

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021424\
Data File : VN081026.D
Acq On : 14 Feb 2024 15:34
Operator : JC\MD
Sample : VN0214MBL01
Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VN0214MBL01

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\82N021424W.M
Title : SW846 8260

Signal : TIC: VN081026.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.665	116	121	131	rVB	92645	177086	5.16%	1.070%
2	2.806	141	145	153	rVB	25855	56376	1.64%	0.341%
3	8.165	1046	1056	1061	rBV	340284	828876	24.15%	5.009%
4	8.224	1061	1066	1076	rVB	689119	1523216	44.38%	9.205%
5	8.582	1117	1127	1139	rBV	327524	748232	21.80%	4.522%
6	9.100	1206	1215	1227	rBV	1167544	2410257	70.22%	14.565%
7	10.571	1453	1465	1473	rBV	1350608	2504198	72.96%	15.133%
8	11.865	1677	1685	1696	rVB	1877330	3432469	100.00%	20.742%
9	12.853	1845	1853	1862	rBV	1109378	1903619	55.46%	11.504%
10	13.794	2005	2013	2028	rVB	1701649	2880803	83.93%	17.409%
11	14.041	2050	2055	2064	rBV8	14147	35266	1.03%	0.213%
12	15.853	2357	2363	2371	rVB7	22874	47649	1.39%	0.288%

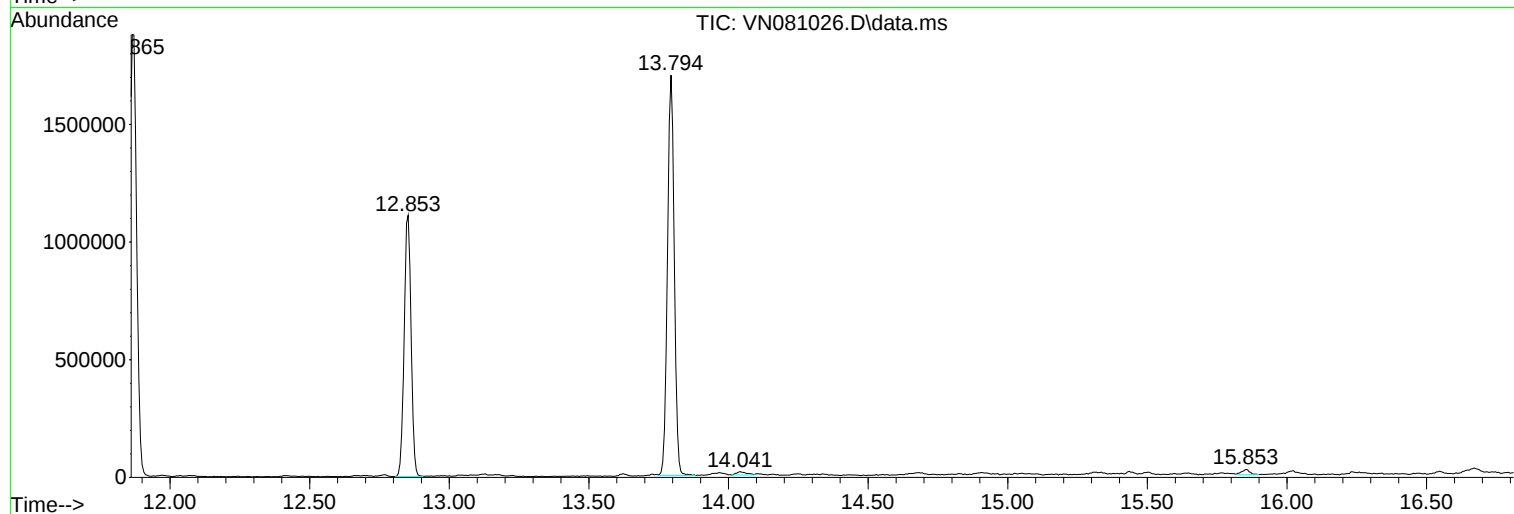
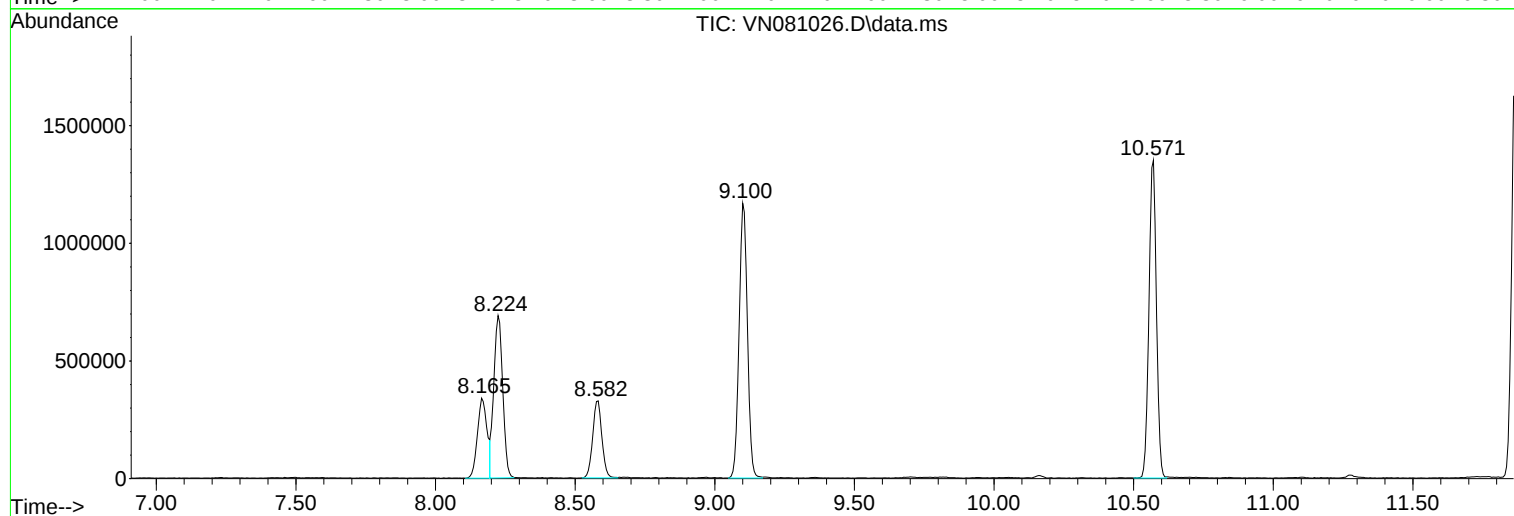
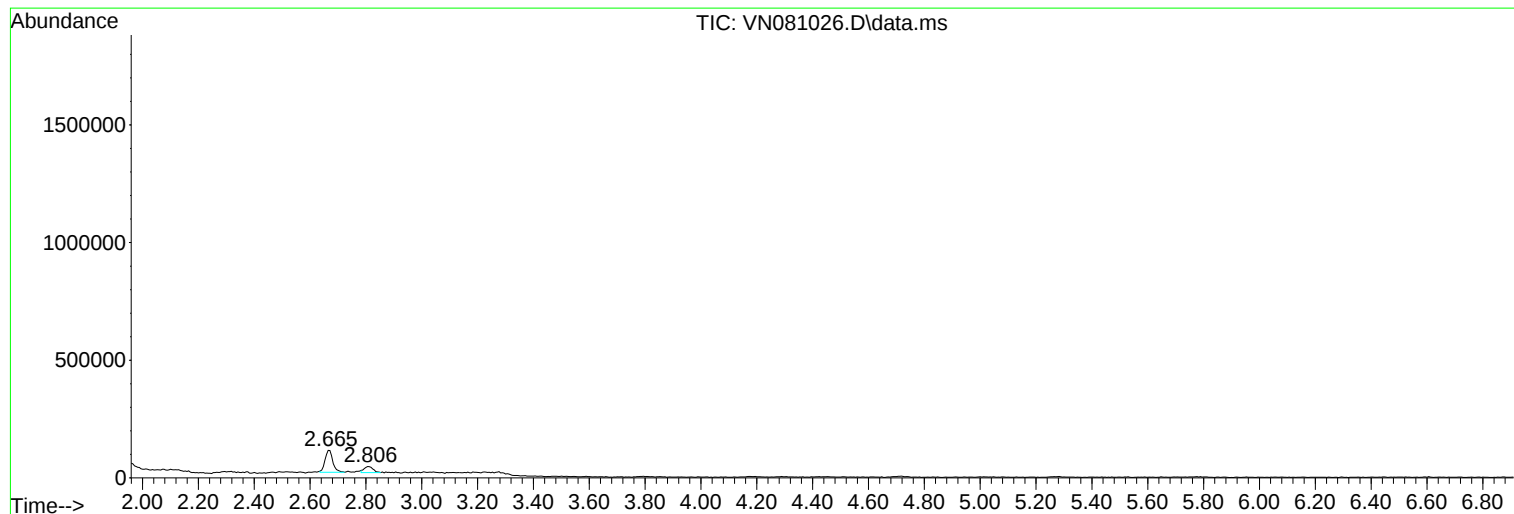
Sum of corrected areas: 16548047

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021424W.M
Quant Title : SW846 8260

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



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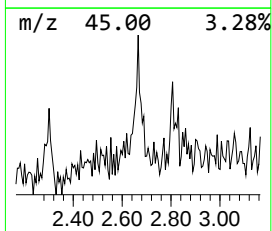
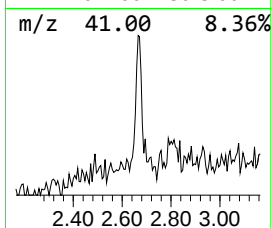
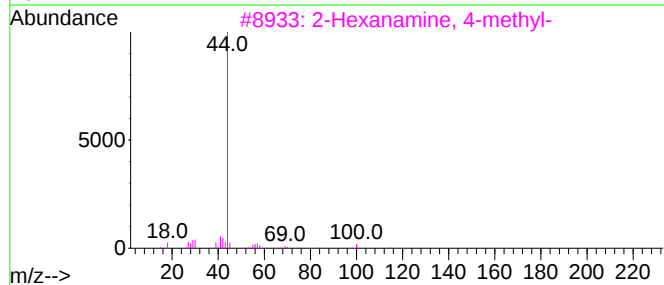
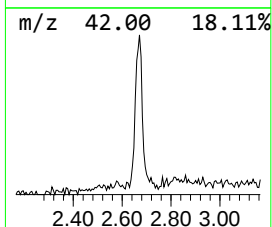
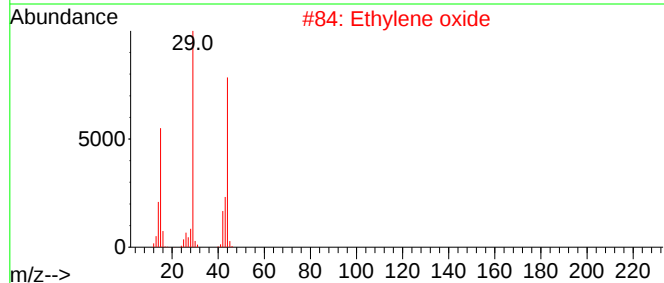
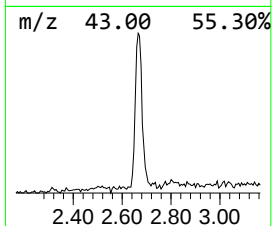
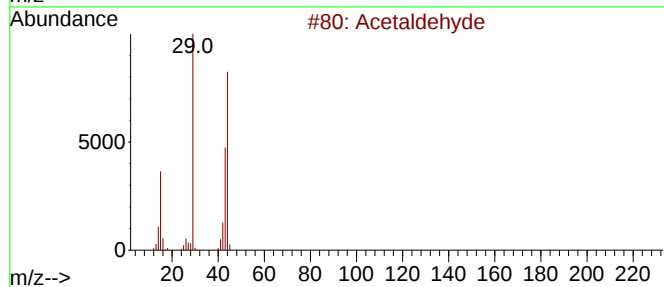
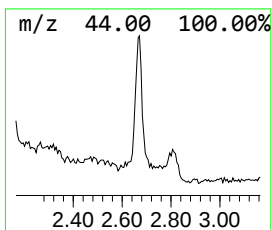
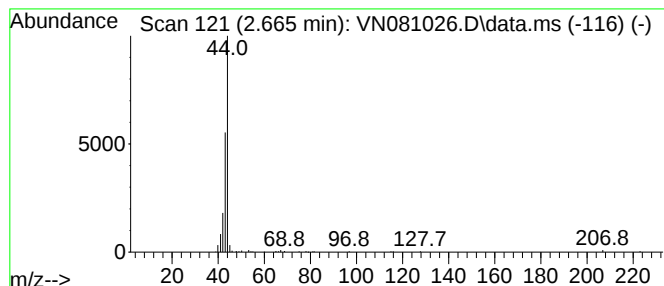
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021424W.M
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Peak Number 1 unknown2.665 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.665	5.81 ug/l	177086	Pentafluorobenzene	8.224

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetaldehyde	44	C2H4O	000075-07-0	56
2			Ethylene oxide	44	C2H4O	000075-21-8	9
3			2-Hexanamine, 4-methyl-	115	C7H17N	000105-41-9	5
4			Tuaminoheptane	115	C7H17N	000123-82-0	5
5			Propane	44	C3H8	000074-98-6	4



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown2.665	2.665	5.8	ug/l	177086	1	8.224	1523220	50.0