

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN053023\
 Data File : VN077959.D
 Acq On : 30 May 2023 10:52
 Operator : JC\MD
 Sample : VN0530MBL01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VN0530MBL01

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_N\methods\82N051523W.M

Title : SW846 8260

Signal : TIC: VN077959.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.818	140	147	156	rVB2	76874	191055	9.06%	1.828%
2	8.177	1047	1058	1062	rBV	291087	705289	33.45%	6.746%
3	8.230	1062	1067	1078	rVB	509537	1126235	53.41%	10.773%
4	8.588	1119	1128	1138	rBV	294676	681067	32.30%	6.515%
5	9.106	1207	1216	1232	rBV	759768	1529567	72.54%	14.631%
6	10.577	1457	1466	1482	rBV	1106995	2108468	100.00%	20.168%
7	11.871	1677	1686	1699	rBV	962270	1753112	83.15%	16.769%
8	12.859	1845	1854	1863	rBV	655198	1150020	54.54%	11.000%
9	13.800	2007	2014	2023	rVB	662102	1164975	55.25%	11.144%
10	15.647	2324	2328	2333	rVB3	13823	23301	1.11%	0.223%
11	15.853	2357	2363	2369	rVB5	11692	21211	1.01%	0.203%

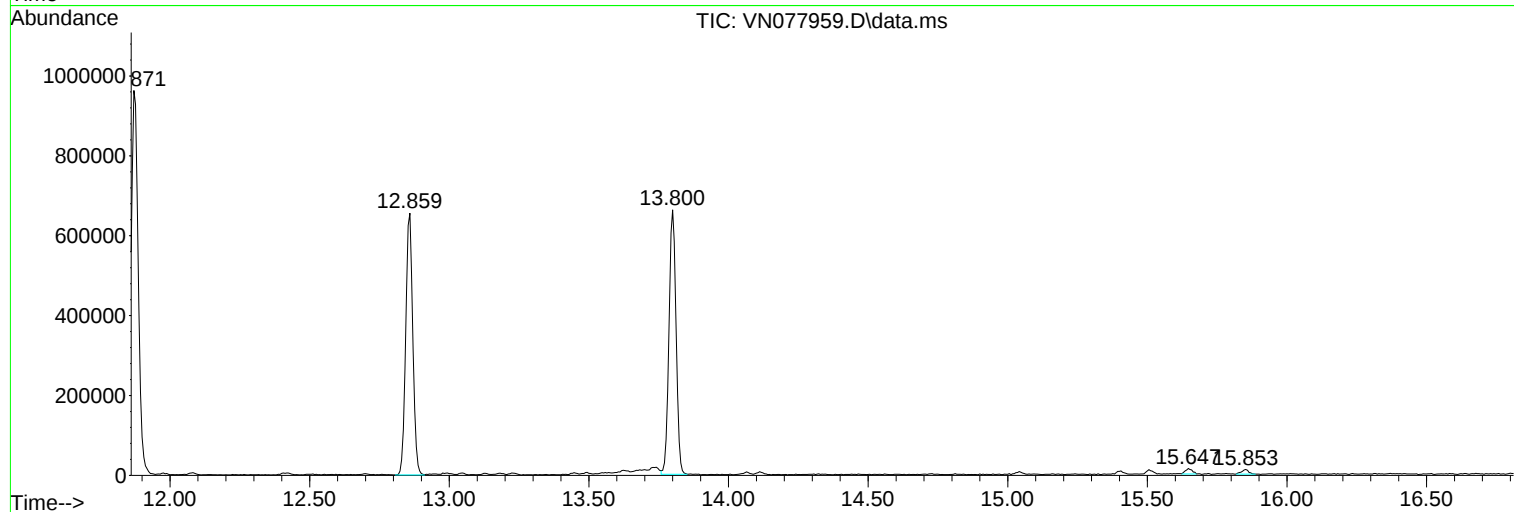
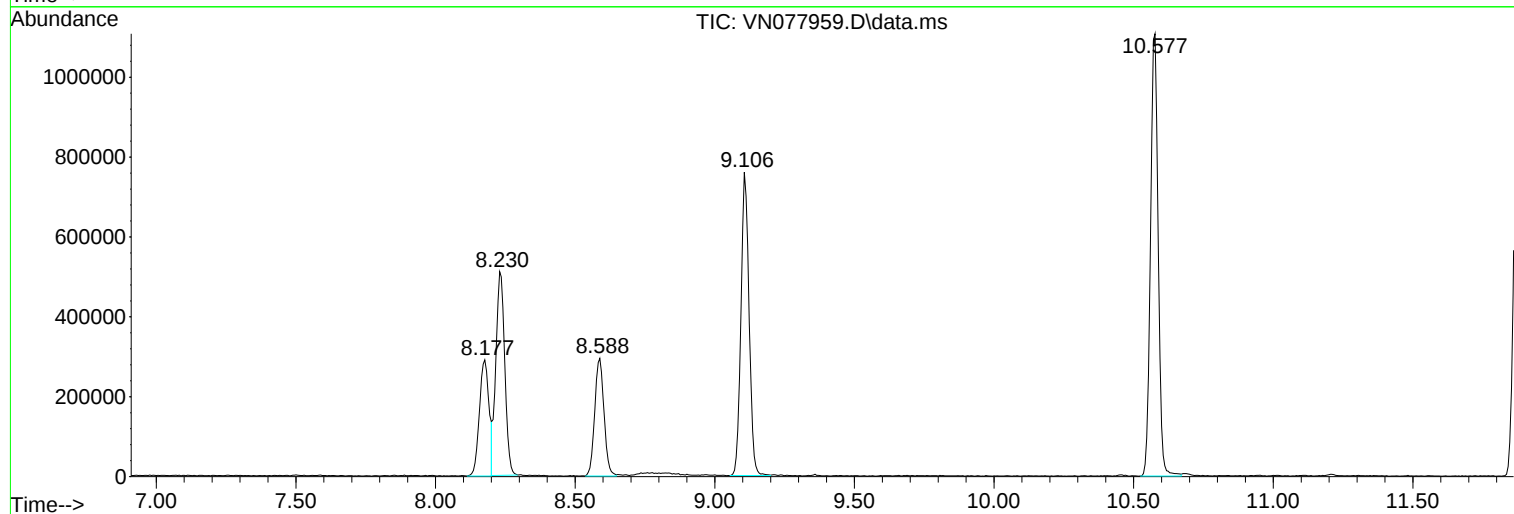
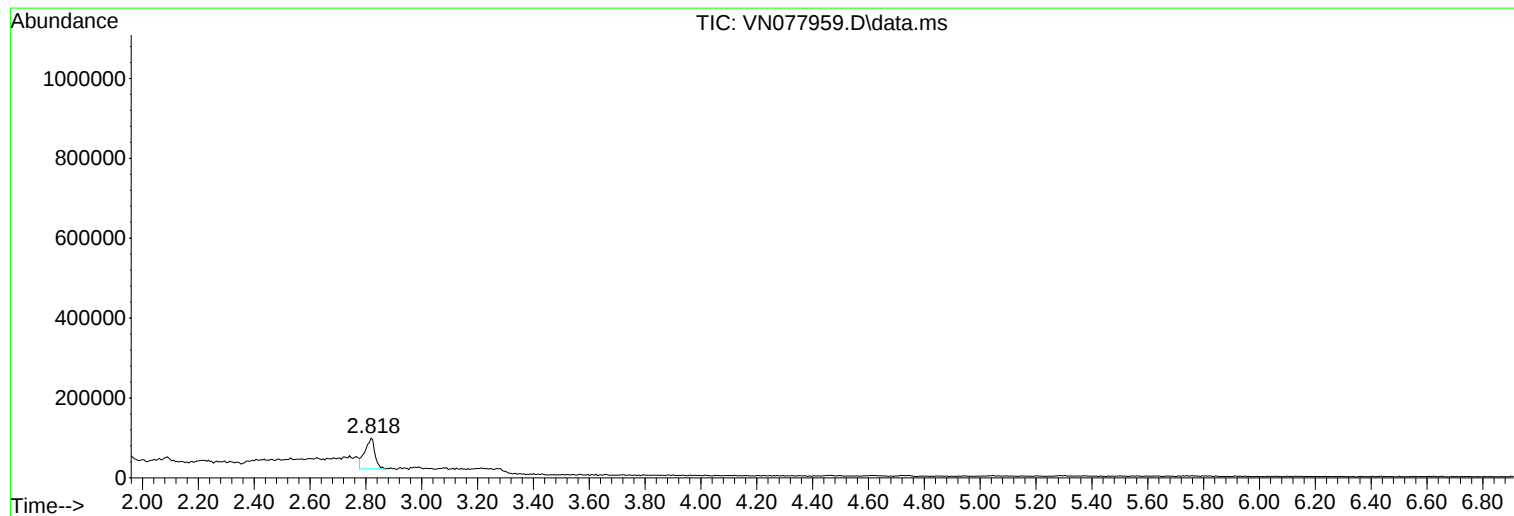
Sum of corrected areas: 10454300

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051523W.M
Quant Title : SW846 8260

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



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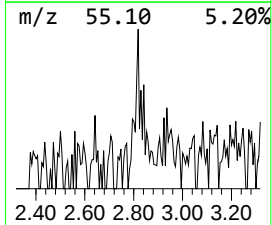
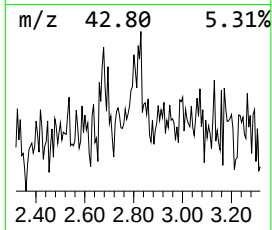
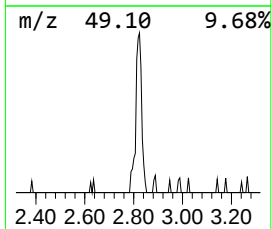
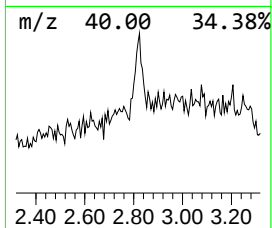
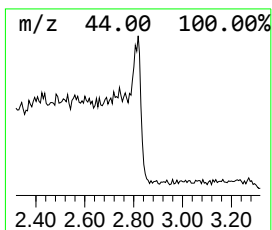
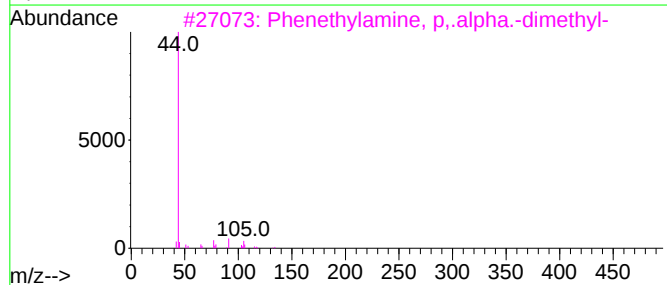
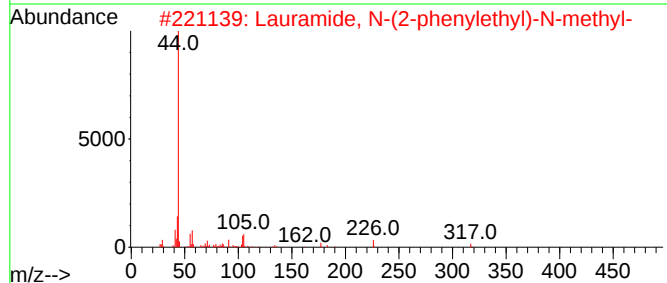
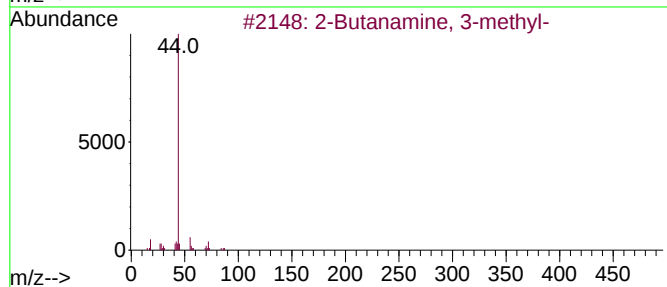
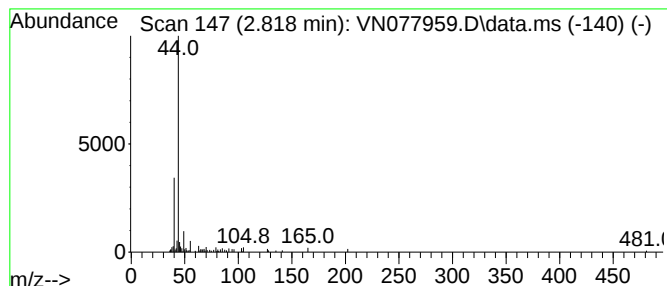
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051523W.M
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Peak Number 1 unknown2.818 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.818	8.48 ug/l	191055	Pentafluorobenzene	8.236

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanamine, 3-methyl-			87	C5H13N	000598-74-3	42
2	Lauramide, N-(2-phenylethyl)-N-m...			317	C21H35NO	1010490-83-4	38
3	Phenethylamine, p,.alpha.-dimethyl-			149	C10H15N	000064-11-9	38
4	Actinobolin			300	C13H20N2O6	024397-89-5	36
5	2-Pentanamine			87	C5H13N	063493-28-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown2.818	2.818	8.5	ug/l	191055	1	8.236	1126240	50.0