

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN032425\
 Data File : VN086074.D
 Acq On : 24 Mar 2025 16:00
 Operator : JC\MD
 Sample : Q1606-01
 Misc : 5.08g/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CHRT24743

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\82N031825W.M

Title : SW846 8260

Signal : TIC: VN086074.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.295	52	58	63	rVB	7666	12266	1.24%	0.228%
2	2.865	139	155	156	rBV8	23762	81454	8.22%	1.513%
3	8.159	1046	1055	1060	rBV	149028	360838	36.42%	6.702%
4	8.218	1060	1065	1075	rVB	251644	536551	54.15%	9.965%
5	8.577	1114	1126	1135	rBV2	147597	345188	34.84%	6.411%
6	8.783	1157	1161	1174	rVB4	6051	19433	1.96%	0.361%
7	9.094	1205	1214	1223	rBV	357318	728745	73.55%	13.534%
8	10.565	1454	1464	1471	rBV	549806	990876	100.00%	18.403%
9	10.624	1471	1474	1483	rVB	16409	28749	2.90%	0.534%
10	11.865	1676	1685	1696	rBV	476450	861772	86.97%	16.005%
11	12.847	1845	1852	1862	rBV	365625	620691	62.64%	11.528%
12	13.788	2005	2012	2026	rBV	455162	786521	79.38%	14.607%
13	15.641	2321	2327	2332	rVB3	5511	11350	1.15%	0.211%

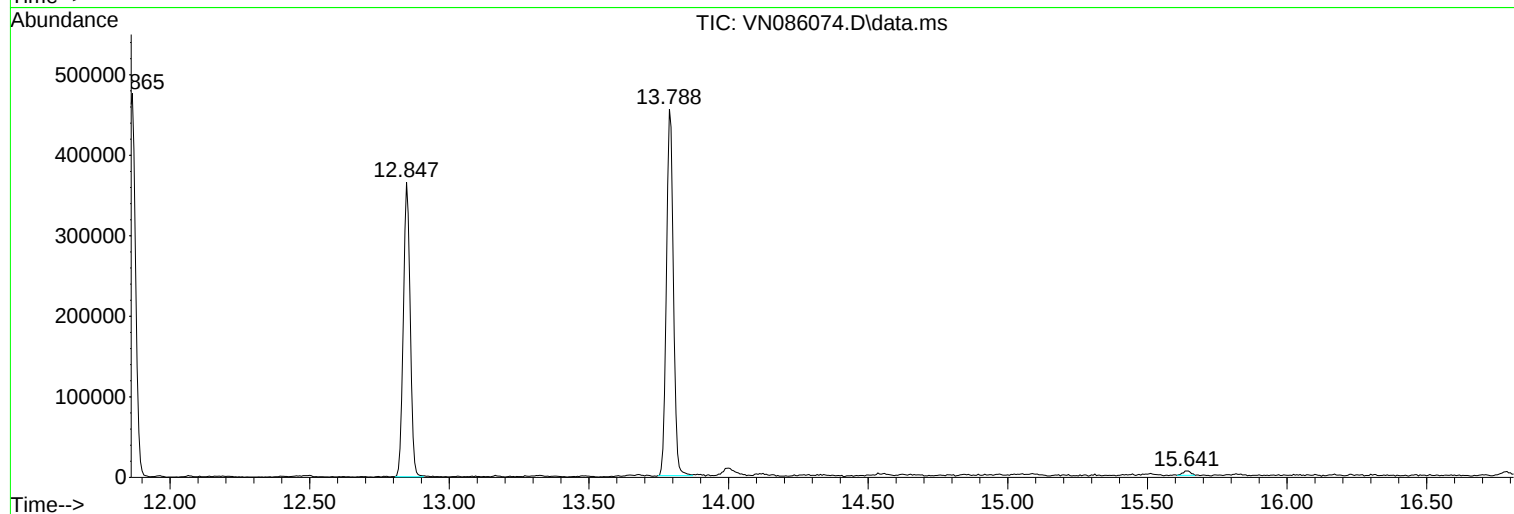
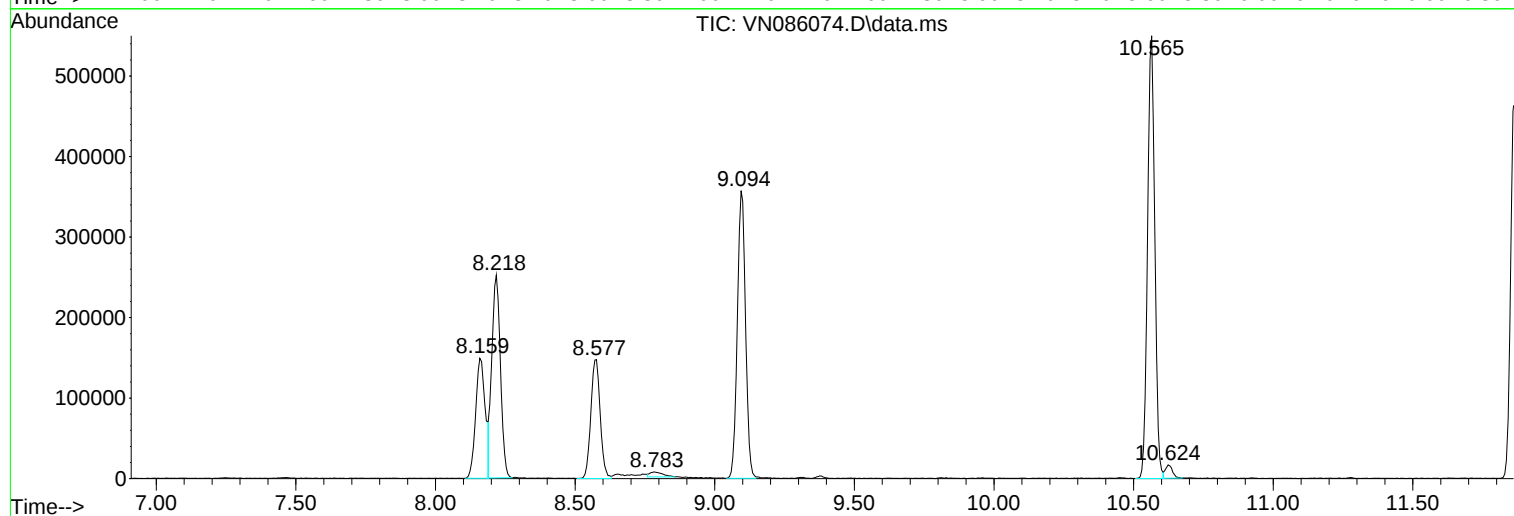
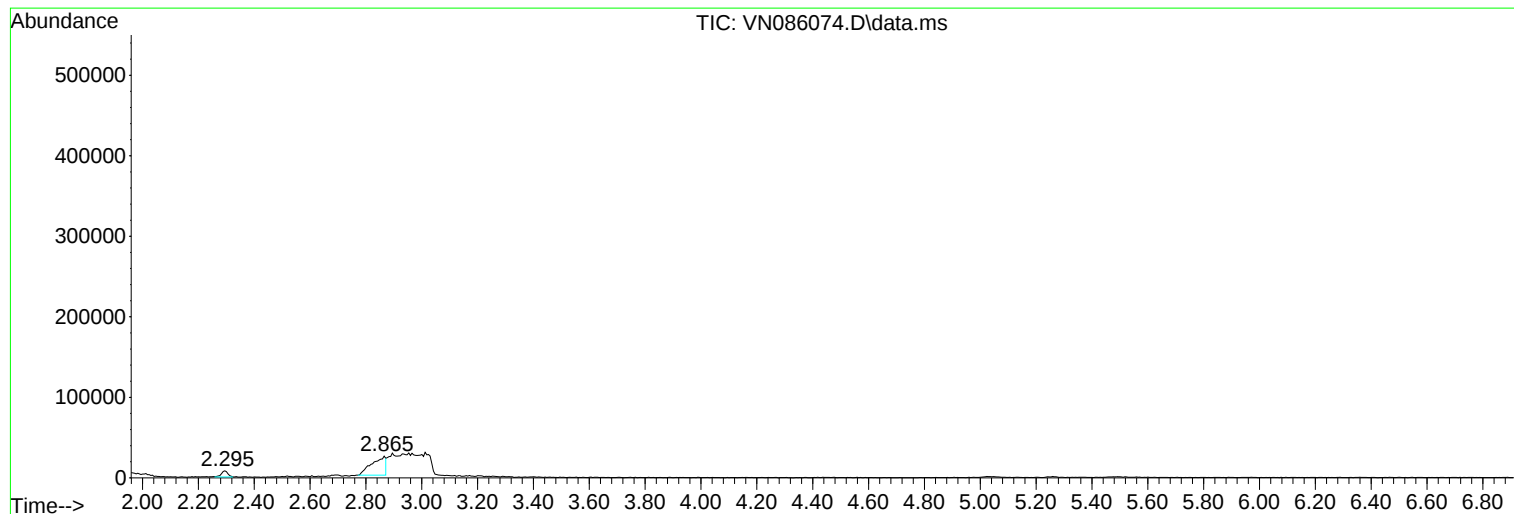
Sum of corrected areas: 5384434

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N031825W.M
Quant Title : SW846 8260

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



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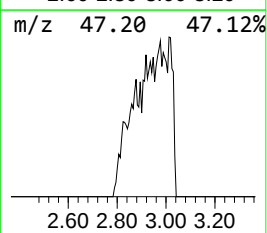
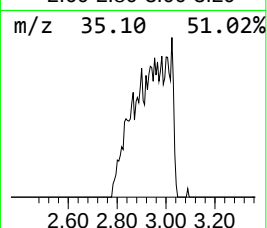
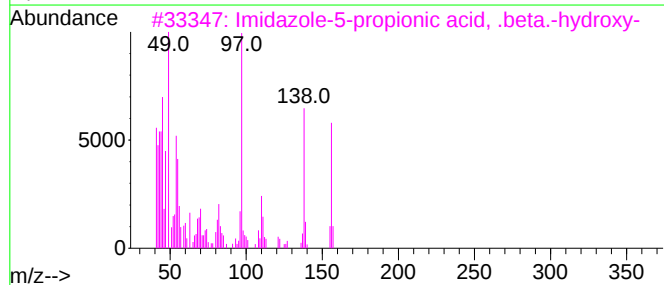
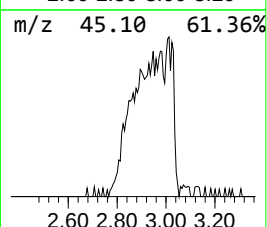
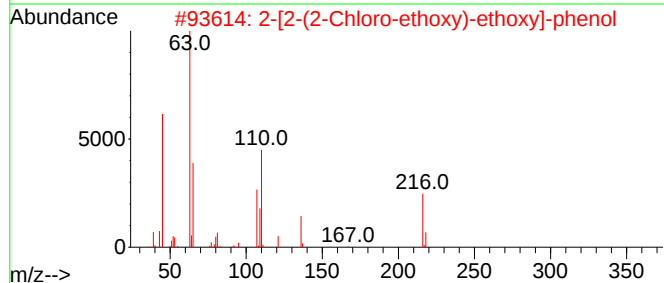
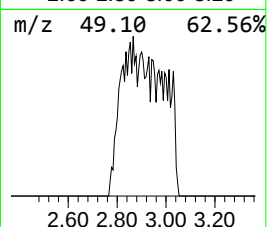
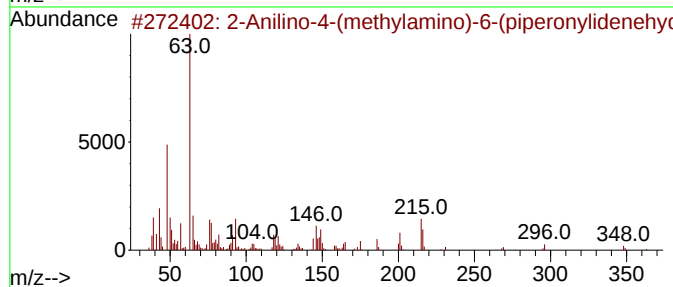
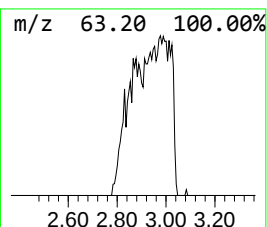
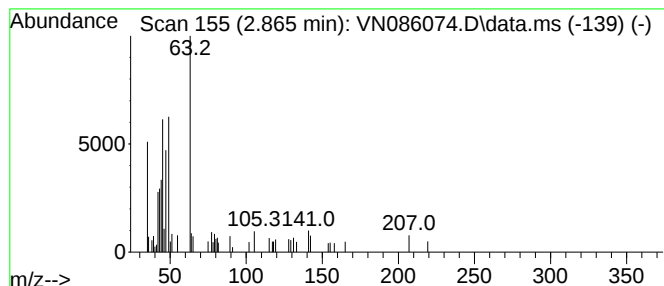
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Peak Number 1 unknown2.865 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.865	7.59 ug/l	81454	Pentafluorobenzene	8.218

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Anilino-4-(methylamino)-6-(pip...	363	C18H17N7O2	1000225-97-3	32		
2	2-[2-(2-Chloro-ethoxy)-ethoxy]-p...	216	C10H13ClO3	1000317-46-1	25		
3	Imidazole-5-propionic acid, .bet...	156	C6H8N2O3	1000129-47-8	16		
4	Methane, chloromethoxy-	80	C2H5ClO	000107-30-2	9		
5	Bromochloroacetaldehyde	156	C2H2BrClO	098136-99-3	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown2.865	2.865	7.6	ug/l	81454	1	8.218	536551	50.0