602	rBV	888800	1463214	82.53%	15.014%
12	12.408	1745	1753	1762	rBV	644802	1023946	57.76%	10.507%
13	13.346	1900	1907	1915	rBV	729312	1106480	62.41%	11.353%
14	13.548	1934	1940	1947	rBV	48391	76333	4.31%	0.783%
15	13.676	1955	1961	1971	rBV4	14834	26925	1.52%	0.276%
16	13.889	1990	1996	2002	rBV2	79973	132922	7.50%	1.364%
17	13.950	2002	2006	2010	rVV3	11936	18677	1.05%	0.192%
18	14.005	2010	2015	2020	rVB	45156	67305	3.80%	0.691%
19	14.133	2030	2036	2041	rBV3	10893	20011	1.13%	0.205%
20	14.218	2045	2050	2056	rVB	39195	59698	3.37%	0.613%
21	14.608	2107	2114	2119	rBV	53279	87991	4.96%	0.903%
22	14.864	2149	2156	2159	rBV3	14640	28433	1.60%	0.292%
23	14.895	2159	2161	2165	rVV	14509	22655	1.28%	0.232%
24	14.968	2167	2173	2179	rVB3	31557	68429	3.86%	0.702%
25	15.206	2206	2212	2216	rBV2	21688	37065	2.09%	0.380%
26	15.279	2216	2224	2237	rVB3	20450	63561	3.59%	0.652%
27	15.431	2243	2249	2254	rBV	13228	25470	1.44%	0.261%
28	15.590	2268	2275	2285	rVB5	12090	34996	1.97%	0.359%
29	15.785	2301	2307	2312	rBV5	8431	19907	1.12%	0.204%
30	15.931	2323	2331	2341	rBV2	22076	48722	2.75%	0.500%
31	16.376	2397	2404	2410	rBV9	7292	19235	1.08%	0.197%

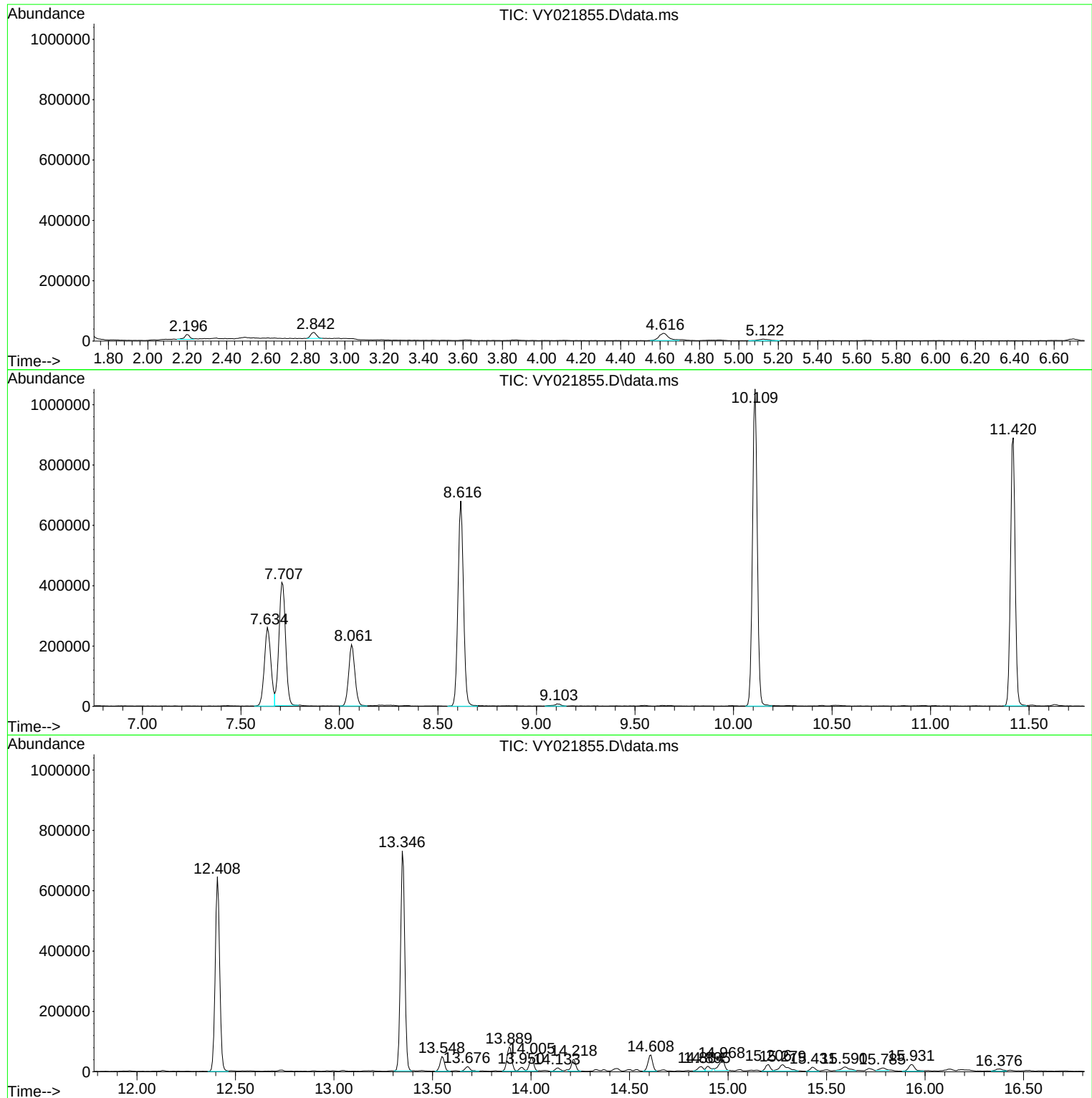
Sum of corrected areas: 9745800

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
Data File : VY021855.D
Acq On : 11 Apr 2025 13:33
Operator : SY/MD
Sample : Q1761-01
Misc : 7.82g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
GST3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
Data File : VY021855.D
Acq On : 11 Apr 2025 13:33
Operator : SY/MD
Sample : Q1761-01
Misc : 7.82g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
GST3

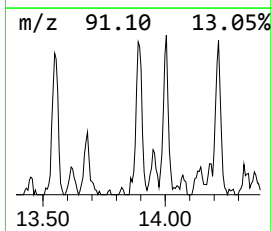
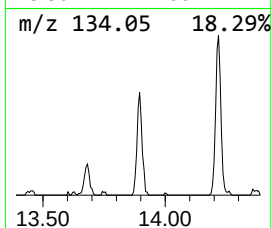
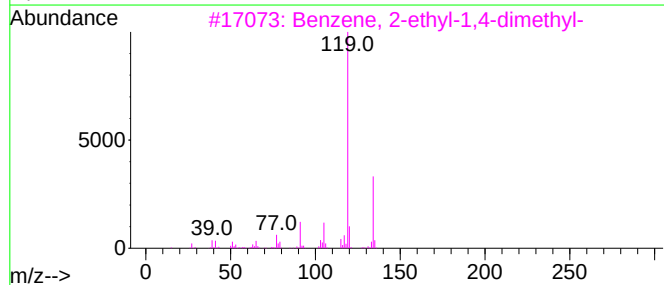
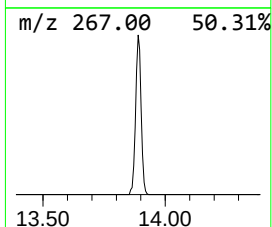
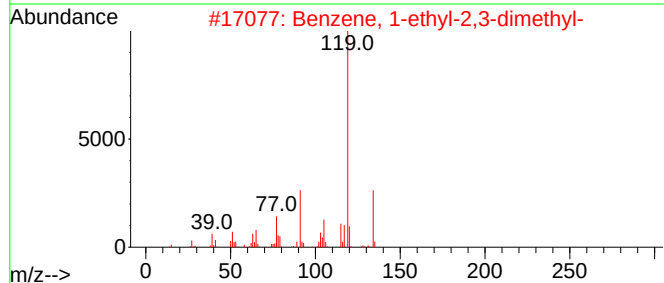
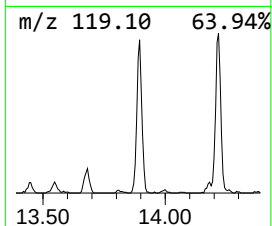
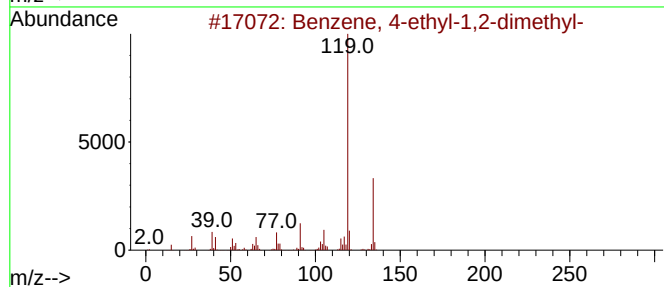
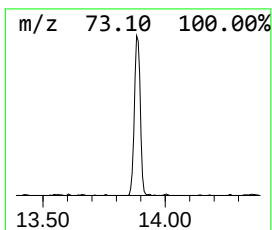
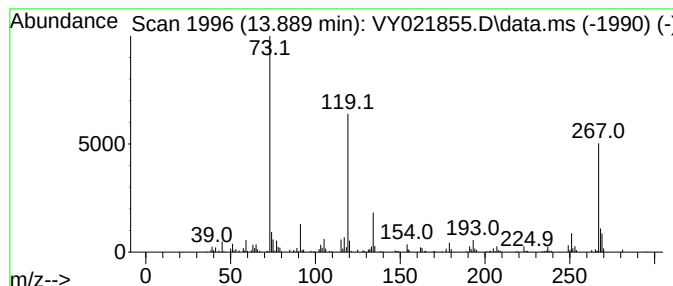
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.889	6.01 ug/l	132922	1,4-Dichlorobenzene-d4	13.346

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	74
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	72
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	70
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	38
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
Data File : VY021855.D
Acq On : 11 Apr 2025 13:33
Operator : SY/MD
Sample : Q1761-01
Misc : 7.82g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
GST3

A

B

C

D

E

F

G

H

I

J

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 4-ethy...	13.889	6.0	ug/l	132922	4	13.346	1106480	50.0

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
 Data File : VY021851.D
 Acq On : 11 Apr 2025 11:35
 Operator : SY/MD
 Sample : VY0411SBL01
 Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 VY0411SBL01

A
B
C
D
E
F
G
H
I
J

Quant Time: Apr 12 02:45:44 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M
 Quant Title : SW846 8260
 QLast Update : Fri Mar 28 02:30:29 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	7.713	168	321354	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.615	114	579974	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.420	117	477999	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.352	152	174550	50.000	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	8.067	65	163798	47.044	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	=	94.080%
35) Dibromofluoromethane	7.634	113	185078	49.194	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	=	98.380%
50) Toluene-d8	10.109	98	678908	47.435	ug/l	0.00
Spiked Amount	50.000	Range	58 - 134	Recovery	=	94.860%
62) 4-Bromofluorobenzene	12.407	95	189286	39.099	ug/l	0.00
Spiked Amount	50.000	Range	30 - 143	Recovery	=	78.200%

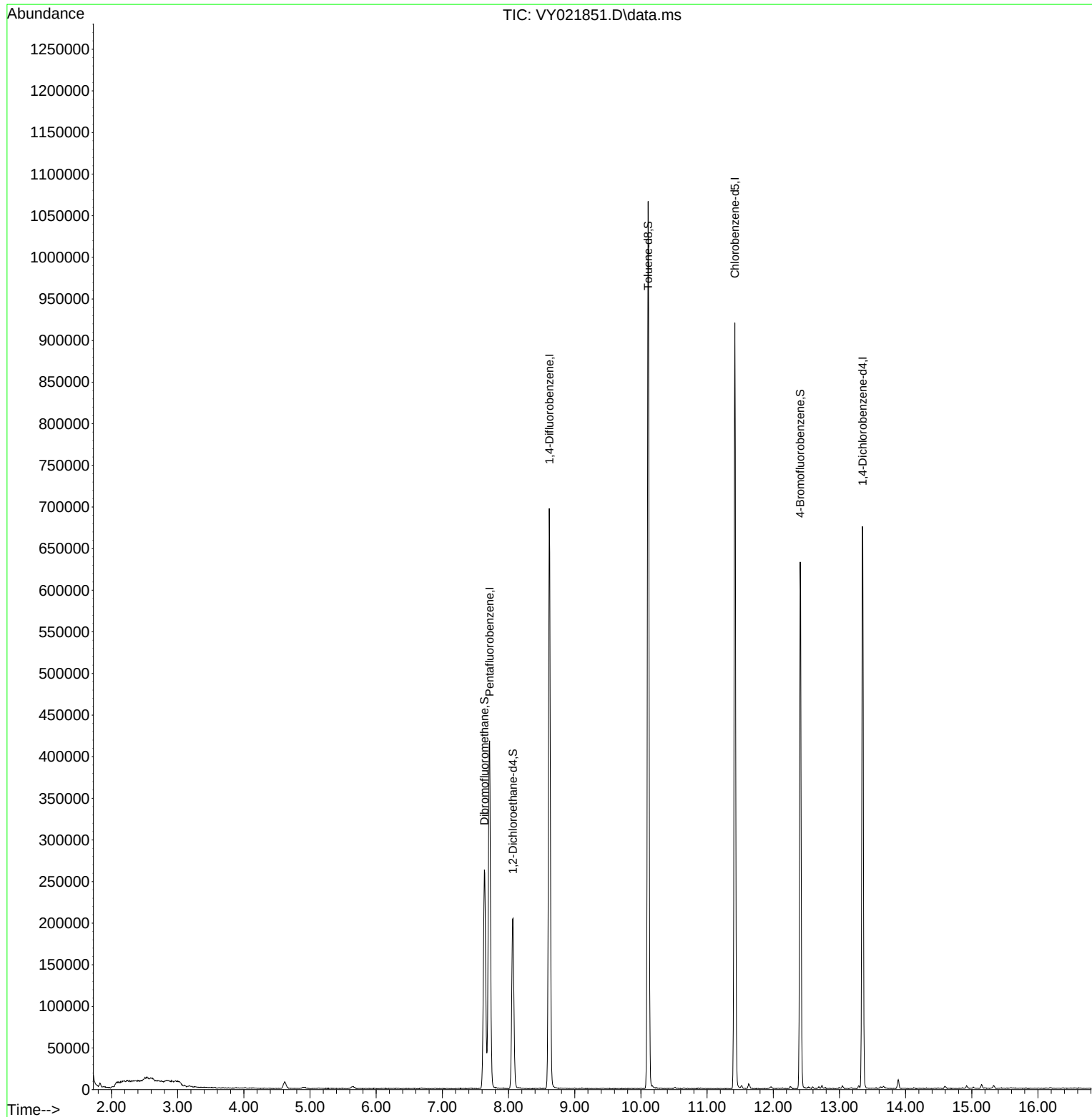
Target Compounds Qvalue

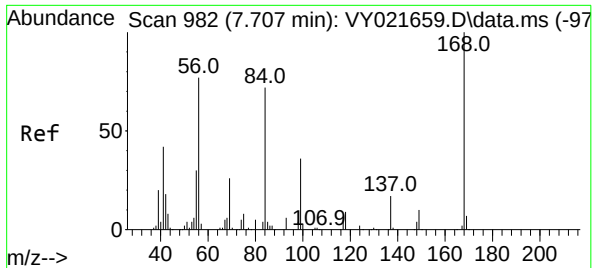
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
Data File : VY021851.D
Acq On : 11 Apr 2025 11:35
Operator : SY/MD
Sample : VY0411SBL01
Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
VY0411SBL01

Quant Time: Apr 12 02:45:44 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M
Quant Title : SW846 8260
QLast Update : Fri Mar 28 02:30:29 2025
Response via : Initial Calibration

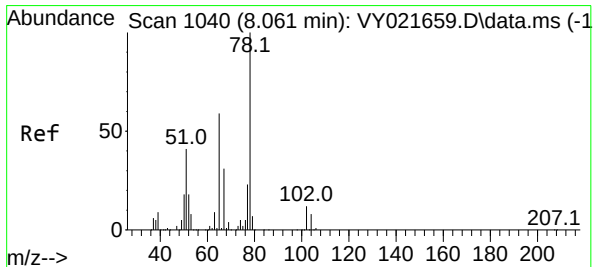
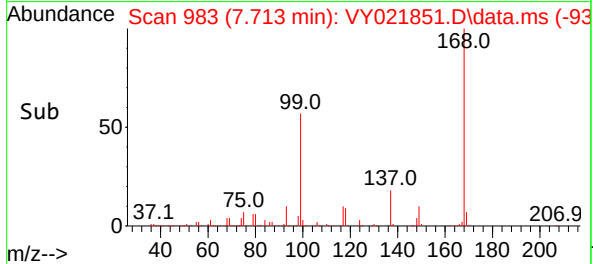
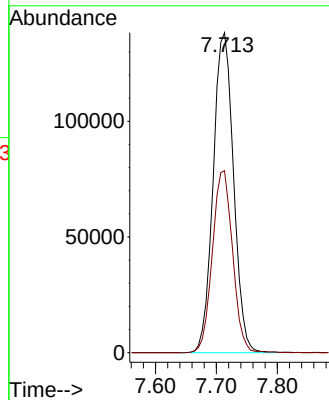
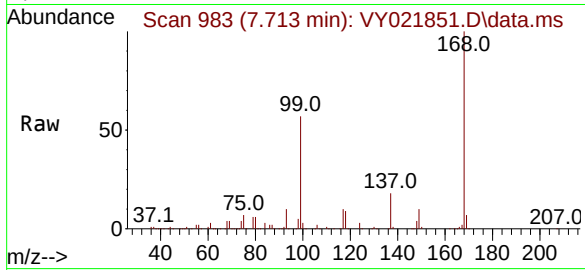




#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 7.713 min Scan# 99
Delta R.T. 0.006 min
Lab File: VY021851.D
Acq: 11 Apr 2025 11:35

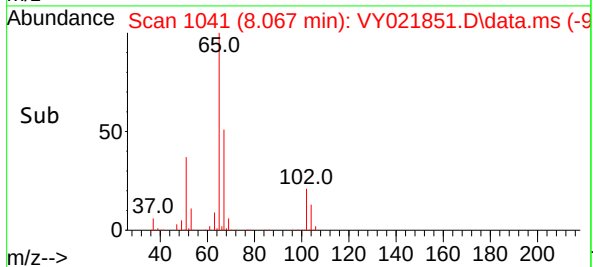
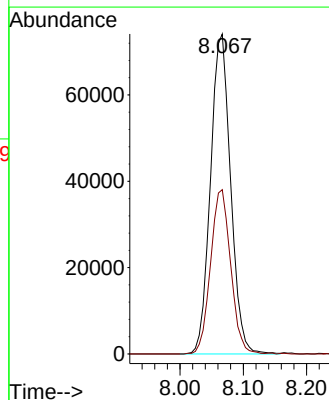
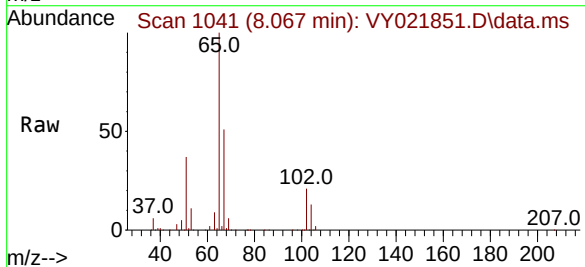
Instrument :
MSVOA_Y
ClientSampleId :
VY0411SBL01

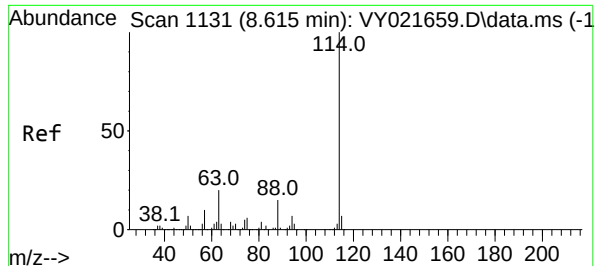
Tgt Ion:168 Resp: 321354
Ion Ratio Lower Upper
168 100
99 56.8 46.7 70.1



#33
1,2-Dichloroethane-d4
Concen: 47.044 ug/l
RT: 8.067 min Scan# 1041
Delta R.T. 0.006 min
Lab File: VY021851.D
Acq: 11 Apr 2025 11:35

Tgt Ion: 65 Resp: 163798
Ion Ratio Lower Upper
65 100
67 51.2 0.0 104.6





#34

1,4-Difluorobenzene

Concen: 50.000 ug/l

RT: 8.615 min Scan# 1131

Delta R.T. 0.000 min

Lab File: VY021851.D

Acq: 11 Apr 2025 11:35

Instrument :

MSVOA_Y

ClientSampleId :

VY0411SBL01

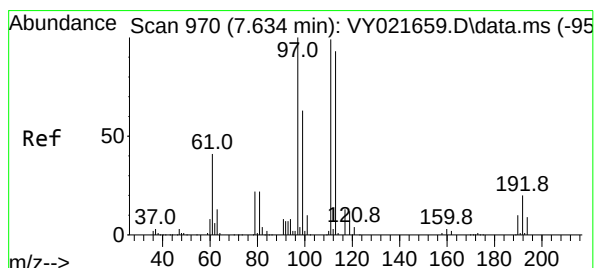
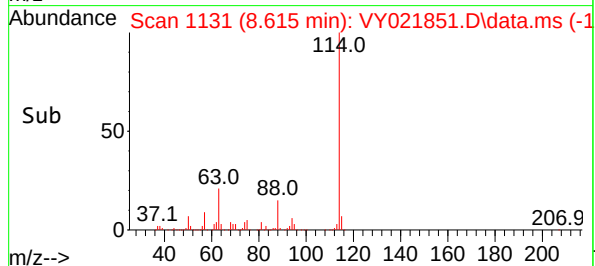
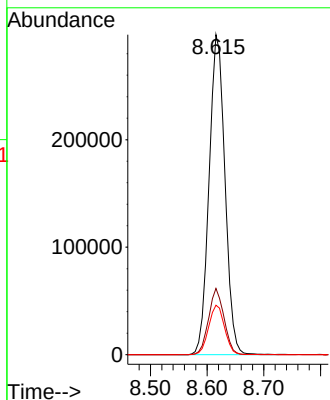
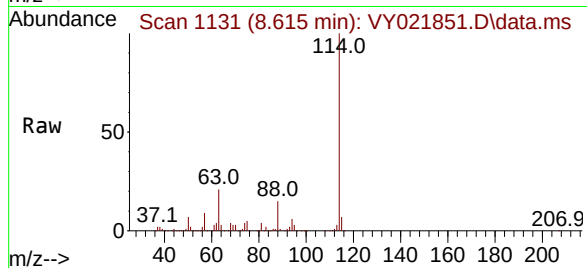
Tgt Ion:114 Resp: 579974

Ion Ratio Lower Upper

114 100

63 20.7 0.0 40.2

88 15.5 0.0 29.8



#35

Dibromofluoromethane

Concen: 49.194 ug/l

RT: 7.634 min Scan# 970

Delta R.T. 0.000 min

Lab File: VY021851.D

Acq: 11 Apr 2025 11:35

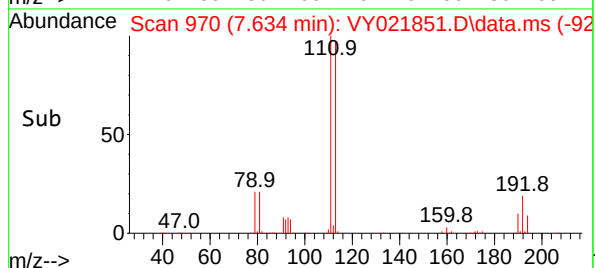
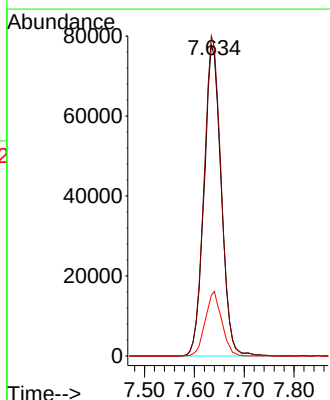
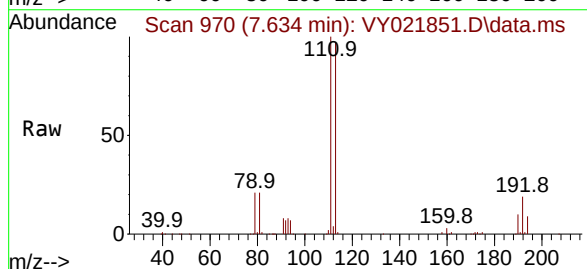
Tgt Ion:113 Resp: 185078

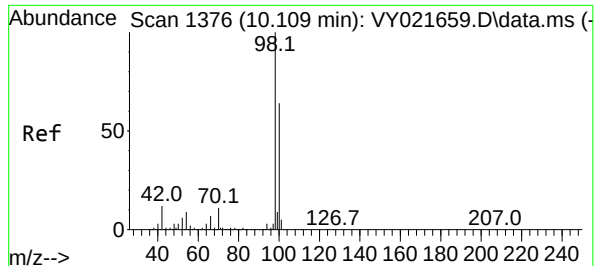
Ion Ratio Lower Upper

113 100

111 103.1 82.6 123.8

192 20.0 16.2 24.4





#50

Toluene-d8

Concen: 47.435 ug/l

RT: 10.109 min Scan# 1376

Delta R.T. 0.000 min

Lab File: VY021851.D

Acq: 11 Apr 2025 11:35

Instrument :

MSVOA_Y

ClientSampleId :

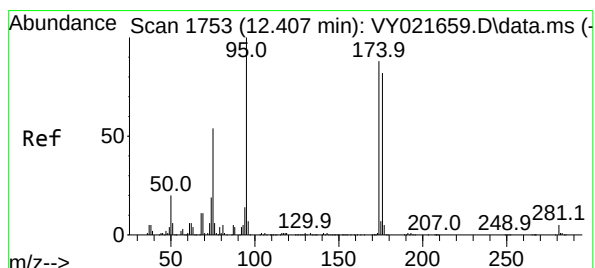
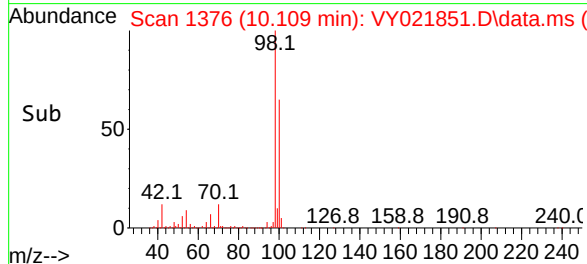
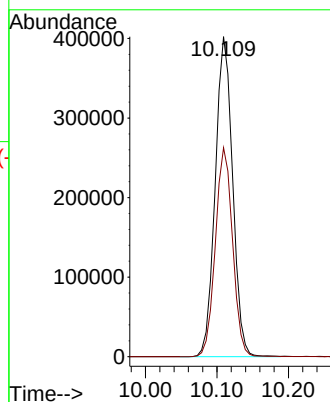
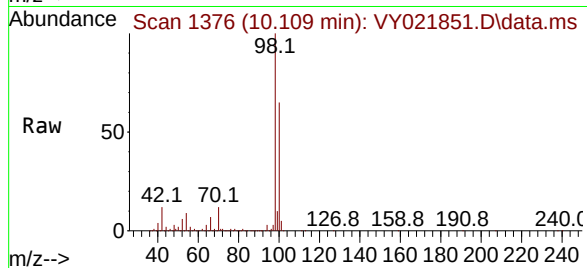
VY0411SBL01

Tgt Ion: 98 Resp: 678908

Ion Ratio Lower Upper

98 100

100 64.7 52.1 78.1



#62

4-Bromofluorobenzene

Concen: 39.099 ug/l

RT: 12.407 min Scan# 1753

Delta R.T. 0.000 min

Lab File: VY021851.D

Acq: 11 Apr 2025 11:35

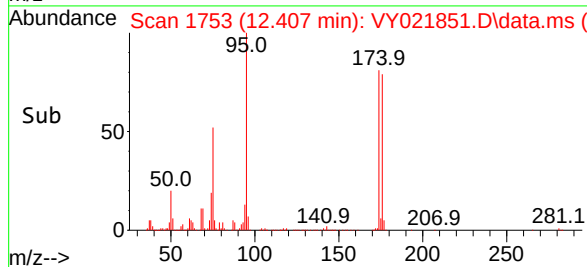
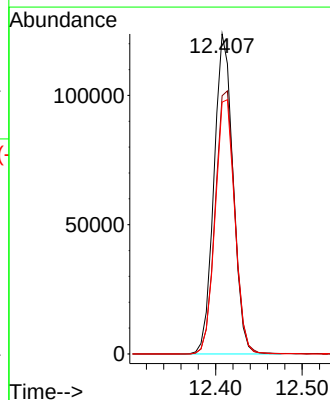
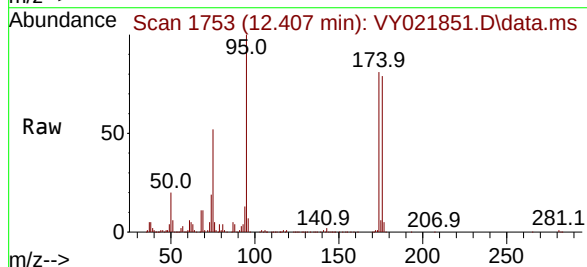
Tgt Ion: 95 Resp: 189286

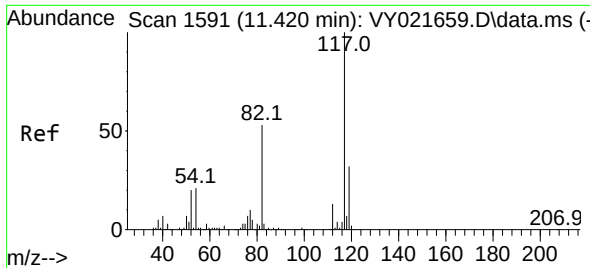
Ion Ratio Lower Upper

95 100

174 84.8 0.0 170.8

176 81.7 0.0 162.0





#63

Chlorobenzene-d5

Concen: 50.000 ug/l

RT: 11.420 min Scan# 1

Delta R.T. 0.000 min

Lab File: VY021851.D

Acq: 11 Apr 2025 11:35

Instrument :

MSVOA_Y

ClientSampleId :

VY0411SBL01

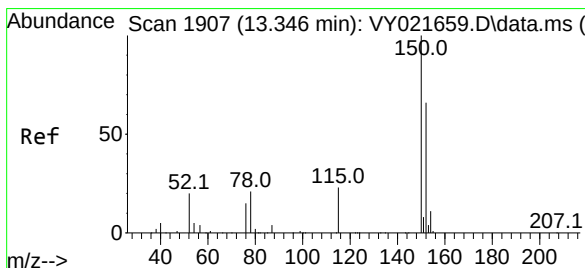
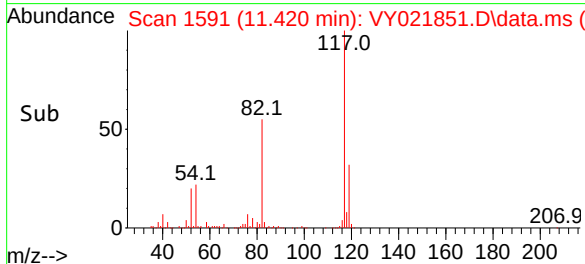
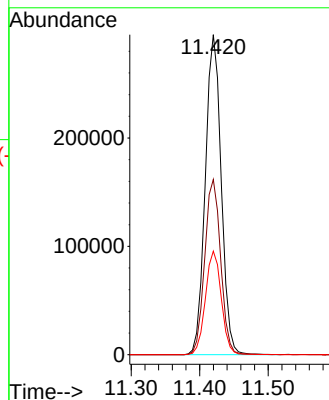
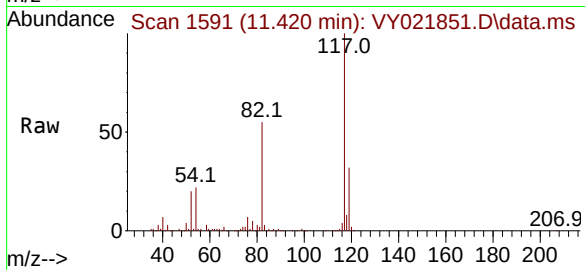
Tgt Ion:117 Resp: 477999

Ion Ratio Lower Upper

117 100

82 54.8 42.8 64.2

119 32.3 25.7 38.5



#72

1,4-Dichlorobenzene-d4

Concen: 50.000 ug/l

RT: 13.352 min Scan# 1908

Delta R.T. 0.006 min

Lab File: VY021851.D

Acq: 11 Apr 2025 11:35

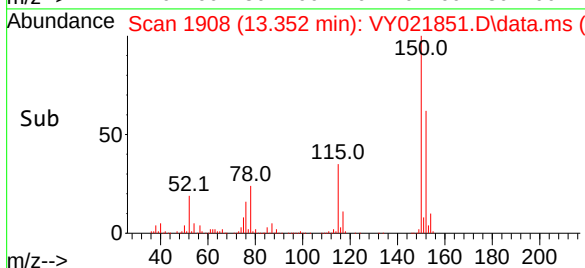
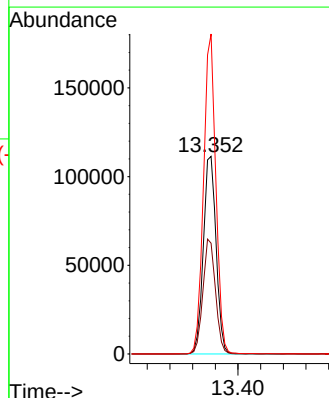
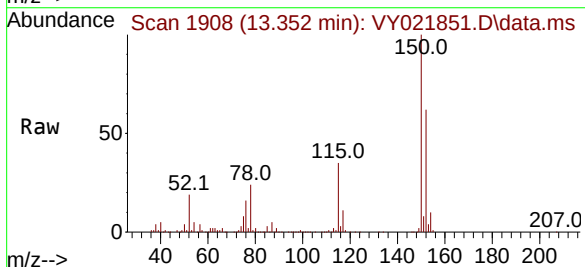
Tgt Ion:152 Resp: 174550

Ion Ratio Lower Upper

152 100

115 57.2 28.8 86.5

150 155.5 0.0 348.4



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
Data File : VY021851.D
Acq On : 11 Apr 2025 11:35
Operator : SY/MD
Sample : VY0411SBL01
Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
VY0411SBL01

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J

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M

Title : SW846 8260

Signal : TIC: VY021851.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.616	466	475	485	rBV4	8002	23227	1.29%	0.265%
2	7.634	961	970	976	rBV	262818	621888	34.48%	7.084%
3	7.713	976	983	994	rVB	416671	967826	53.66%	11.024%
4	8.067	1030	1041	1052	rBV	204614	464921	25.78%	5.296%
5	8.615	1121	1131	1145	rBV	697303	1360201	75.42%	15.493%
6	10.109	1368	1376	1385	rBV	1065866	1803588	100.00%	20.544%
7	11.420	1583	1591	1603	rBV	920013	1490961	82.67%	16.983%
8	12.407	1746	1753	1763	rBV	632880	985592	54.65%	11.226%
9	13.346	1901	1907	1918	rVB	674977	1061129	58.83%	12.087%

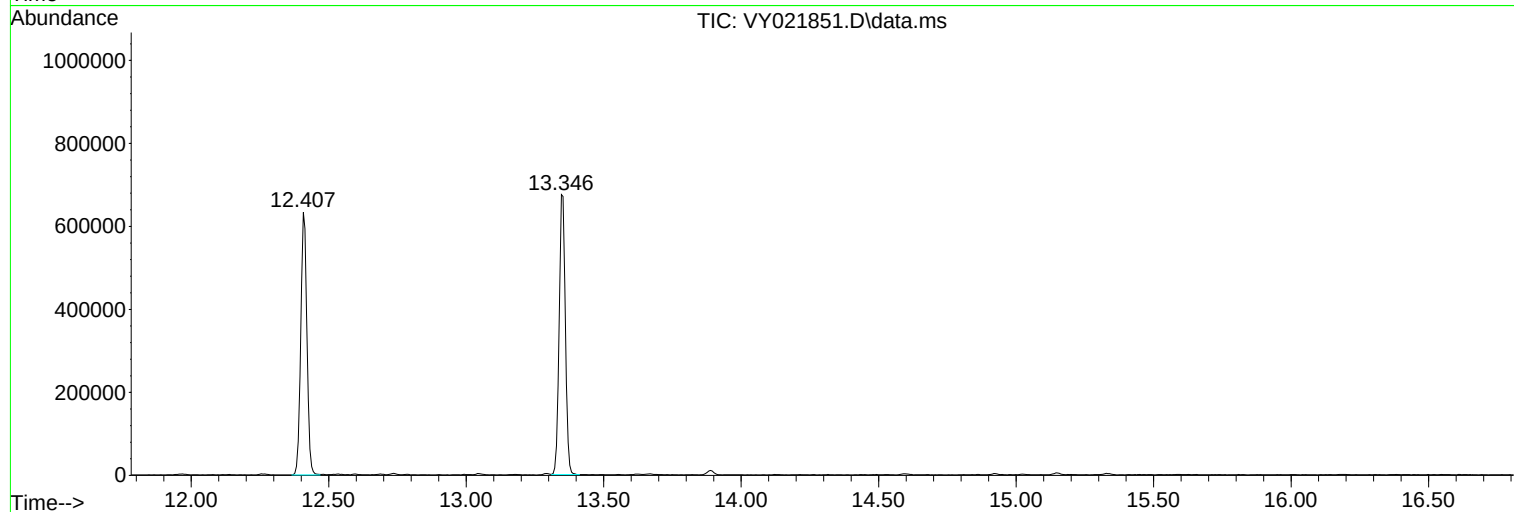
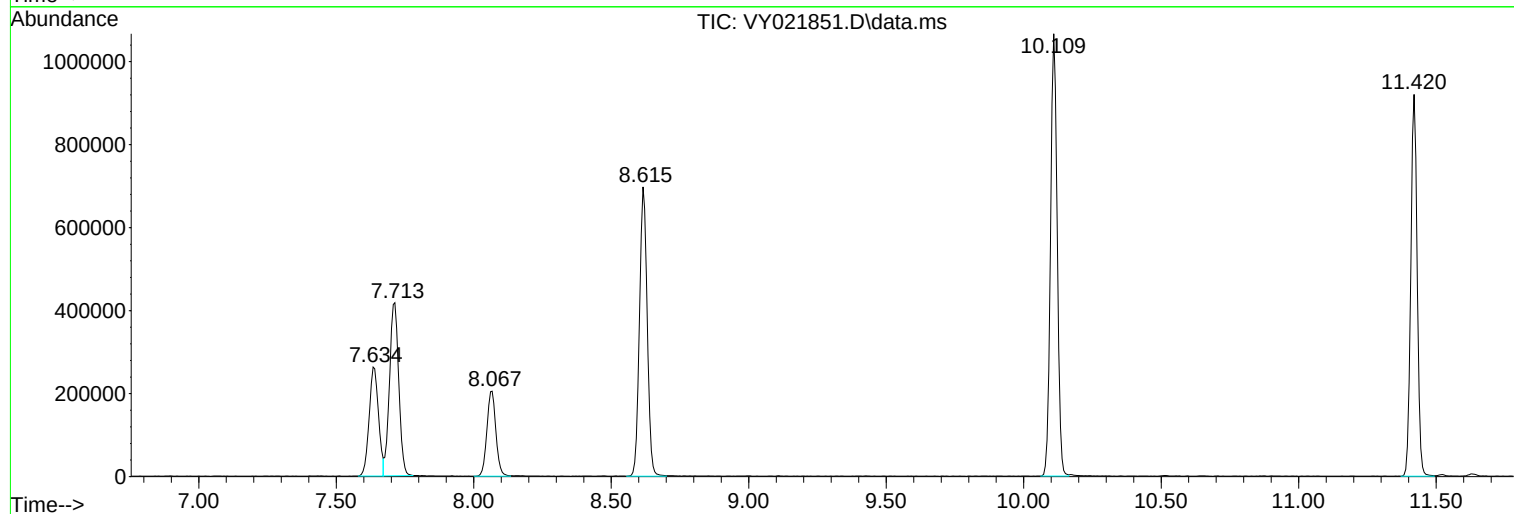
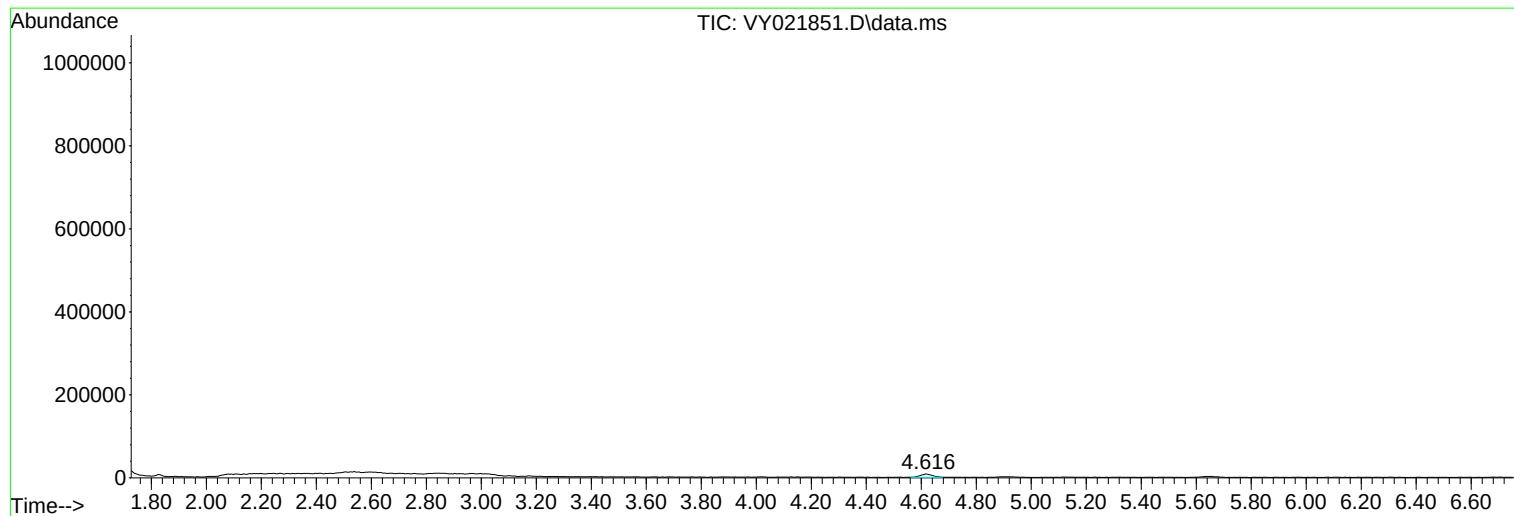
Sum of corrected areas: 8779333

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
Data File : VY021851.D
Acq On : 11 Apr 2025 11:35
Operator : SY/MD
Sample : VY0411SBL01
Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
VY0411SBL01

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
Data File : VY021851.D
Acq On : 11 Apr 2025 11:35
Operator : SY/MD
Sample : VY0411SBL01
Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
VY0411SBL01

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

No Library Search Compounds Detected

A
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J

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
Data File : VY021851.D
Acq On : 11 Apr 2025 11:35
Operator : SY/MD
Sample : VY0411SBL01
Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
VY0411SBL01

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E

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G

H

I

J

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
 Data File : VY021852.D
 Acq On : 11 Apr 2025 12:06
 Operator : SY/MD
 Sample : VY0411SBS01
 Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 VY0411SBS01

Manual Integrations APPROVED

Reviewed By :Mahesh Dadoda 04/14/2025
 Supervised By :Semsettin Yesilyurt 04/14/2025

Quant Time: Apr 12 02:46:09 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M
 Quant Title : SW846 8260
 QLast Update : Fri Mar 28 02:30:29 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Pentafluorobenzene	7.707	168	261639	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.616	114	412441	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.420	117	367570	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.347	152	190311	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.061	65	148170	52.269	ug/l	0.00
Spiked Amount 50.000	Range 50 - 163		Recovery = 104.540%			
35) Dibromofluoromethane	7.634	113	142257	53.172	ug/l	0.00
Spiked Amount 50.000	Range 54 - 147		Recovery = 106.340%			
50) Toluene-d8	10.109	98	547864	53.827	ug/l	0.00
Spiked Amount 50.000	Range 58 - 134		Recovery = 107.660%			
62) 4-Bromofluorobenzene	12.408	95	190322	55.282	ug/l	0.00
Spiked Amount 50.000	Range 30 - 143		Recovery = 110.560%			
Target Compounds						
						Qvalue
2) Dichlorodifluoromethane	1.867	85	47931	17.675	ug/l	100
3) Chloromethane	2.068	50	66874	17.537	ug/l	96
4) Vinyl Chloride	2.202	62	73984	17.156	ug/l	100
5) Bromomethane	2.592	94	55650	18.921	ug/l	92
6) Chloroethane	2.733	64	50881	18.031	ug/l	97
7) Trichlorofluoromethane	3.056	101	102701	18.591	ug/l	99
8) Diethyl Ether	3.458	74	29060	19.344	ug/l	99
9) 1,1,2-Trichlorotrifluo...	3.812	101	56995	19.562	ug/l	99
10) Methyl Iodide	4.007	142	65715	20.559	ug/l	99
11) Tert butyl alcohol	4.866	59	17296	93.935	ug/l #	87
12) 1,1-Dichloroethene	3.787	96	52342	19.073	ug/l	97
13) Acrolein	3.653	56	17310	51.486	ug/l	98
14) Allyl chloride	4.385	41	81064	18.526	ug/l	98
15) Acrylonitrile	5.061	53	60829	96.976	ug/l	100
16) Acetone	3.879	43	58163	100.252	ug/l	97
17) Carbon Disulfide	4.104	76	157086	17.379	ug/l	99
18) Methyl Acetate	4.391	43	29819	20.738	ug/l	97
19) Methyl tert-butyl Ether	5.116	73	135524	19.835	ug/l	98
20) Methylene Chloride	4.610	84	63577	20.080	ug/l	95
21) trans-1,2-Dichloroethene	5.110	96	57161	18.917	ug/l	95
22) Diisopropyl ether	6.019	45	177421	19.329	ug/l	99
23) Vinyl Acetate	5.964	43	503378	93.744	ug/l	97
24) 1,1-Dichloroethane	5.915	63	107900	19.659	ug/l	99
25) 2-Butanone	6.896	43	78937	95.259	ug/l	100
26) 2,2-Dichloropropane	6.890	77	95824	19.627	ug/l	99
27) cis-1,2-Dichloroethene	6.890	96	67782	19.948	ug/l	97
28) Bromochloromethane	7.244	49	45578	19.691	ug/l	95
29) Tetrahydrofuran	7.262	42	49850	93.472	ug/l	97
30) Chloroform	7.421	83	112631	19.730	ug/l	96
31) Cyclohexane	7.701	56	89585	17.905	ug/l	96
32) 1,1,1-Trichloroethane	7.622	97	100880	19.552	ug/l	99
36) 1,1-Dichloropropene	7.835	75	77267	18.961	ug/l	98
37) Ethyl Acetate	6.988	43	36325	18.888	ug/l	100
38) Carbon Tetrachloride	7.817	117	91970	19.705	ug/l	99
39) Methylcyclohexane	9.109	83	91679	18.832	ug/l	98
40) Benzene	8.079	78	240011	19.667	ug/l	98

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
 Data File : VY021852.D
 Acq On : 11 Apr 2025 12:06
 Operator : SY/MD
 Sample : VY0411SBS01
 Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 VY0411SBS01

Manual Integrations
 APPROVED

Reviewed By :Mahesh Dadoda 04/14/2025
 Supervised By :Semsettin Yesilyurt 04/14/2025

Quant Time: Apr 12 02:46:09 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M
 Quant Title : SW846 8260
 QLast Update : Fri Mar 28 02:30:29 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	7.220	41	19246m	18.843	ug/l	
42) 1,2-Dichloroethane	8.158	62	68214	19.827	ug/l	99
43) Isopropyl Acetate	8.201	43	69884	19.174	ug/l #	85
44) Trichloroethene	8.866	130	61686	19.838	ug/l	94
45) 1,2-Dichloropropane	9.140	63	56361	19.511	ug/l	98
46) Dibromomethane	9.231	93	31831	19.119	ug/l	97
47) Bromodichloromethane	9.426	83	86291	20.003	ug/l	97
48) Methyl methacrylate	9.219	41	31606	18.895	ug/l	98
49) 1,4-Dioxane	9.237	88	7141	400.396	ug/l	98
51) 4-Methyl-2-Pentanone	10.000	43	184332	97.188	ug/l	100
52) Toluene	10.170	92	152947	20.087	ug/l	98
53) t-1,3-Dichloropropene	10.396	75	74193	19.363	ug/l	98
54) cis-1,3-Dichloropropene	9.859	75	89915	19.891	ug/l	96
55) 1,1,2-Trichloroethane	10.573	97	43135	20.133	ug/l	96
56) Ethyl methacrylate	10.438	69	53948	19.720	ug/l	97
57) 1,3-Dichloropropane	10.719	76	75191	20.321	ug/l	100
58) 2-Chloroethyl Vinyl ether	9.713	63	136203	102.058	ug/l	98
59) 2-Hexanone	10.762	43	122817	97.231	ug/l	100
60) Dibromochloromethane	10.914	129	59311	20.276	ug/l	99
61) 1,2-Dibromoethane	11.018	107	41010	20.405	ug/l	99
64) Tetrachloroethene	10.646	164	73872	22.066	ug/l	98
65) Chlorobenzene	11.444	112	166748	19.790	ug/l	97
66) 1,1,1,2-Tetrachloroethane	11.518	131	59523	19.911	ug/l	98
67) Ethyl Benzene	11.518	91	281537	19.485	ug/l	100
68) m/p-Xylenes	11.627	106	218365	39.468	ug/l	99
69) o-Xylene	11.957	106	102058	19.983	ug/l	99
70) Styrene	11.969	104	168475	19.785	ug/l	99
71) Bromoform	12.133	173	34728	20.365	ug/l #	99
73) Isopropylbenzene	12.255	105	267864	18.950	ug/l	99
74) N-amyl acetate	12.072	43	60183	18.142	ug/l	98
75) 1,1,2,2-Tetrachloroethane	12.505	83	45580	17.935	ug/l	99
76) 1,2,3-Trichloropropane	12.560	75	38417m	20.364	ug/l	
77) Bromobenzene	12.536	156	65230	19.220	ug/l	99
78) n-propylbenzene	12.597	91	326537	19.215	ug/l	100
79) 2-Chlorotoluene	12.682	91	186879	18.933	ug/l	100
80) 1,3,5-Trimethylbenzene	12.737	105	224041	19.393	ug/l	99
81) trans-1,4-Dichloro-2-b...	12.304	75	14546	17.767	ug/l	99
82) 4-Chlorotoluene	12.780	91	195026	18.953	ug/l	98
83) tert-Butylbenzene	12.999	119	193537	18.809	ug/l	100
84) 1,2,4-Trimethylbenzene	13.048	105	225265	19.729	ug/l	99
85) sec-Butylbenzene	13.176	105	288462	19.183	ug/l	99
86) p-Isopropyltoluene	13.292	119	240050	19.314	ug/l	100
87) 1,3-Dichlorobenzene	13.292	146	129499	19.367	ug/l	99
88) 1,4-Dichlorobenzene	13.371	146	128644	19.473	ug/l	96
89) n-Butylbenzene	13.621	91	215648	18.762	ug/l	100
90) Hexachloroethane	13.883	117	51012	18.730	ug/l	96
91) 1,2-Dichlorobenzene	13.657	146	117128	20.123	ug/l	99
92) 1,2-Dibromo-3-Chloropr...	14.273	75	8235	20.907	ug/l	93
93) 1,2,4-Trichlorobenzene	14.919	180	66752	21.147	ug/l	99
94) Hexachlorobutadiene	15.023	225	41048	20.822	ug/l	99
95) Naphthalene	15.145	128	108783	20.792	ug/l	100
96) 1,2,3-Trichlorobenzene	15.328	180	57607	21.402	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
Data File : VY021852.D
Acq On : 11 Apr 2025 12:06
Operator : SY/MD
Sample : VY0411SBS01
Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
VY0411SBS01

Manual Integrations
APPROVED

Reviewed By :Mahesh Dadoda 04/14/2025
Supervised By :Semsettin Yesilyurt 04/14/2025

Quant Time: Apr 12 02:46:09 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M
Quant Title : SW846 8260
QLast Update : Fri Mar 28 02:30:29 2025
Response via : Initial Calibration

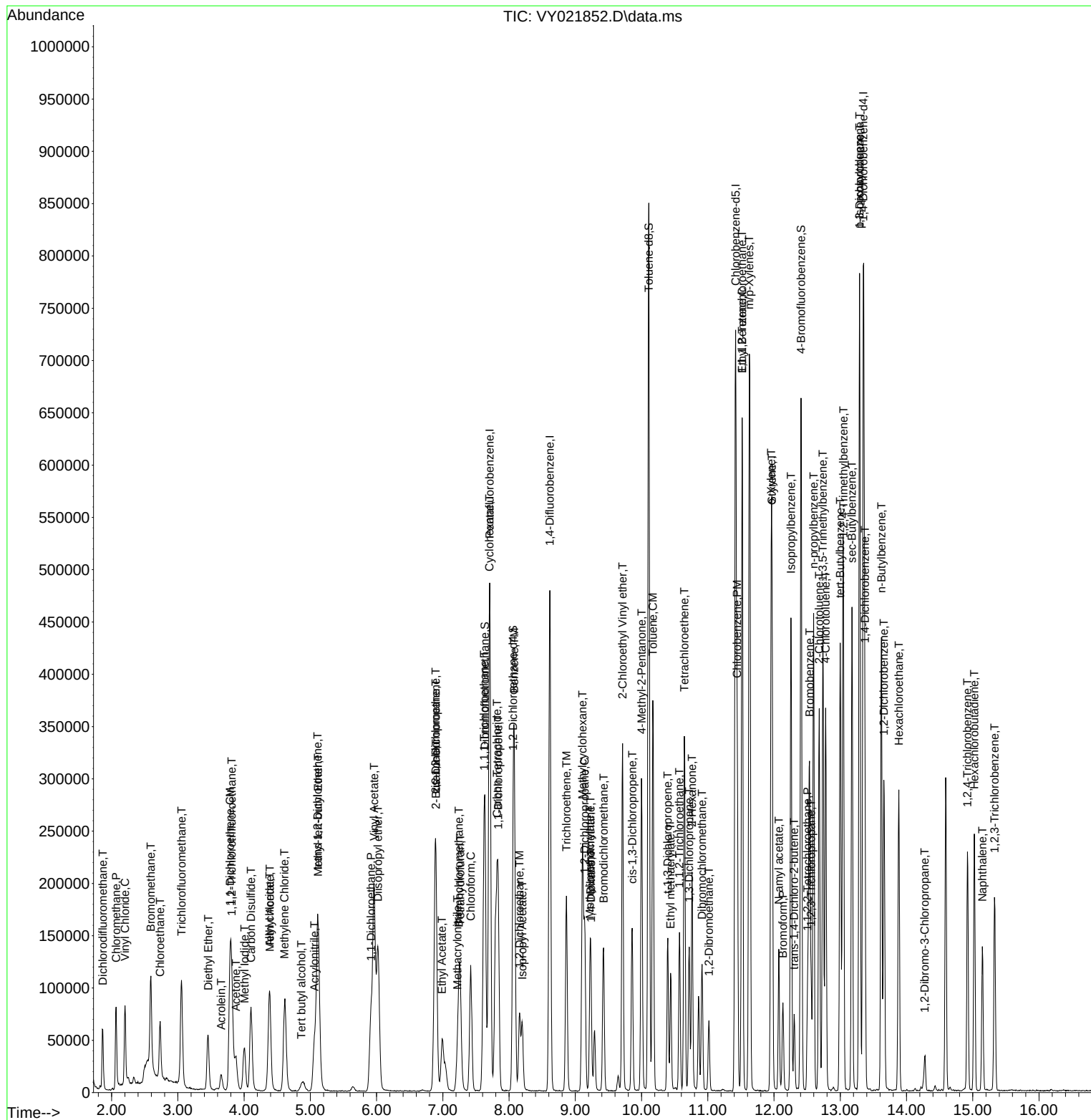
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

Instrument :
MSVOA_Y
ClientSampleId :
VY0411SBS01

Manual Integrations APPROVED

Reviewed By :Mahesh Dadoda 04/14/2025
Supervised By :Semsettin Yesilyurt 04/14/2025



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
 Data File : VY021853.D
 Acq On : 11 Apr 2025 12:28
 Operator : SY/MD
 Sample : VY0411SBSD01
 Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 VY0411SBSD01

Manual Integrations APPROVED

Reviewed By :Mahesh Dadoda 04/14/2025
 Supervised By :Semsettin Yesilyurt 04/14/2025

Quant Time: Apr 12 02:47:01 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M
 Quant Title : SW846 8260
 QLast Update : Fri Mar 28 02:30:29 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	7.707	168	261813	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.616	114	409914	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.414	117	362626	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.347	152	188335	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.061	65	143603	50.624	ug/l	0.00
Spiked Amount 50.000	Range 50 - 163		Recovery = 101.240%			
35) Dibromofluoromethane	7.634	113	143350	53.910	ug/l	0.00
Spiked Amount 50.000	Range 54 - 147		Recovery = 107.820%			
50) Toluene-d8	10.103	98	537312	53.116	ug/l	0.00
Spiked Amount 50.000	Range 58 - 134		Recovery = 106.240%			
62) 4-Bromofluorobenzene	12.408	95	186390	54.474	ug/l	0.00
Spiked Amount 50.000	Range 30 - 143		Recovery = 108.940%			
Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.861	85	49656	18.299	ug/l	99
3) Chloromethane	2.068	50	62767	16.449	ug/l	97
4) Vinyl Chloride	2.202	62	73201	16.963	ug/l	97
5) Bromomethane	2.592	94	51528	17.508	ug/l	100
6) Chloroethane	2.733	64	48298	17.104	ug/l	99
7) Trichlorofluoromethane	3.056	101	104991	18.993	ug/l	98
8) Diethyl Ether	3.458	74	29054	19.327	ug/l	96
9) 1,1,2-Trichlorotrifluo...	3.818	101	56846	19.498	ug/l	96
10) Methyl Iodide	4.001	142	67929	21.238	ug/l	99
11) Tert butyl alcohol	4.866	59	16952	92.006	ug/l #	90
12) 1,1-Dichloroethene	3.787	96	54499	19.846	ug/l	94
13) Acrolein	3.647	56	15884	47.214	ug/l	99
14) Allyl chloride	4.379	41	82696	18.887	ug/l	98
15) Acrylonitrile	5.061	53	59158	94.249	ug/l	99
16) Acetone	3.873	43	49207	84.759	ug/l	95
17) Carbon Disulfide	4.104	76	162030	17.914	ug/l	97
18) Methyl Acetate	4.379	43	31345	21.785	ug/l	98
19) Methyl tert-butyl Ether	5.122	73	133728	19.559	ug/l	97
20) Methylene Chloride	4.610	84	64347	20.358	ug/l	96
21) trans-1,2-Dichloroethene	5.110	96	58865	19.468	ug/l	97
22) Diisopropyl ether	6.019	45	182652	19.886	ug/l	98
23) Vinyl Acetate	5.958	43	498605	92.794	ug/l	98
24) 1,1-Dichloroethane	5.915	63	109510	19.939	ug/l	98
25) 2-Butanone	6.890	43	75132	90.607	ug/l	98
26) 2,2-Dichloropropane	6.884	77	97186	19.893	ug/l	99
27) cis-1,2-Dichloroethene	6.890	96	68636	20.185	ug/l	96
28) Bromochloromethane	7.238	49	41634	17.975	ug/l	92
29) Tetrahydrofuran	7.262	42	49127	92.055	ug/l	98
30) Chloroform	7.421	83	114067	19.968	ug/l	98
31) Cyclohexane	7.701	56	91074	18.191	ug/l	90
32) 1,1,1-Trichloroethane	7.616	97	103851	20.115	ug/l	98
36) 1,1-Dichloropropene	7.835	75	78621	19.412	ug/l	98
37) Ethyl Acetate	6.982	43	35272	18.453	ug/l	98
38) Carbon Tetrachloride	7.817	117	93678	20.194	ug/l	97
39) Methylcyclohexane	9.110	83	92467	19.111	ug/l	98
40) Benzene	8.079	78	246182	20.297	ug/l	100

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
 Data File : VY021853.D
 Acq On : 11 Apr 2025 12:28
 Operator : SY/MD
 Sample : VY0411SBS01
 Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 VY0411SBS01

Manual Integrations
 APPROVED

Reviewed By :Mahesh Dadoda 04/14/2025
 Supervised By :Semsettin Yesilyurt 04/14/2025

Quant Time: Apr 12 02:47:01 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M
 Quant Title : SW846 8260
 QLast Update : Fri Mar 28 02:30:29 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	7.220	41	21263	20.947	ug/l #	100
42) 1,2-Dichloroethane	8.158	62	66951	19.580	ug/l	100
43) Isopropyl Acetate	8.195	43	67897	18.744	ug/l	98
44) Trichloroethene	8.866	130	64547	20.886	ug/l	100
45) 1,2-Dichloropropane	9.140	63	57185	19.918	ug/l	94
46) Dibromomethane	9.231	93	33009	19.949	ug/l	99
47) Bromodichloromethane	9.420	83	85927	20.041	ug/l	99
48) Methyl methacrylate	9.219	41	32878	19.776	ug/l	98
49) 1,4-Dioxane	9.225	88	7355	414.937	ug/l	97
51) 4-Methyl-2-Pentanone	10.000	43	178028	94.443	ug/l	98
52) Toluene	10.170	92	155125	20.499	ug/l	100
53) t-1,3-Dichloropropene	10.396	75	75693	19.876	ug/l	98
54) cis-1,3-Dichloropropene	9.853	75	87863	19.557	ug/l	97
55) 1,1,2-Trichloroethane	10.573	97	43853	20.595	ug/l	98
56) Ethyl methacrylate	10.439	69	54410	20.012	ug/l	96
57) 1,3-Dichloropropane	10.719	76	73973	20.115	ug/l	98
58) 2-Chloroethyl Vinyl ether	9.707	63	136529	102.933	ug/l	98
59) 2-Hexanone	10.762	43	117310	93.444	ug/l	99
60) Dibromochloromethane	10.908	129	59548	20.483	ug/l	99
61) 1,2-Dibromoethane	11.012	107	39785	19.918	ug/l	95
64) Tetrachloroethene	10.646	164	73897	22.375	ug/l	97
65) Chlorobenzene	11.444	112	169509	20.392	ug/l	99
66) 1,1,1,2-Tetrachloroethane	11.518	131	61208	20.753	ug/l	99
67) Ethyl Benzene	11.518	91	285855	20.054	ug/l	99
68) m/p-Xylenes	11.627	106	222848	40.827	ug/l	99
69) o-Xylene	11.957	106	103070	20.457	ug/l	100
70) Styrene	11.969	104	172929	20.585	ug/l	98
71) Bromoform	12.133	173	34263	20.366	ug/l #	100
73) Isopropylbenzene	12.255	105	276145	19.741	ug/l	100
74) N-amyl acetate	12.072	43	57979	17.661	ug/l	95
75) 1,1,2,2-Tetrachloroethane	12.505	83	45218	17.980	ug/l	100
76) 1,2,3-Trichloropropane	12.554	75	35519m	19.026	ug/l	
77) Bromobenzene	12.530	156	66160	19.699	ug/l	97
78) n-propylbenzene	12.597	91	331933	19.738	ug/l	99
79) 2-Chlorotoluene	12.682	91	194696	19.932	ug/l	99
80) 1,3,5-Trimethylbenzene	12.737	105	227007	19.856	ug/l	99
81) trans-1,4-Dichloro-2-b...	12.304	75	14585	18.001	ug/l	97
82) 4-Chlorotoluene	12.780	91	200035	19.643	ug/l	98
83) tert-Butylbenzene	12.999	119	201861	19.823	ug/l	100
84) 1,2,4-Trimethylbenzene	13.042	105	228682	20.238	ug/l	100
85) sec-Butylbenzene	13.176	105	294858	19.814	ug/l	98
86) p-Isopropyltoluene	13.292	119	249819	20.311	ug/l	99
87) 1,3-Dichlorobenzene	13.286	146	134109	20.267	ug/l	100
88) 1,4-Dichlorobenzene	13.365	146	131817	20.163	ug/l	98
89) n-Butylbenzene	13.615	91	224734	19.757	ug/l	99
90) Hexachloroethane	13.883	117	52557	19.500	ug/l	98
91) 1,2-Dichlorobenzene	13.657	146	115715	20.089	ug/l	99
92) 1,2-Dibromo-3-Chloropr...	14.273	75	7745	19.869	ug/l	99
93) 1,2,4-Trichlorobenzene	14.919	180	68380	21.890	ug/l	99
94) Hexachlorobutadiene	15.023	225	43077	22.081	ug/l	99
95) Naphthalene	15.145	128	111977	21.627	ug/l	100
96) 1,2,3-Trichlorobenzene	15.328	180	59827	22.460	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041125\
Data File : VY021853.D
Acq On : 11 Apr 2025 12:28
Operator : SY/MD
Sample : VY0411SBSD01
Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
VY0411SBSD01

Manual Integrations
APPROVED

Reviewed By :Mahesh Dadoda 04/14/2025
Supervised By :Semsettin Yesilyurt 04/14/2025

Quant Time: Apr 12 02:47:01 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032725S.M
Quant Title : SW846 8260
QLast Update : Fri Mar 28 02:30:29 2025
Response via : Initial Calibration

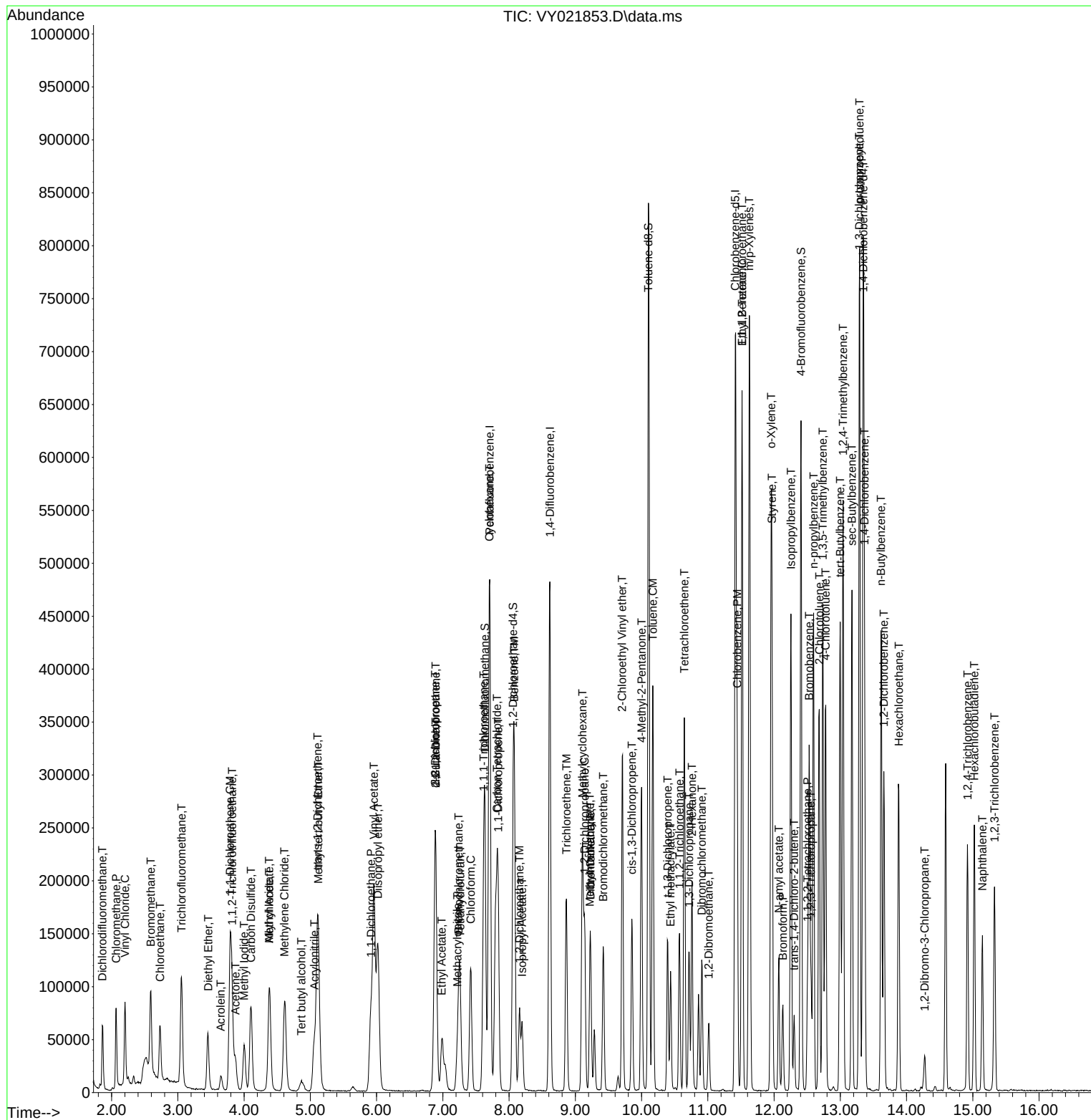
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

Instrument :
MSVOA_Y
ClientSampleId :
VY0411SBSD01

Manual Integrations APPROVED

Reviewed By :Mahesh Dadoda 04/14/2025
Supervised By :Semsettin Yesilyurt 04/14/2025



Manual Integration Report

Sequence:	vy032725	Instrument	MSVOA_y
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
VSTDICC005	VY021656.D	1,2,3-Trichloropropane	Romaben	3/28/2025 9:14:28 AM	MMDadoda	3/28/2025 9:32:34 AM	Peak Integrated by Software
VSTDICC005	VY021656.D	Methacrylonitrile	Romaben	3/28/2025 9:14:28 AM	MMDadoda	3/28/2025 9:32:34 AM	Peak Integrated by Software
VSTDICC010	VY021657.D	1,2,3-Trichloropropane	Romaben	3/28/2025 9:14:32 AM	MMDadoda	3/28/2025 9:32:37 AM	Peak Integrated by Software
VSTDICC010	VY021657.D	Methacrylonitrile	Romaben	3/28/2025 9:14:32 AM	MMDadoda	3/28/2025 9:32:37 AM	Peak Integrated by Software
VSTDICC020	VY021658.D	1,2,3-Trichloropropane	Romaben	3/28/2025 9:15:20 AM	MMDadoda	3/28/2025 9:32:41 AM	Peak Integrated by Software
VSTDICCC050	VY021659.D	1,2,3-Trichloropropane	Romaben	3/28/2025 9:14:38 AM	MMDadoda	3/28/2025 9:32:45 AM	Peak Integrated by Software
VSTDICCC050	VY021659.D	Methacrylonitrile	Romaben	3/28/2025 9:14:38 AM	MMDadoda	3/28/2025 9:32:45 AM	Peak Integrated by Software
VSTDICC100	VY021660.D	1,2,3-Trichloropropane	Romaben	3/28/2025 9:15:15 AM	MMDadoda	3/28/2025 9:32:49 AM	Peak Integrated by Software
VSTDICC150	VY021661.D	1,2,3-Trichloropropane	Romaben	3/28/2025 9:14:43 AM	MMDadoda	3/28/2025 9:32:53 AM	Peak Integrated by Software
VSTDICV050	VY021663.D	1,2,3-Trichloropropane	Romaben	3/28/2025 9:15:11 AM	MMDadoda	3/28/2025 9:32:04 AM	Peak Integrated by Software
VSTDCCC050	VY021680.D	1,2,3-Trichloropropane	Romaben	3/28/2025 9:14:58 AM	MMDadoda	3/28/2025 9:32:15 AM	Peak Integrated by Software

Manual Integration Report

Sequence:

VY041125

Instrument

MSVOA_y

Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
VSTDCCC050	VY021850.D	1,2,3-Trichloropropane	MMDadoda	4/14/2025 3:42:39 PM	SAM	4/14/2025 3:45:07 PM	Peak Integrated by Software
VY0411SBS01	VY021852.D	1,2,3-Trichloropropane	MMDadoda	4/14/2025 3:42:40 PM	SAM	4/14/2025 3:45:07 PM	Peak Integrated by Software
VY0411SBS01	VY021852.D	Methacrylonitrile	MMDadoda	4/14/2025 3:42:40 PM	SAM	4/14/2025 3:45:07 PM	Peak Integrated by Software
VY0411SBSD01	VY021853.D	1,2,3-Trichloropropane	MMDadoda	4/14/2025 3:42:42 PM	SAM	4/14/2025 3:45:08 PM	Peak Integrated by Software
VSTDCCC050	VY021863.D	1,2,3-Trichloropropane	MMDadoda	4/14/2025 3:42:44 PM	SAM	4/14/2025 3:45:09 PM	Peak Integrated by Software

Instrument ID: **MSVOA_Y**

Daily Analysis Runlog For Sequence/QCBatch ID # VY032725

Review By	Maresh Dadoda	Review On	3/28/2025 3:29:57 PM
Supervise By	Semsettin Yesilyurt	Supervise On	3/28/2025 3:30:48 PM
SubDirectory	VY032725	HP Acquire Method	MSVOA_Y
		HP Processing Method	82y032725s.m
STD. NAME	STD REF.#		
Tune/Reschk	VP133499		
Initial Calibration Stds	VP133500,VP133501,VP133502,VP133503,VP133505,VP133508		
CCC	VP133510		
Internal Standard/PEM	VP131783		
ICV/I.BLK	VP133509		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

Sr#	SampleId	Data File Name	Date-Time	Operator	Status
1	BFB	VY021655.D	27 Mar 2025 09:23	SY/MD	Ok
2	VSTDIC005	VY021656.D	27 Mar 2025 10:08	SY/MD	Ok,M
3	VSTDIC010	VY021657.D	27 Mar 2025 10:31	SY/MD	Ok,M
4	VSTDIC020	VY021658.D	27 Mar 2025 10:54	SY/MD	Ok,M
5	VSTDIC050	VY021659.D	27 Mar 2025 11:16	SY/MD	Ok,M
6	VSTDIC100	VY021660.D	27 Mar 2025 11:39	SY/MD	Ok,M
7	VSTDIC150	VY021661.D	27 Mar 2025 12:02	SY/MD	Ok,M
8	VIBLK	VY021662.D	27 Mar 2025 12:35	SY/MD	Ok
9	VSTDICV050	VY021663.D	27 Mar 2025 13:15	SY/MD	Ok,M
10	VY0327SBL01	VY021664.D	27 Mar 2025 13:59	SY/MD	Ok
11	VY0327SBS01	VY021665.D	27 Mar 2025 14:38	SY/MD	Ok,M
12	VY0327SBS01	VY021666.D	27 Mar 2025 15:00	SY/MD	Ok,M
13	Q1661-02	VY021667.D	27 Mar 2025 15:24	SY/MD	Not Ok
14	Q1645-03RE	VY021668.D	27 Mar 2025 15:48	SY/MD	Confirms
15	Q1662-03	VY021669.D	27 Mar 2025 16:11	SY/MD	Not Ok
16	Q1662-07	VY021670.D	27 Mar 2025 16:35	SY/MD	ReRun
17	Q1661-02	VY021671.D	27 Mar 2025 16:58	SY/MD	Not Ok
18	Q1650-02	VY021672.D	27 Mar 2025 17:21	SY/MD	Not Ok
19	Q1654-01	VY021673.D	27 Mar 2025 17:45	SY/MD	Ok
20	Q1656-01	VY021674.D	27 Mar 2025 18:08	SY/MD	Not Ok
21	Q1648-09	VY021675.D	27 Mar 2025 18:32	SY/MD	Ok

Instrument ID: **MSVOA_Y**

Daily Analysis Runlog For Sequence/QCBatch ID # VY032725

Review By	Maresh Dadoda	Review On	3/28/2025 3:29:57 PM		
Supervise By	Semsettin Yesilyurt	Supervise On	3/28/2025 3:30:48 PM		
SubDirectory	VY032725	HP Acquire Method	MSVOA_Y	HP Processing Method	82y032725s.m
STD. NAME	STD REF.#				
Tune/Reschk	VP133499				
Initial Calibration Stds	VP133500,VP133501,VP133502,VP133503,VP133505,VP133508				
CCC	VP133510				
Internal Standard/PEM	VP131783				
ICV/I.BLK	VP133509				
Surrogate Standard					
MS/MSD Standard					
LCS Standard					

22	Q1648-07	VY021676.D	27 Mar 2025 18:55	SY/MD	Ok
23	Q1648-05	VY021677.D	27 Mar 2025 19:19	SY/MD	Ok
24	Q1648-01	VY021678.D	27 Mar 2025 19:42	SY/MD	Ok
25	Q1648-03	VY021679.D	27 Mar 2025 20:05	SY/MD	Ok
26	VSTDCCC050	VY021680.D	27 Mar 2025 20:28	SY/MD	Ok,M

M : Manual Integration

Instrument ID: **MSVOA_Y**

Daily Analysis Runlog For Sequence/QC Batch ID # VY041125

Review By	Mahesh Dadoda	Review On	4/14/2025 3:42:48 PM		
Supervise By	Semsettin Yesilyurt	Supervise On	4/14/2025 3:45:15 PM		
SubDirectory	VY041125	HP Acquire Method	MSVOA_Y	HP Processing Method	82y032725s.m
STD. NAME		STD REF.#			
Tune/Reschk Initial Calibration Stds		VP133637			
CCC		VP133638,VP133639			
Internal Standard/PEM		VP131783			
ICV/I.BLK					
Surrogate Standard					
MS/MSD Standard					
LCS Standard					

Sr#	SampleId	Data File Name	Date-Time	Operator	Status
1	BFB	VY021849.D	11 Apr 2025 08:46	SY/MD	Ok
2	VSTDCCC050	VY021850.D	11 Apr 2025 09:18	SY/MD	Ok,M
3	VY0411SBL01	VY021851.D	11 Apr 2025 11:35	SY/MD	Ok
4	VY0411SBS01	VY021852.D	11 Apr 2025 12:06	SY/MD	Ok,M
5	VY0411SBSD01	VY021853.D	11 Apr 2025 12:28	SY/MD	Ok,M
6	Q1756-03	VY021854.D	11 Apr 2025 13:10	SY/MD	ReRun
7	Q1761-01	VY021855.D	11 Apr 2025 13:33	SY/MD	Ok
8	Q1781-03	VY021856.D	11 Apr 2025 13:57	SY/MD	Ok
9	Q1781-07	VY021857.D	11 Apr 2025 14:20	SY/MD	Ok
10	Q1781-11	VY021858.D	11 Apr 2025 14:44	SY/MD	Ok
11	Q1768-01	VY021859.D	11 Apr 2025 15:07	SY/MD	ReRun
12	Q1779-07	VY021860.D	11 Apr 2025 15:31	SY/MD	Ok
13	Q1783-01	VY021861.D	11 Apr 2025 15:54	SY/MD	ReRun
14	Q1784-01	VY021862.D	11 Apr 2025 16:18	SY/MD	ReRun
15	VSTDCCC050	VY021863.D	11 Apr 2025 16:40	SY/MD	Ok,M

M : Manual Integration

Instrument ID: MSVOA_Y

Daily Analysis Runlog For Sequence/QC Batch ID # VY032725

Review By	Mahesh Dadoda	Review On	3/28/2025 3:29:57 PM		
Supervise By	Semsettin Yesilyurt	Supervise On	3/28/2025 3:30:48 PM		
SubDirectory	VY032725	HP Acquire Method	MSVOA_Y	HP Processing Method	82y032725s.m
STD. NAME		STD REF.#			
Tune/Reschk		VP133499			
Initial Calibration Stds		VP133500,VP133501,VP133502,VP133503,VP133505,VP133508			
CCC		VP133510			
Internal Standard/PEM		VP131783			
ICV/I.BLK		VP133509			
Surrogate Standard					
MS/MSD Standard					
LCS Standard					

Sr#	SampleID	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	BFB	BFB	VY021655.D	27 Mar 2025 09:23		SY/MD	Ok
2	VSTDIC005	VSTDIC005	VY021656.D	27 Mar 2025 10:08		SY/MD	Ok,M
3	VSTDIC010	VSTDIC010	VY021657.D	27 Mar 2025 10:31		SY/MD	Ok,M
4	VSTDIC020	VSTDIC020	VY021658.D	27 Mar 2025 10:54	Comp.#20 is on Linear Regression	SY/MD	Ok,M
5	VSTDIC050	VSTDIC050	VY021659.D	27 Mar 2025 11:16		SY/MD	Ok,M
6	VSTDIC100	VSTDIC100	VY021660.D	27 Mar 2025 11:39		SY/MD	Ok,M
7	VSTDIC150	VSTDIC150	VY021661.D	27 Mar 2025 12:02		SY/MD	Ok,M
8	VIBLK	VIBLK	VY021662.D	27 Mar 2025 12:35		SY/MD	Ok
9	VSTDICV050	ICVVY032725	VY021663.D	27 Mar 2025 13:15		SY/MD	Ok,M
10	VY0327SBL01	VY0327SBL01	VY021664.D	27 Mar 2025 13:59		SY/MD	Ok
11	VY0327SBS01	VY0327SBS01	VY021665.D	27 Mar 2025 14:38		SY/MD	Ok,M
12	VY0327SBS01	VY0327SBS01	VY021666.D	27 Mar 2025 15:00		SY/MD	Ok,M
13	Q1661-02	351	VY021667.D	27 Mar 2025 15:24	vial-B Not purge	SY/MD	Not Ok
14	Q1645-03RE	TP-4-VOCRE	VY021668.D	27 Mar 2025 15:48	vial-B Internal Standard Fail	SY/MD	Confirms
15	Q1662-03	TP-6-VOC	VY021669.D	27 Mar 2025 16:11	vial-A NOT PURGE	SY/MD	Not Ok
16	Q1662-07	TP-7-VOC	VY021670.D	27 Mar 2025 16:35	vial-A Internal Standard Fail	SY/MD	ReRun
17	Q1661-02	351	VY021671.D	27 Mar 2025 16:58	vial-A NOT PURGE	SY/MD	Not Ok

Instrument ID: MSVOA_Y

Daily Analysis Runlog For Sequence/QC Batch ID # VY032725

Review By	Mahesh Dadoda	Review On	3/28/2025 3:29:57 PM		
Supervise By	Semsettin Yesilyurt	Supervise On	3/28/2025 3:30:48 PM		
SubDirectory	VY032725	HP Acquire Method	MSVOA_Y	HP Processing Method	82y032725s.m
STD. NAME	STD REF.#				
Tune/Reschk	VP133499				
Initial Calibration Stds	VP133500,VP133501,VP133502,VP133503,VP133505,VP133508				
CCC	VP133510				
Internal Standard/PEM	VP131783				
ICV/I.BLK	VP133509				
Surrogate Standard					
MS/MSD Standard					
LCS Standard					

18	Q1650-02	STOCK-PILE-BIN2	VY021672.D	27 Mar 2025 17:21	vial-A NOT PURGE	SY/MD	Not Ok
19	Q1654-01	RT3407	VY021673.D	27 Mar 2025 17:45	vial-A Internal Standard Fail	SY/MD	Ok
20	Q1656-01	VNJ239	VY021674.D	27 Mar 2025 18:08	vial-A NOT PURGE	SY/MD	Not Ok
21	Q1648-09	TP-7	VY021675.D	27 Mar 2025 18:32	vial-A	SY/MD	Ok
22	Q1648-07	TP-6	VY021676.D	27 Mar 2025 18:55	vial-A	SY/MD	Ok
23	Q1648-05	TP-3	VY021677.D	27 Mar 2025 19:19	vial-A	SY/MD	Ok
24	Q1648-01	TP-4	VY021678.D	27 Mar 2025 19:42	vial-A	SY/MD	Ok
25	Q1648-03	TP-5	VY021679.D	27 Mar 2025 20:05	vial-A	SY/MD	Ok
26	VSTDCCC050	VSTDCCC050EC	VY021680.D	27 Mar 2025 20:28		SY/MD	Ok,M

M : Manual Integration

Instrument ID: MSVOA_Y

Daily Analysis Runlog For Sequence/QC Batch ID # VY041125

Review By	Mahesh Dadoda	Review On	4/14/2025 3:42:48 PM		
Supervise By	Semsettin Yesilyurt	Supervise On	4/14/2025 3:45:15 PM		
SubDirectory	VY041125	HP Acquire Method	MSVOA_Y	HP Processing Method	82y032725s.m
STD. NAME		STD REF.#			
Tune/Reschk Initial Calibration Stds		VP133637			
CCC		VP133638,VP133639			
Internal Standard/PEM		VP131783			
ICV/I.BLK					
Surrogate Standard					
MS/MSD Standard					
LCS Standard					

Sr#	Sampleld	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	BFB	BFB	VY021849.D	11 Apr 2025 08:46		SY/MD	Ok
2	VSTDCCC050	VSTDCCC050	VY021850.D	11 Apr 2025 09:18		SY/MD	Ok,M
3	VY0411SBL01	VY0411SBL01	VY021851.D	11 Apr 2025 11:35		SY/MD	Ok
4	VY0411SBS01	VY0411SBS01	VY021852.D	11 Apr 2025 12:06		SY/MD	Ok,M
5	VY0411SBSD01	VY0411SBSD01	VY021853.D	11 Apr 2025 12:28		SY/MD	Ok,M
6	Q1756-03	TP-9-VOC	VY021854.D	11 Apr 2025 13:10	vial-A Internal Standard Fail ;Surrogate fail	SY/MD	ReRun
7	Q1761-01	GST3	VY021855.D	11 Apr 2025 13:33	vial-A	SY/MD	Ok
8	Q1781-03	WC-13-VOC	VY021856.D	11 Apr 2025 13:57	vial-A	SY/MD	Ok
9	Q1781-07	WC-12-VOC	VY021857.D	11 Apr 2025 14:20	vial-A	SY/MD	Ok
10	Q1781-11	WC-11-VOC	VY021858.D	11 Apr 2025 14:44	vial-A	SY/MD	Ok
11	Q1768-01	CRUSHED STONE	VY021859.D	11 Apr 2025 15:07	vial-A Internal Standard Fail ;Surrogate fail	SY/MD	ReRun
12	Q1779-07	TP-23-VOC	VY021860.D	11 Apr 2025 15:31	vial-A	SY/MD	Ok
13	Q1783-01	SU-03-41125	VY021861.D	11 Apr 2025 15:54	vial-A Internal Standard Fail	SY/MD	ReRun
14	Q1784-01	OR-01-41125	VY021862.D	11 Apr 2025 16:18	vial-A Internal Standard Fail	SY/MD	ReRun
15	VSTDCCC050	VSTDCCC050EC	VY021863.D	11 Apr 2025 16:40		SY/MD	Ok,M

M : Manual Integration



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LAB CHRONICLE

OrderID:	Q1761	OrderDate:	4/9/2025 2:47:00 PM
Client:	G Environmental	Project:	Stockton
Contact:	Gary Landis	Location:	L31,VOA Ref. #2 Soil

LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
Q1761-01	GST3	SOIL	VOC-TCLVOA-10	8260D	04/08/25		04/11/25	04/09/25