

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121224\
Data File : VX044263.D
Acq On : 12 Dec 2024 19:24
Operator : JC/MD
Sample : P5261-17
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
RE120D1-20241210

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: VX044263.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.611	69	86	88	rBV3	20162	88278	12.66%	2.319%
2	2.319	197	202	209	rVB4	4632	8789	1.26%	0.231%
3	5.385	694	705	722	rBV	78446	235348	33.75%	6.181%
4	5.544	722	731	746	rVV	123726	358223	51.37%	9.408%
5	5.952	786	798	812	rBV2	97375	264749	37.97%	6.953%
6	6.757	919	930	946	rBV	210101	512056	73.43%	13.448%
7	7.123	982	990	1003	rBV2	43464	101915	14.62%	2.677%
8	8.647	1233	1240	1255	rBV	443650	697308	100.00%	18.314%
9	10.049	1465	1470	1481	rBV	412555	598686	85.86%	15.724%
10	11.079	1634	1639	1651	rBV	317915	426425	61.15%	11.200%
11	12.018	1788	1793	1809	rBV	390681	515759	73.96%	13.546%

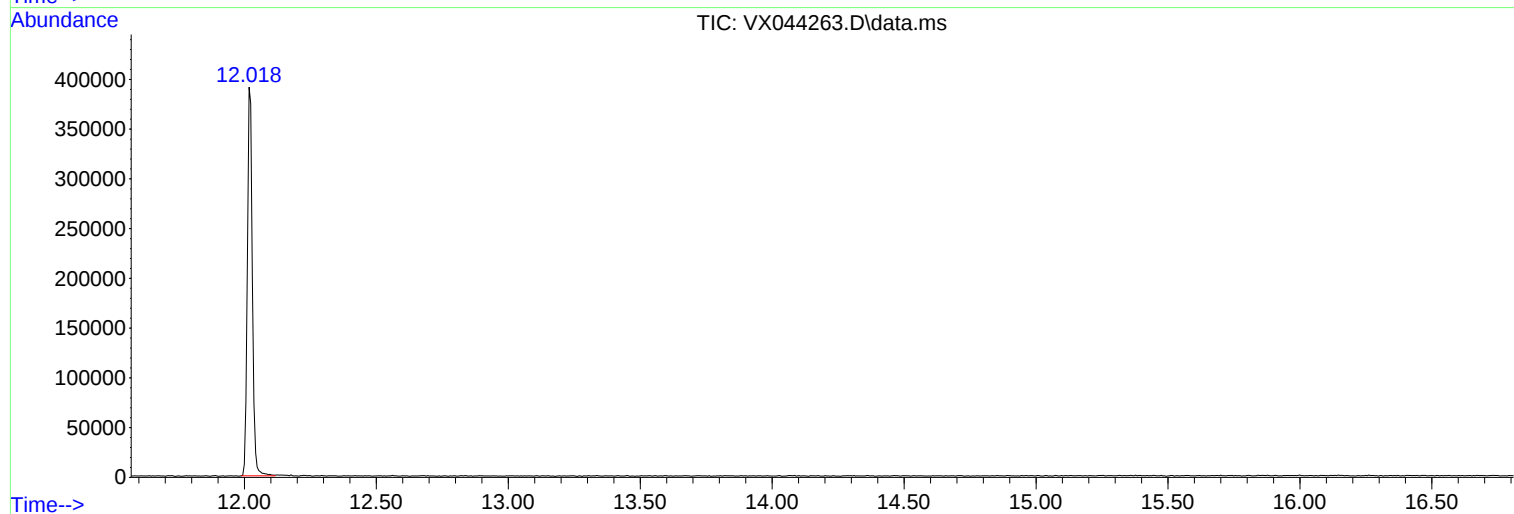
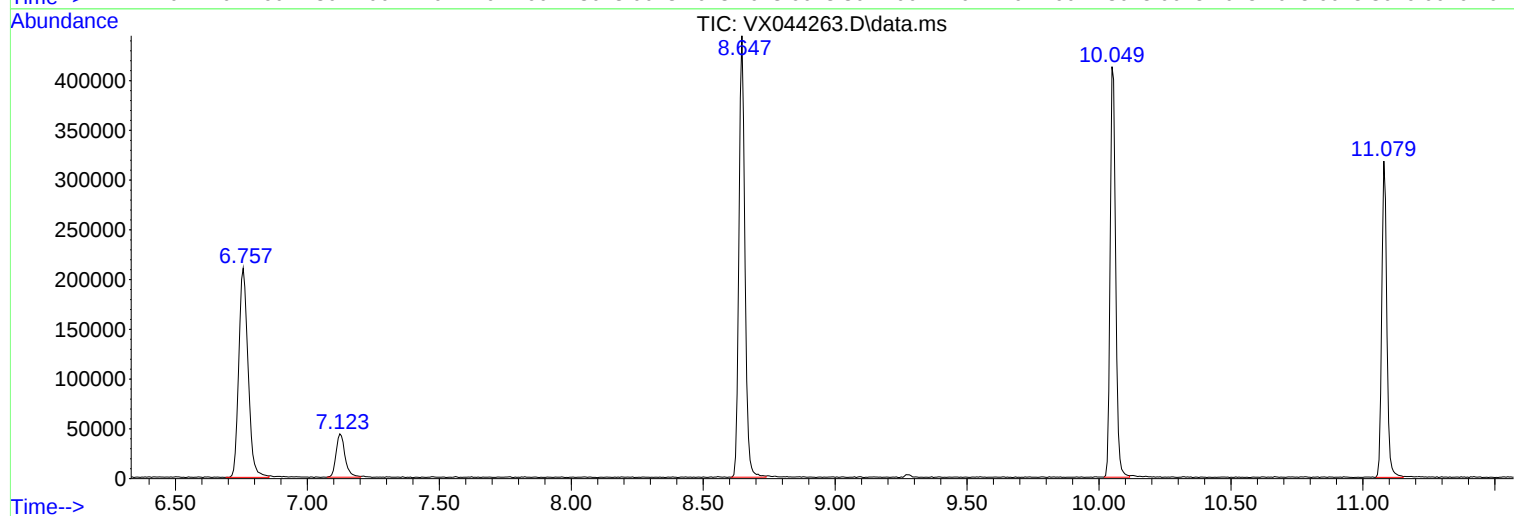
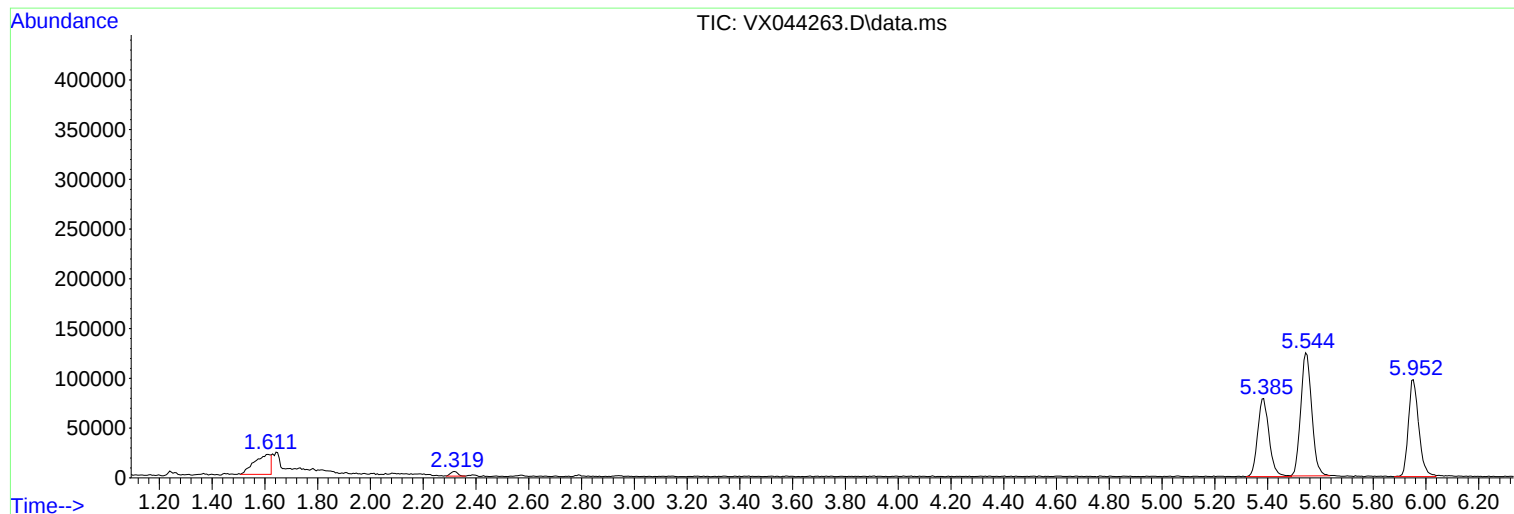
Sum of corrected areas: 3807536

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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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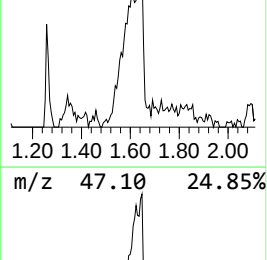
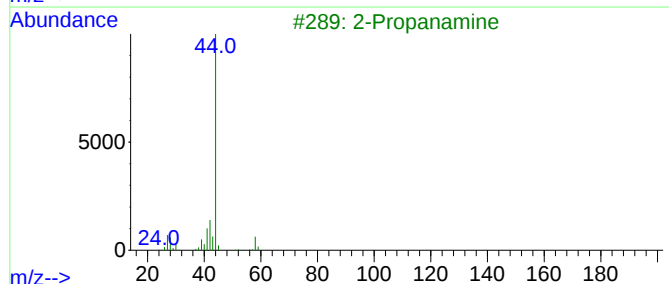
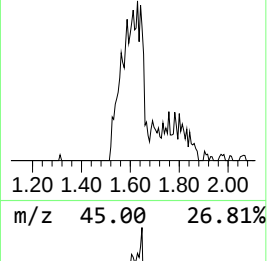
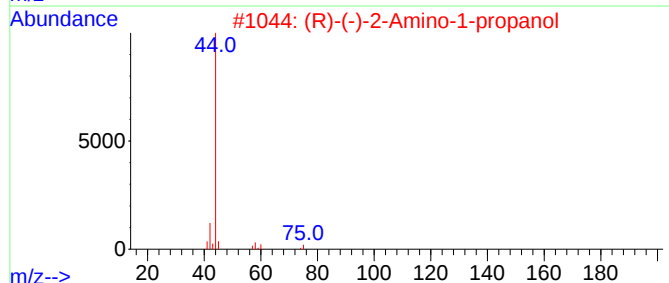
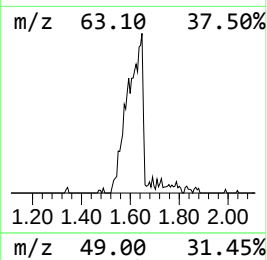
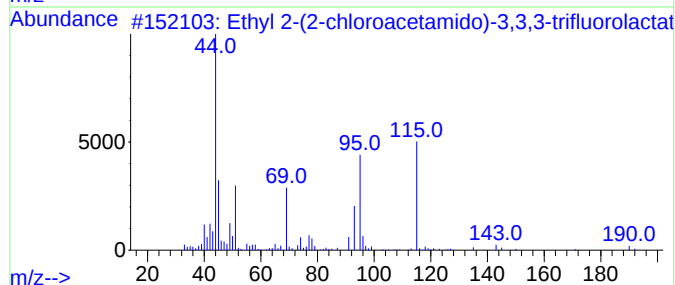
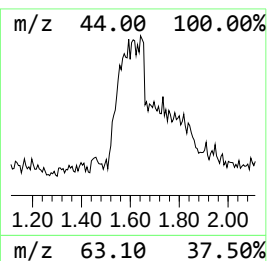
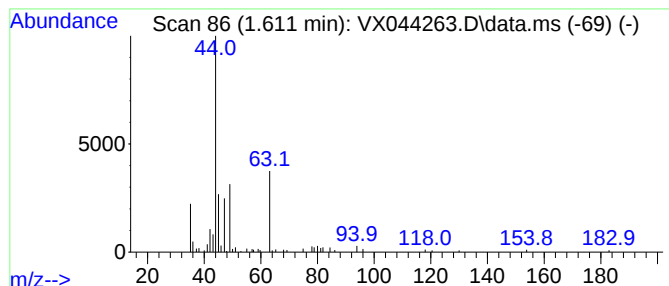
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Peak Number 1 unknown1.611 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.611	12.32 ug/l	88278	Pentafluorobenzene	5.544

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethyl 2-(2-chloroacetamido)-3,3,...	263	C7H9ClF3NO4	1000224-74-2	10
2			(R)-(-)-2-Amino-1-propanol	75	C3H9NO	035320-23-1	9
3			2-Propanamine	59	C3H9N	000075-31-0	9
4			1,2-Benzenediol, 4-(2-amino-1-hy...	183	C9H13NO3	006539-57-7	9
5			Paradrine	151	C9H13NO	000103-86-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.611	1.611	12.3	ug/l	88278	1	5.544	358223	50.0