

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121224\
Data File : VX044251.D
Acq On : 12 Dec 2024 14:46
Operator : JC/MD
Sample : P5261-20 100X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
DUP-02-20241210

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\82X121124W.M

Title : SW846 8260

Signal : TIC: VX044251.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.617	67	87	88	rBV3	19621	87393	12.27%	2.332%
2	2.940	295	304	309	rBV4	3877	8860	1.24%	0.236%
3	5.379	694	704	717	rBV3	77979	236294	33.17%	6.304%
4	5.544	721	731	744	rVV3	126043	364716	51.20%	9.730%
5	5.952	789	798	816	rBV	96081	268599	37.70%	7.166%
6	6.757	919	930	951	rBV	214798	525398	73.75%	14.017%
7	8.647	1233	1240	1258	rBV	448158	712370	100.00%	19.005%
8	10.049	1465	1470	1480	rBV	420401	608237	85.38%	16.227%
9	11.079	1634	1639	1655	rVB	336206	433832	60.90%	11.574%
10	12.018	1788	1793	1805	rBV	401395	502571	70.55%	13.408%

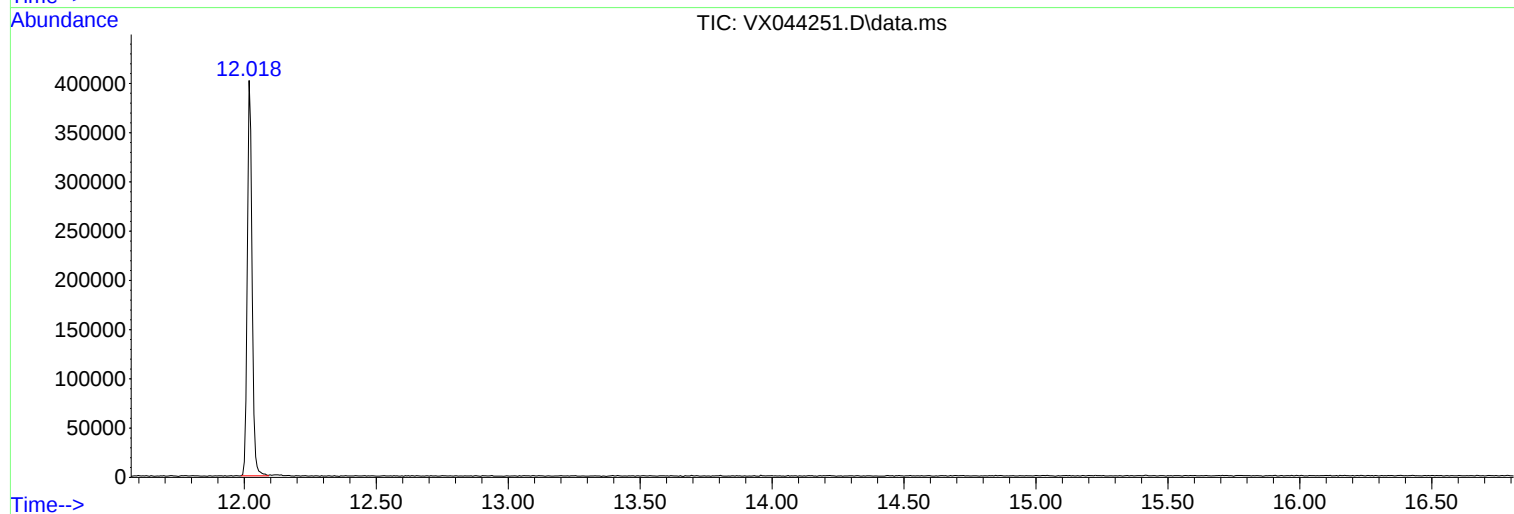
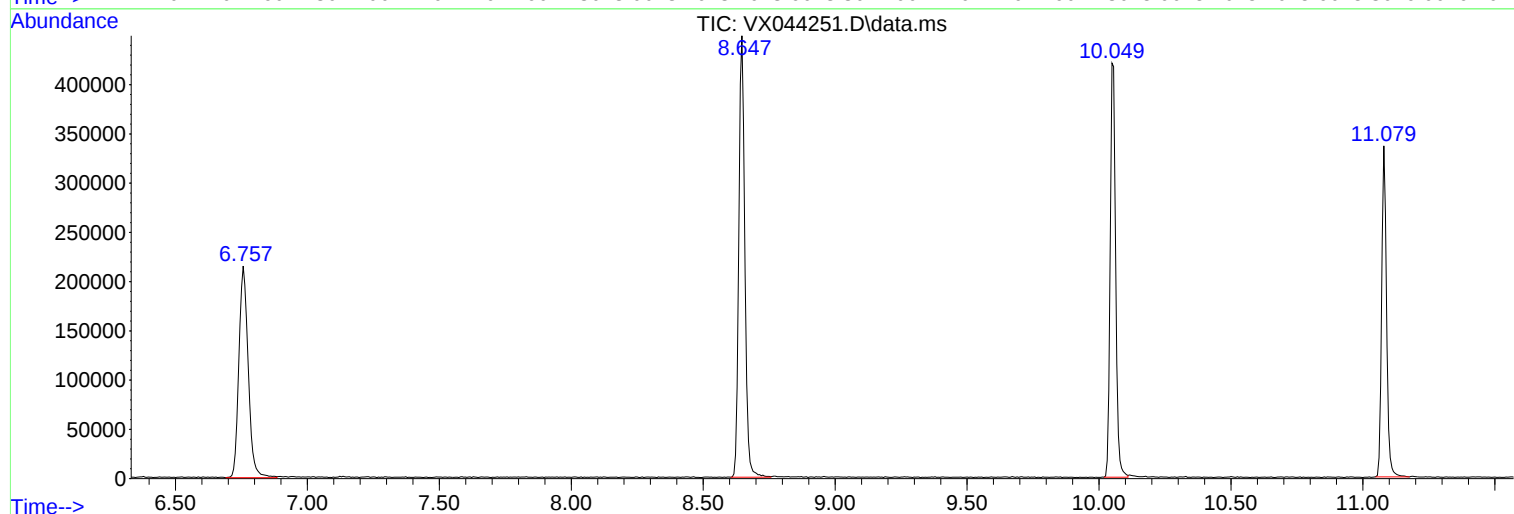
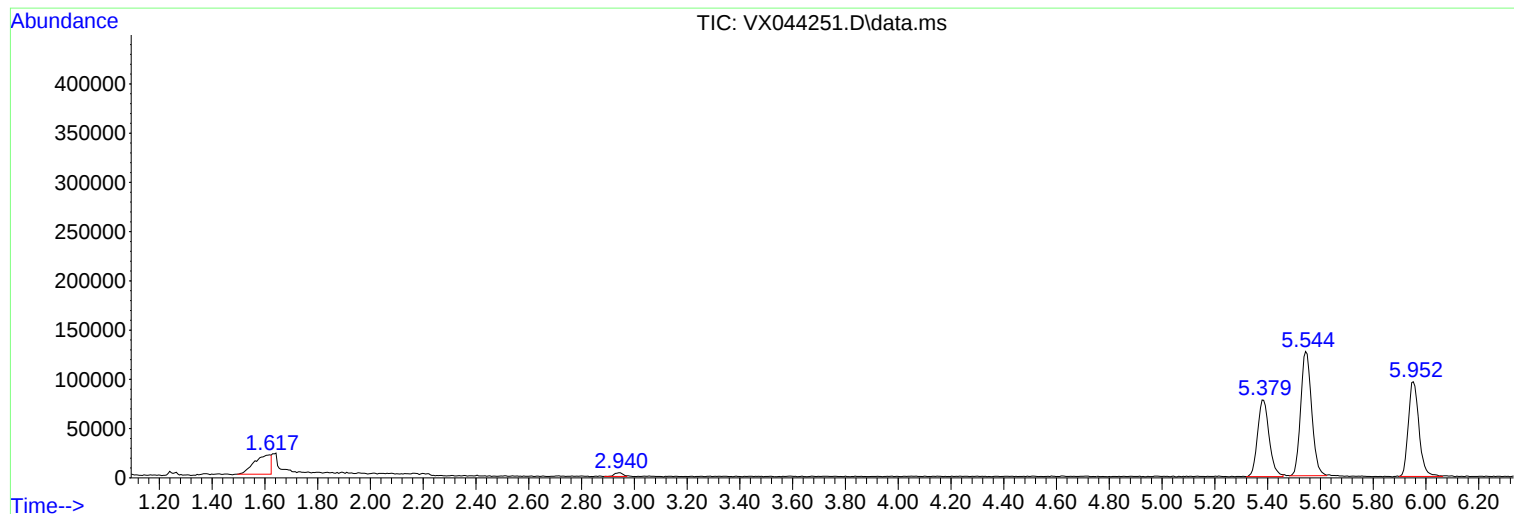
Sum of corrected areas: 3748270

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121124W.M
Quant Title : SW846 8260

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



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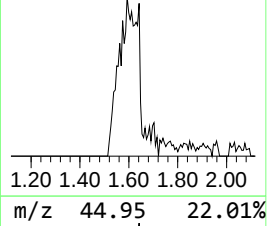
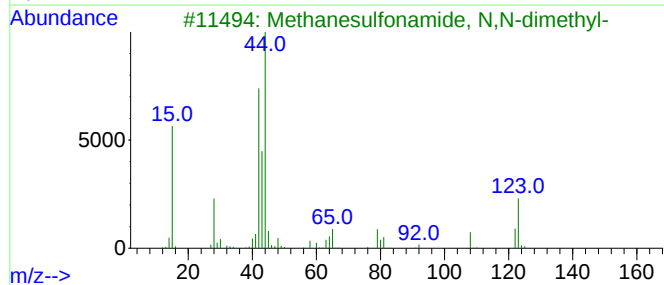
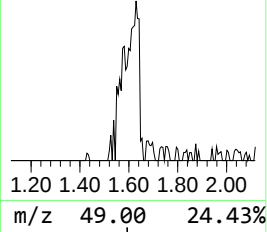
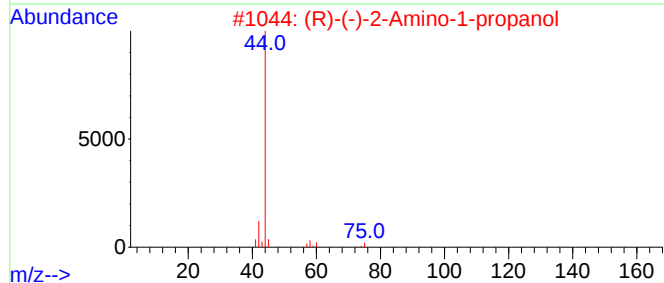
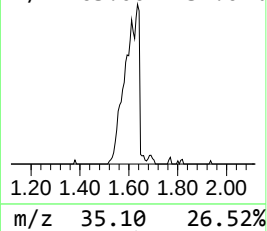
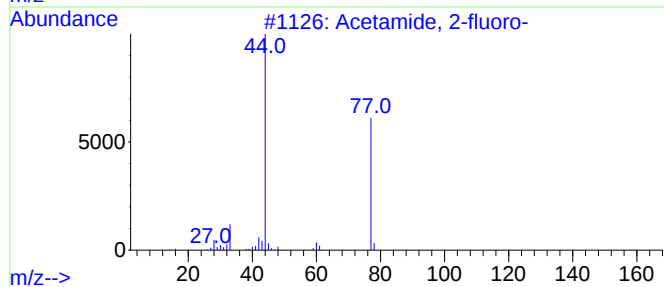
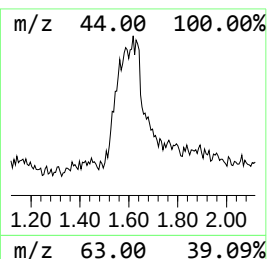
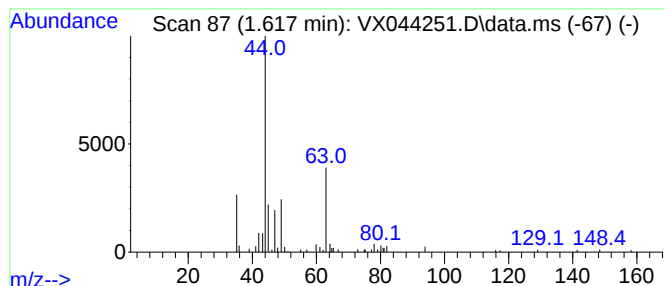
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Peak Number 1 unknown1.617 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.617	11.98 ug/l	87393	Pentafluorobenzene	5.544

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetamide, 2-fluoro-	77	C2H4FNO	000640-19-7	10
2			(R)-(-)-2-Amino-1-propanol	75	C3H9NO	035320-23-1	9
3			Methanesulfonamide, N,N-dimethyl-	123	C3H9NO2S	000918-05-8	9
4			Acetic acid, [(aminocarbonyl)ami...	132	C3H4N2O4	000585-05-7	9
5			Ethyl 2-(2-chloroacetamido)-3,3,...	263	C7H9ClF3NO4	1000224-74-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.617	1.617	12.0	ug/l	87393	1	5.544	364716	50.0