

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070825\
Data File : VX046917.D
Acq On : 08 Jul 2025 13:45
Operator : JC/MD
Sample : IBLK
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
IBLK

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_X\Method\82X070225W.M

Title : SW846 8260

Signal : TIC: VX046917.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.385	693	705	722	rBV3	297036	944027	33.89%	5.086%
2	5.550	722	732	758	rVB	463376	1460809	52.45%	7.870%
3	5.958	786	799	819	rBV	329260	961875	34.53%	5.182%
4	6.757	921	930	956	rBV	778822	2028245	72.82%	10.927%
5	8.647	1233	1240	1259	rBV	1741176	2785246	100.00%	15.005%
6	9.574	1384	1392	1414	rBV	733270	1610569	57.83%	8.677%
7	10.049	1464	1470	1486	rBV	1771543	2441094	87.64%	13.151%
8	11.079	1633	1639	1652	rBV	1472939	1861502	66.83%	10.029%
9	11.384	1683	1689	1717	rBV	1352750	2127561	76.39%	11.462%
10	12.018	1788	1793	1809	rVB	1872305	2340736	84.04%	12.611%

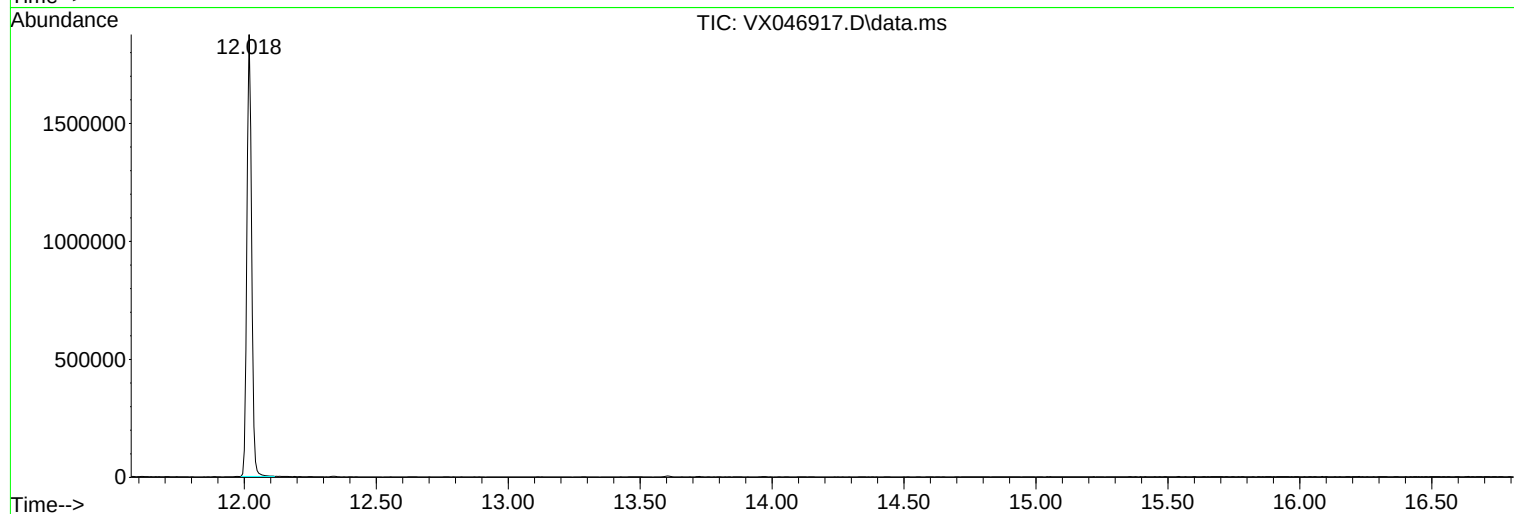
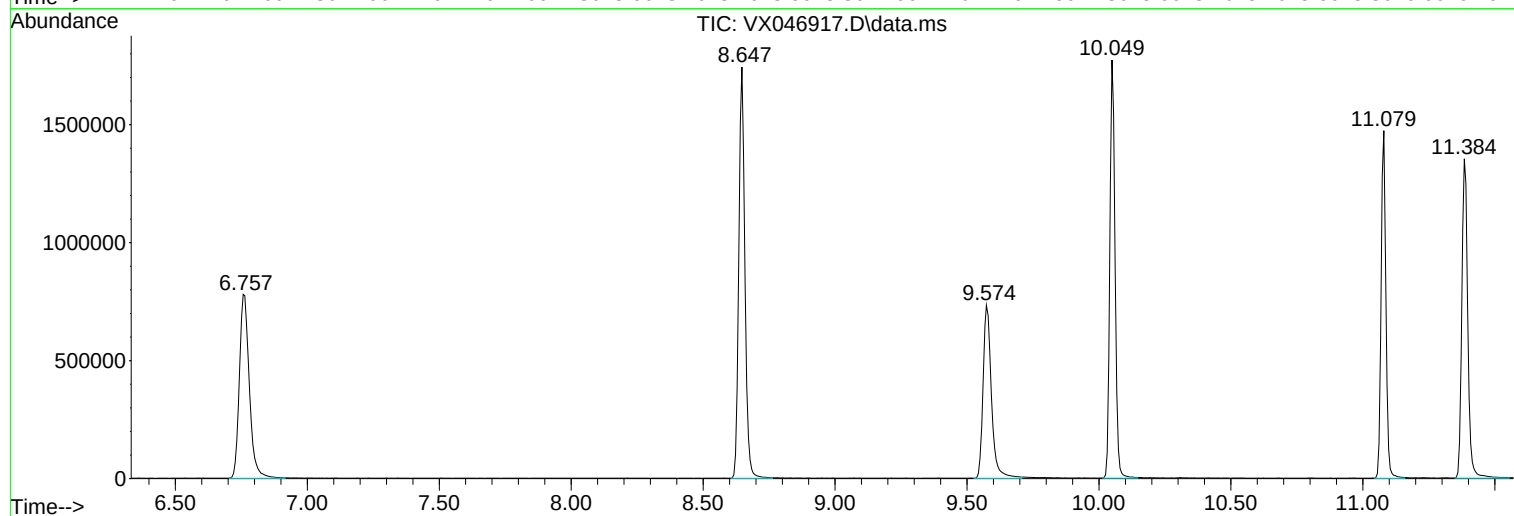
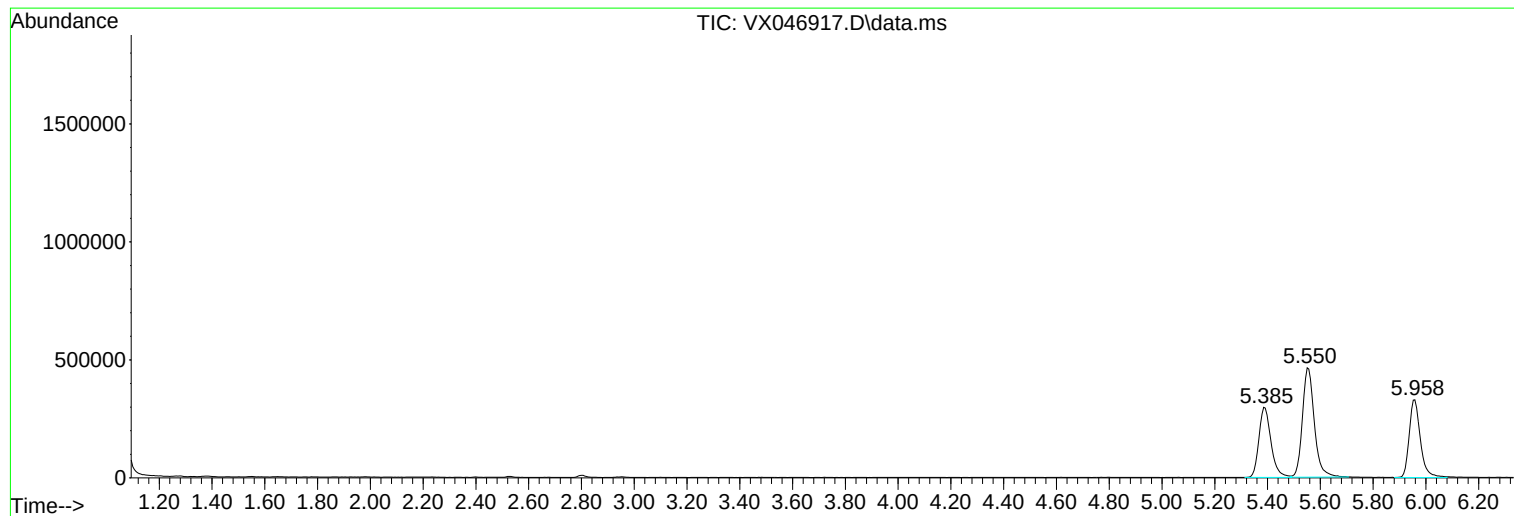
Sum of corrected areas: 18561664

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X070225W.M
Quant Title : SW846 8260

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



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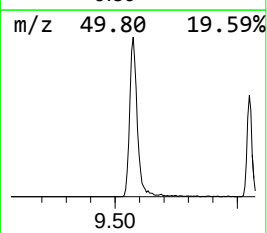
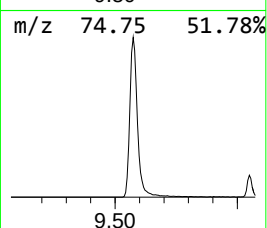
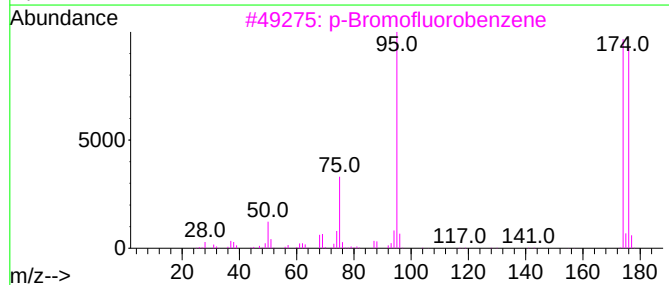
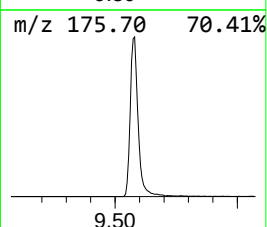
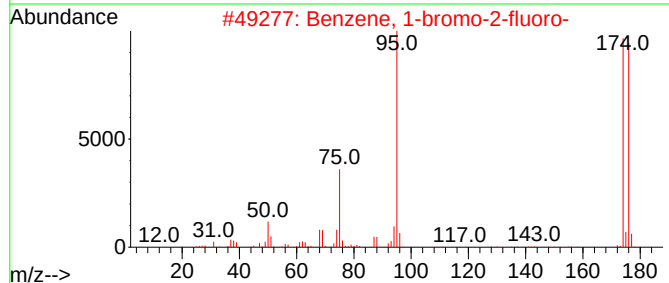
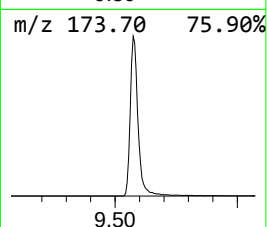
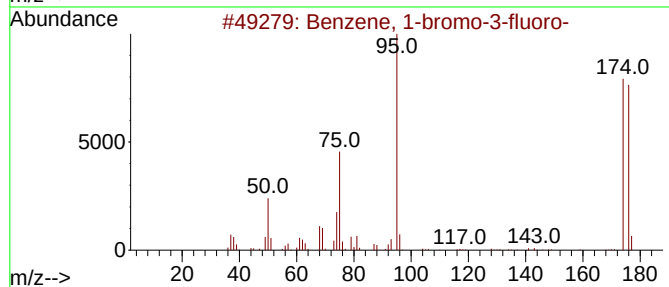
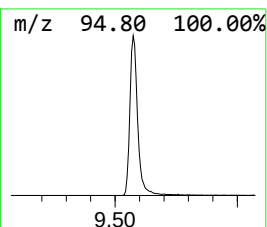
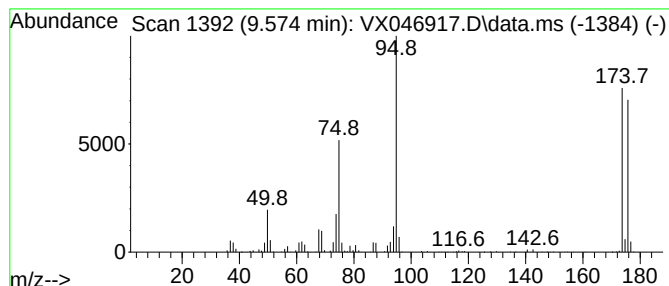
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X070225W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-bromo-3-fluoro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.574	32.99 ug/l	1610570	Chlorobenzene-d5	10.049

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-3-fluoro-	174	C6H4BrF	001073-06-9	97
2			Benzene, 1-bromo-2-fluoro-	174	C6H4BrF	001072-85-1	94
3			p-Bromofluorobenzene	174	C6H4BrF	000460-00-4	87
4			4(1H)-Pyrimidinone, 5-bromo-	174	C4H3BrN2O	019808-30-1	17
5			Benzaldehyde, 2,6-difluoro-3,4-d...	174	C7H4F2O3	152434-98-5	14



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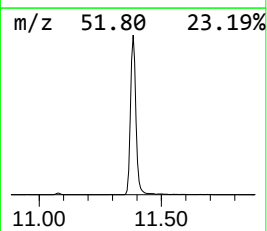
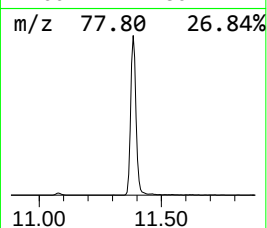
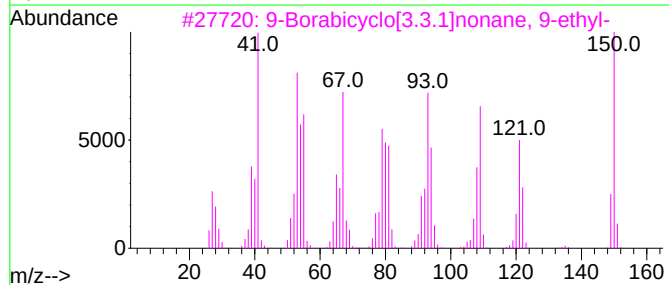
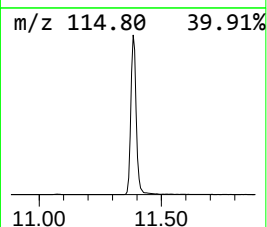
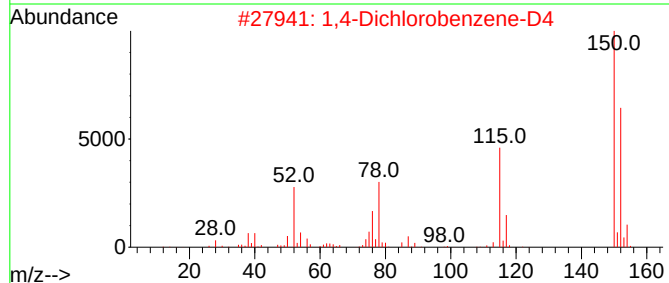
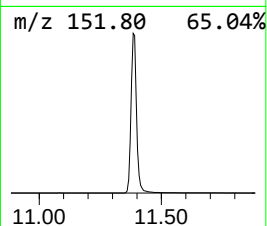
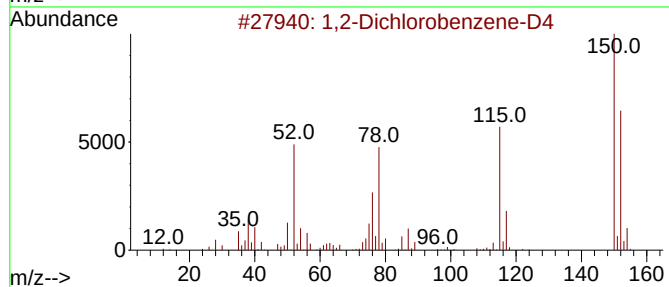
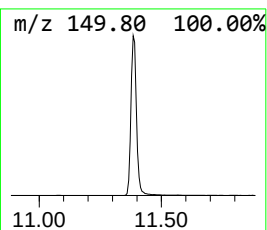
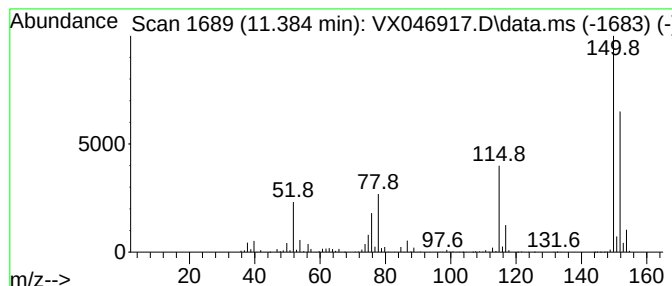
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Peak Number 2 1,2-Dichlorobenzene-D4 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.384	45.45 ug/l	2127560	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Dichlorobenzene-D4	150	C6D4Cl2	002199-69-1	91
2			1,4-Dichlorobenzene-D4	150	C6D4Cl2	003855-82-1	70
3			9-Borabicyclo[3.3.1]nonane, 9-ethyl-	150	C10H19B	052102-17-7	38
4			4-(1H-Pyrrol-1-yl)-1,2,5-oxadiazole	150	C6H6N4O	618428-52-7	28
5			1H,3H,5H,6H,7H-Cyclopenta[d]pyridine	152	C7H8N2O2	005466-00-2	16



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-brom...	9.574	33.0	ug/l	1610570	3	10.049	2441090	50.0
1,2-Dichloroben...	11.384	45.5	ug/l	2127560	4	12.018	2340740	50.0