

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122324\
 Data File : VX044478.D
 Acq On : 23 Dec 2024 14:29
 Operator : JC/MD
 Sample : P5374-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB-190A-GW-718-720

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: VX044478.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.142	6	9	12	rBV	8788	8324	1.09%	0.175%
2	1.612	65	86	87	rBV2	20442	81885	10.76%	1.725%
3	2.319	195	202	209	rBV2	34832	69340	9.11%	1.461%
4	5.379	692	704	717	rBV2	85871	262949	34.55%	5.540%
5	5.544	721	731	748	rVV2	139371	408536	53.67%	8.607%
6	5.952	789	798	813	rBV2	101885	272161	35.76%	5.734%
7	6.757	921	930	945	rBV	227377	562201	73.86%	11.844%
8	7.117	981	989	1001	rBV	263659	581325	76.38%	12.247%
9	8.647	1233	1240	1250	rBV	464059	761134	100.00%	16.035%
10	10.049	1465	1470	1485	rBV	460336	666642	87.59%	14.045%
11	11.079	1634	1639	1651	rBV	373183	483238	63.49%	10.181%
12	12.018	1788	1793	1806	rBV	471601	588879	77.37%	12.406%

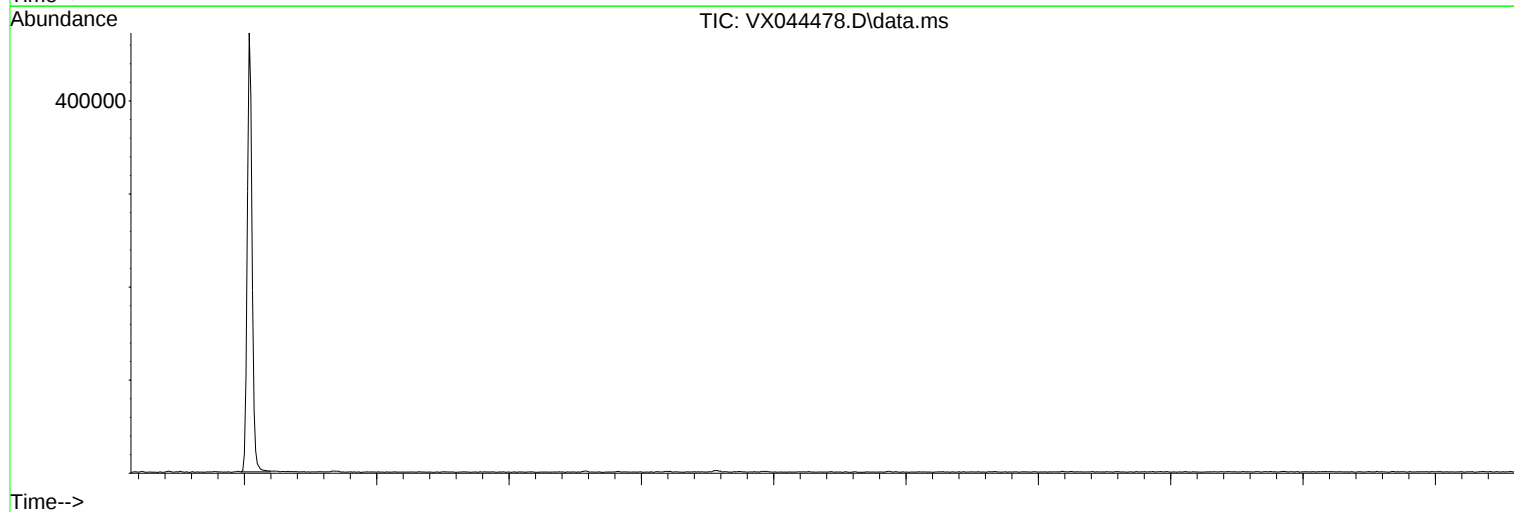
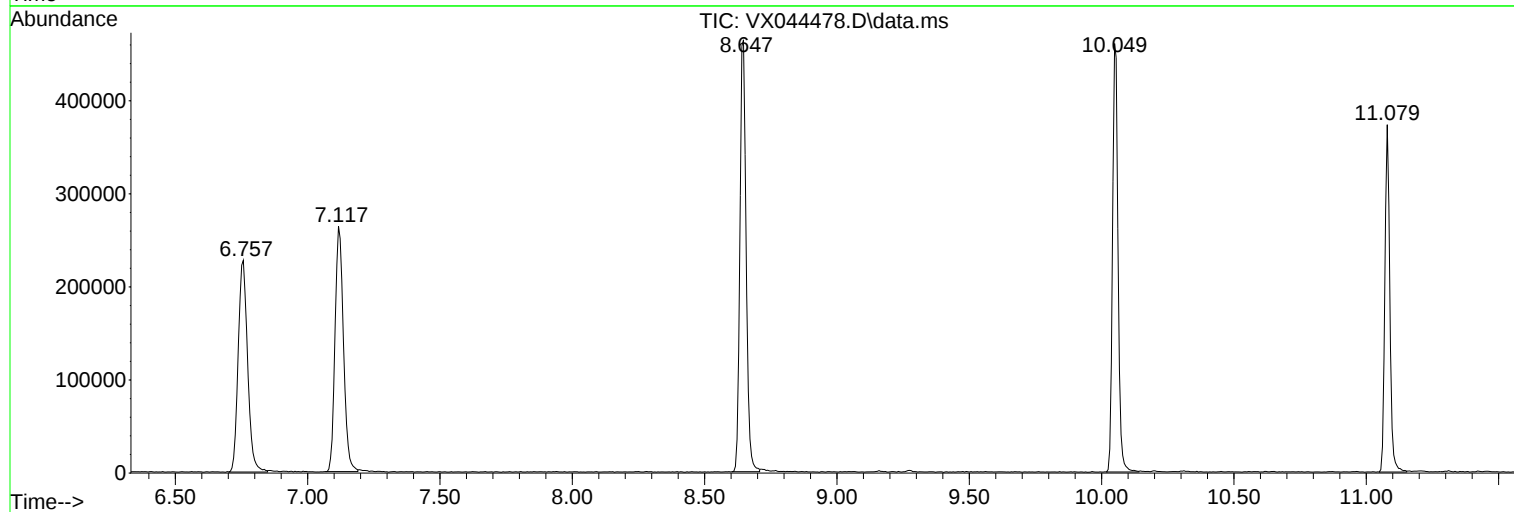
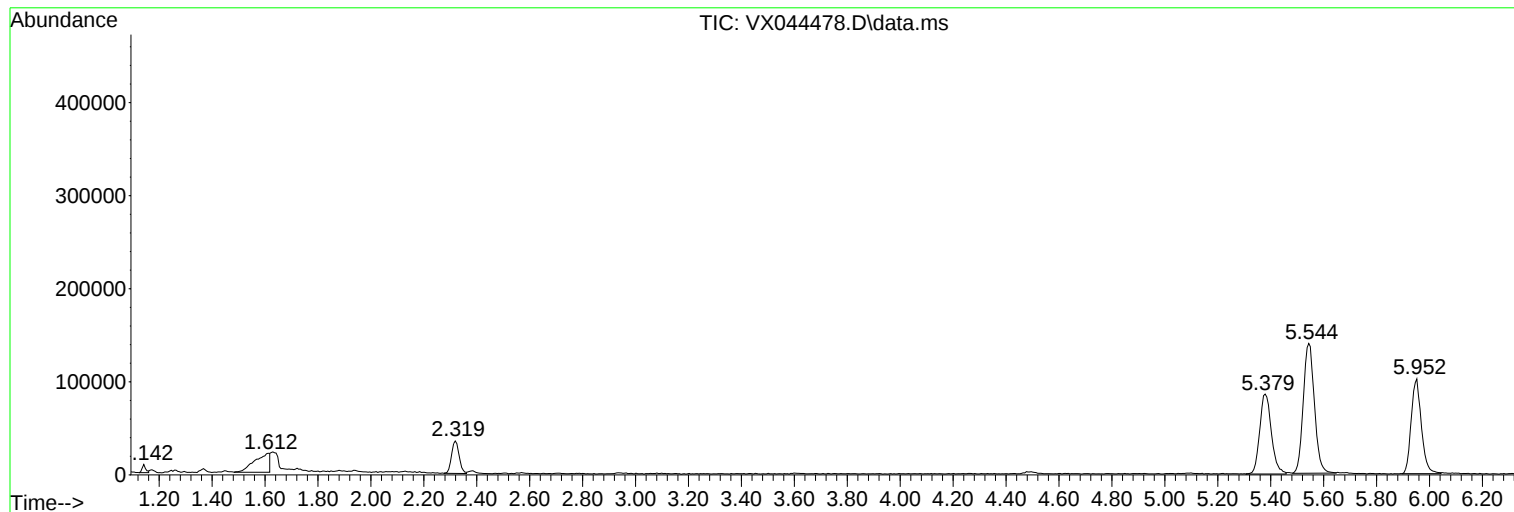
Sum of corrected areas: 4746614

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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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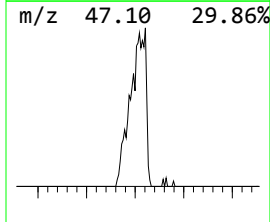
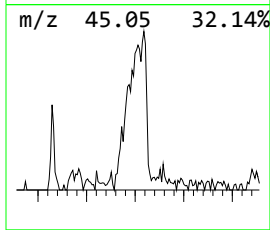
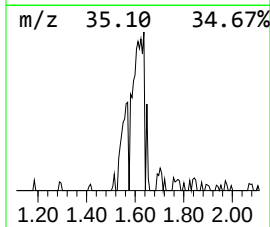
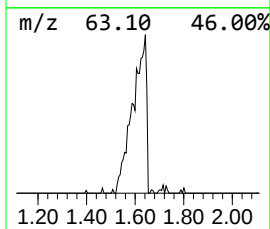
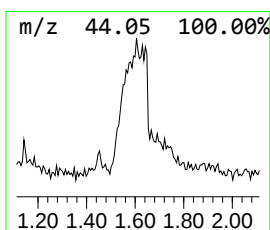
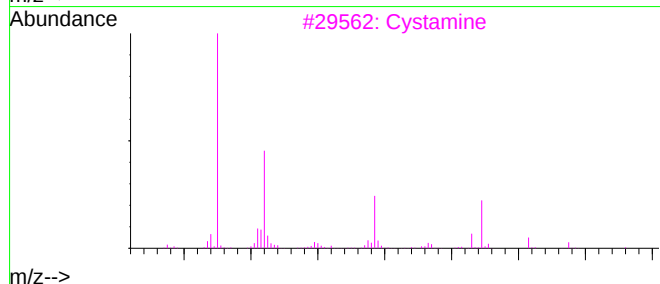
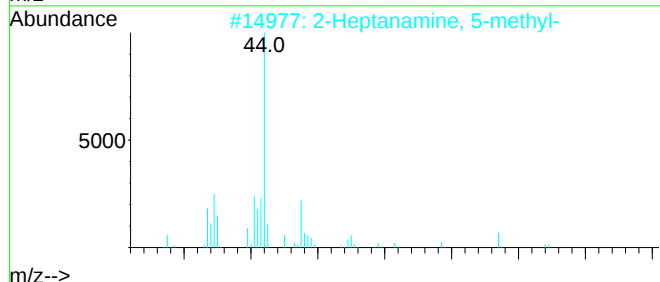
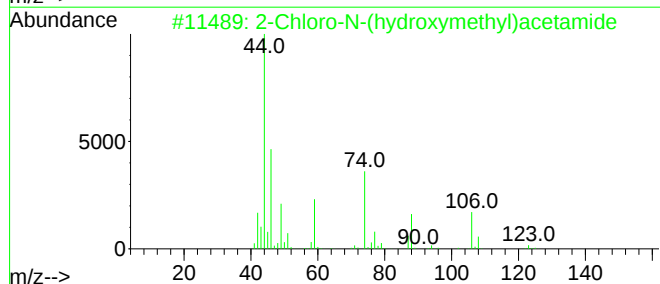
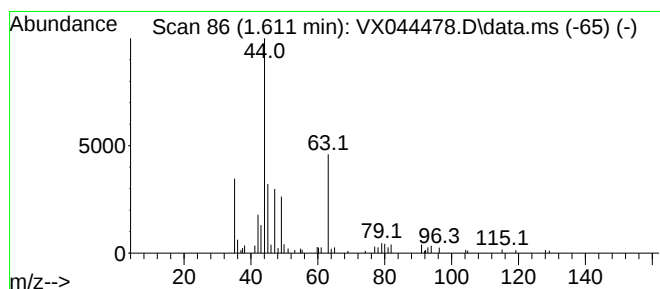
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Peak Number 1 unknown1.611 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.611	10.02 ug/l	81885	Pentafluorobenzene	5.544

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Chloro-N-(hydroxymethyl)acetamide	123	C3H6ClNO2	002832-19-1	17
2	2-Heptanamine, 5-methyl-	129	C8H19N	053907-81-6	9
3	Cystamine	152	C4H12N2S2	000051-85-4	9
4	Acetic acid, [(aminocarbonyl)ami...	132	C3H4N2O4	000585-05-7	9
5	Phenethylamine, p,.alpha.-dimethyl-	149	C10H15N	000064-11-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.611	1.611	10.0	ug/l	81885	1	5.544	408536	50.0