

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030624\
 Data File : VN081330.D
 Acq On : 06 Mar 2024 16:39
 Operator : JC\MD
 Sample : P1683-01 20X
 Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 1132

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\82N030524W.M

Title : SW846 8260

Signal : TIC: VN081330.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.824	139	148	155	rBV2	34376	105009	5.87%	1.114%
2	8.171	1044	1057	1061	rBV	247987	583226	32.62%	6.189%
3	8.224	1061	1066	1079	rVB	444698	991596	55.46%	10.522%
4	8.583	1116	1127	1134	rBV	281000	621617	34.77%	6.596%
5	8.665	1134	1141	1144	rBV4	18713	43362	2.43%	0.460%
6	9.106	1206	1216	1228	rBV	646892	1338688	74.87%	14.205%
7	10.571	1456	1465	1477	rBV	960061	1787985	100.00%	18.973%
8	11.865	1676	1685	1701	rBV	866902	1550711	86.73%	16.455%
9	12.853	1844	1853	1863	rBV	647308	1086348	60.76%	11.528%
10	13.794	2006	2013	2020	rBV	773650	1287350	72.00%	13.661%
11	14.953	2205	2210	2219	rBV8	12787	27863	1.56%	0.296%

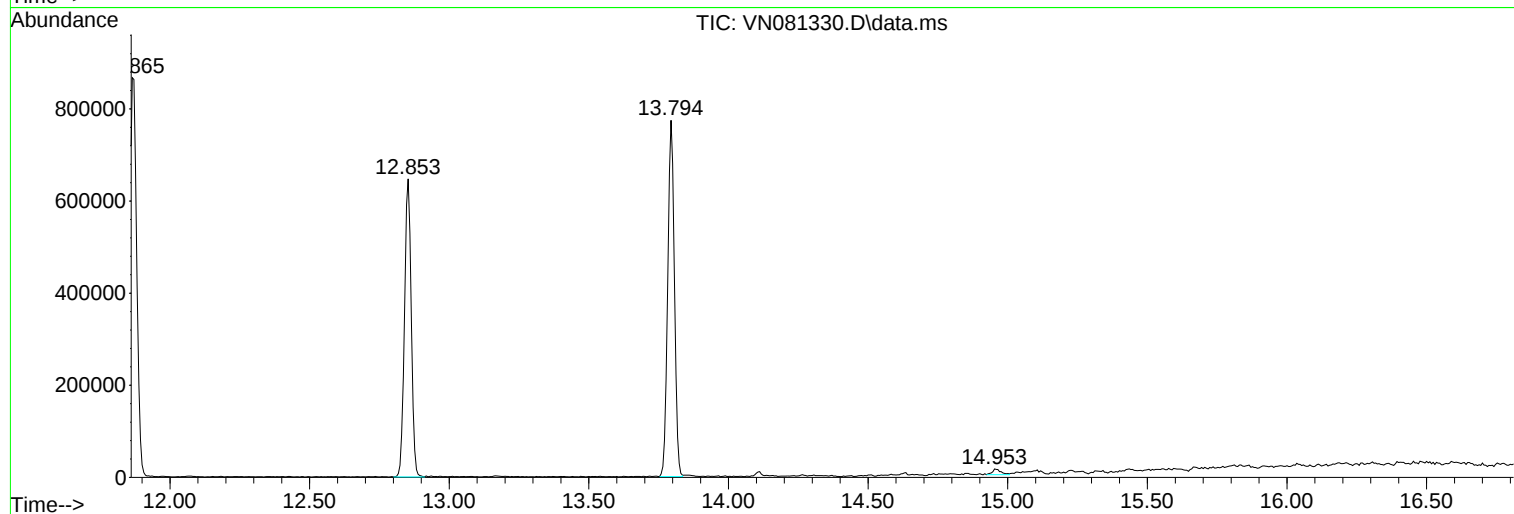
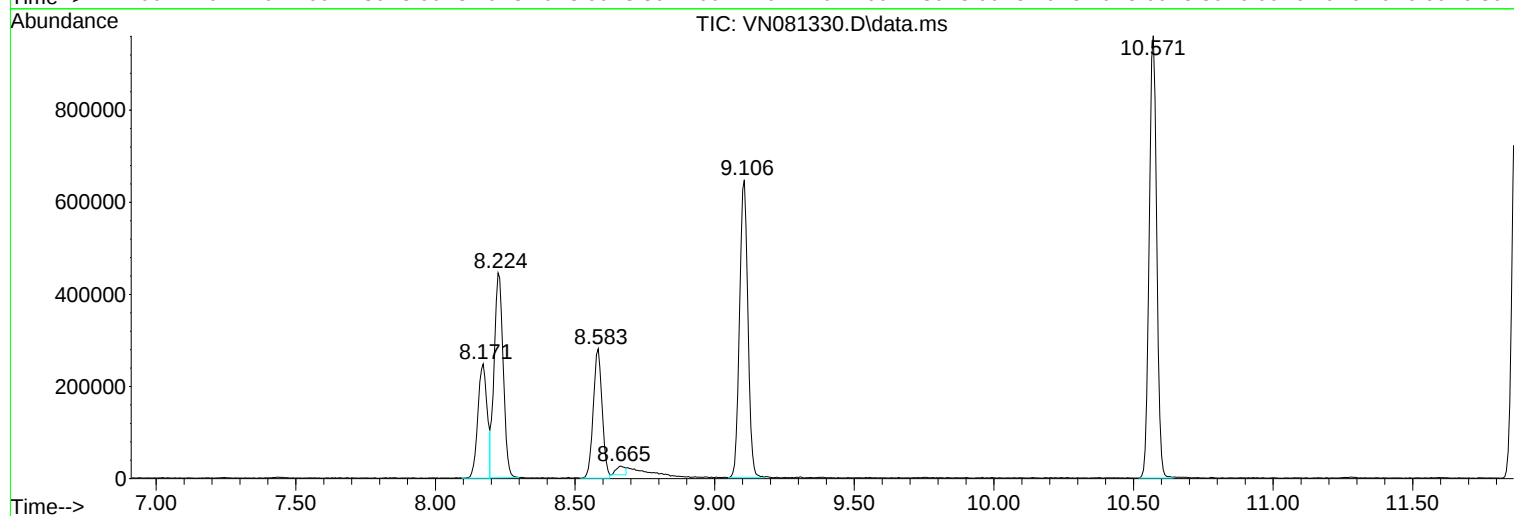
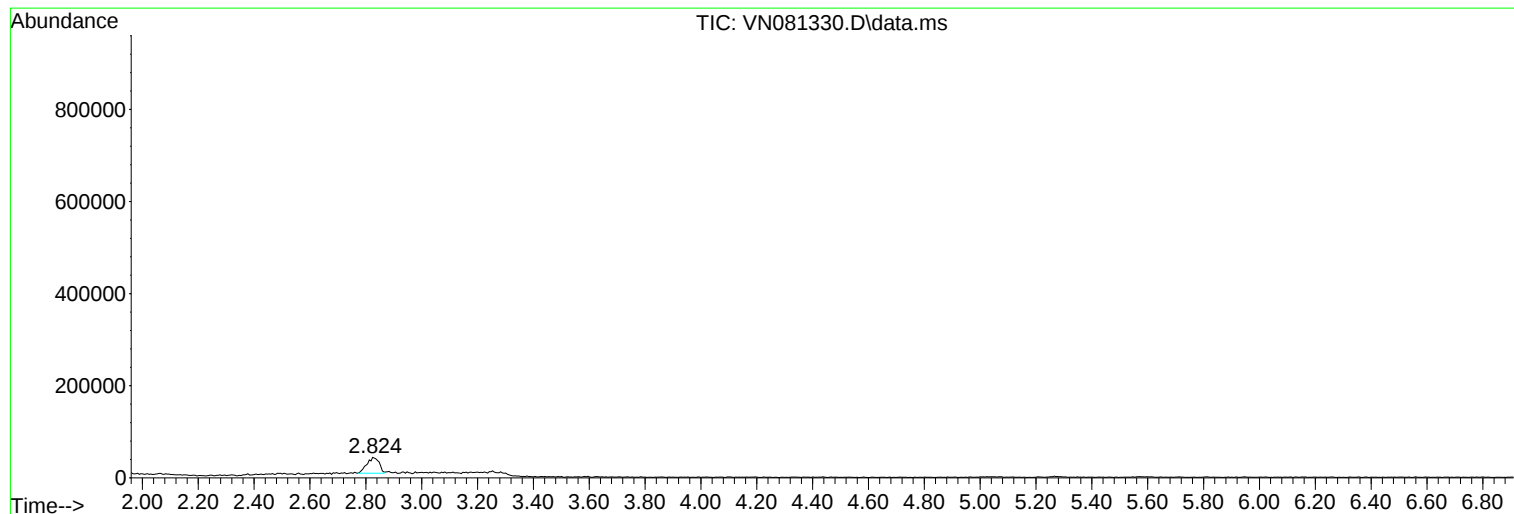
Sum of corrected areas: 9423755

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ClientSampleId :
1132

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030524W.M
Quant Title : SW846 8260

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



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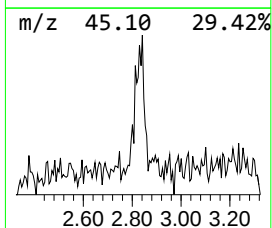
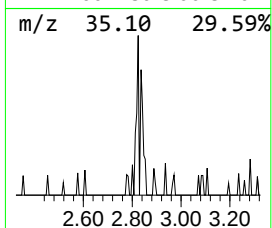
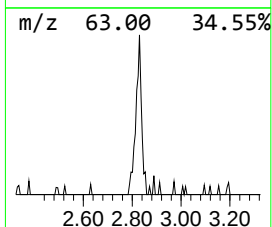
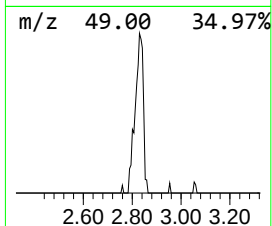
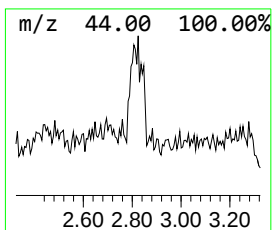
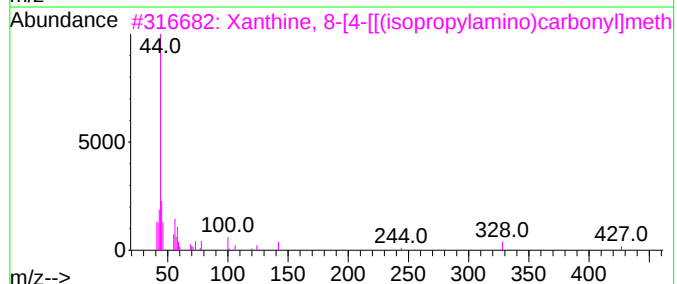
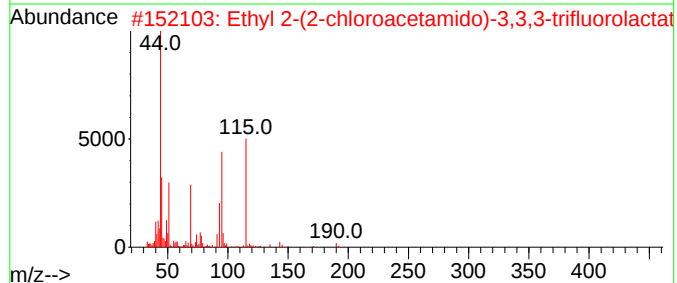
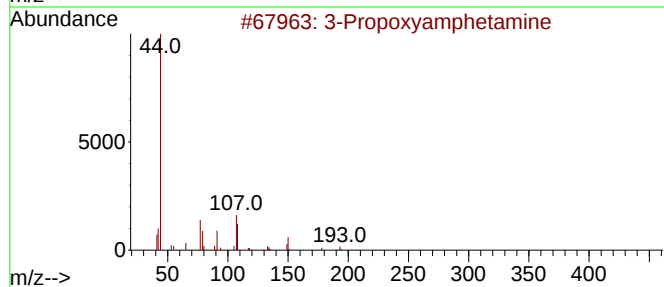
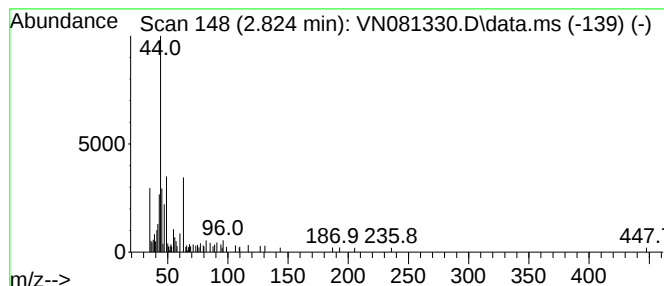
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030524W.M
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown2.824 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.824	5.29 ug/l	105009	Pentafluorobenzene	8.224

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Propoxyamphetamine		193	C12H19NO	1000124-04-4	12
2	Ethyl 2-(2-chloroacetamido)-3,3,...		263	C7H9ClF3NO4	1000224-74-2	12
3	Xanthine, 8-[4-[[(isopropylamino...		427	C22H29N5O4	1000116-67-8	12
4	Phenethylamine, p,.alpha.-dimethyl-		149	C10H15N	000064-11-9	10
5	Amidephrine		244	C10H16N2O3S	003354-67-4	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown2.824	2.824	5.3	ug/l	105009	1	8.224	991596	50.0