

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091922\
Data File : VX031470.D
Acq On : 19 Sep 2022 19:13
Operator : JC/MD
Sample : N4558-04RE
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BP-VPB-RW5-GW-148-150RE

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_X\Method\82X091922W.M

Title : SW846 8260

Signal : TIC: VX031470.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.575	73	80	88	rBV3	6510	16269	2.47%	0.605%
2	2.081	156	163	170	rBV	11372	24367	3.70%	0.907%
3	5.391	694	706	720	rBV	79320	228807	34.70%	8.513%
4	5.556	723	733	745	rVB	74484	203972	30.94%	7.589%
5	5.958	788	799	814	rBV2	87270	228119	34.60%	8.487%
6	6.763	922	931	942	rBV	113170	269668	40.90%	10.033%
7	7.123	983	990	1000	rVB3	5227	12354	1.87%	0.460%
8	8.653	1233	1241	1251	rBV	432384	659341	100.00%	24.532%
9	10.055	1465	1471	1483	rBV	241941	325055	49.30%	12.094%
10	11.085	1634	1640	1651	rBV	325632	423365	64.21%	15.752%
11	12.024	1789	1794	1803	rBV	248394	296409	44.96%	11.028%

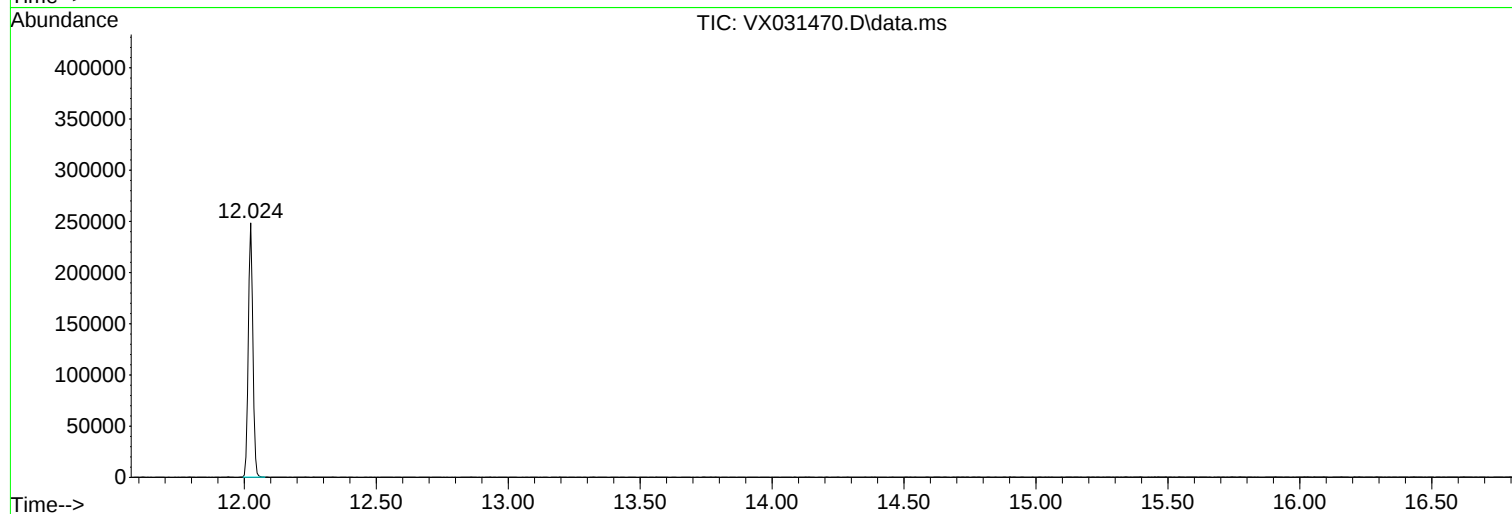
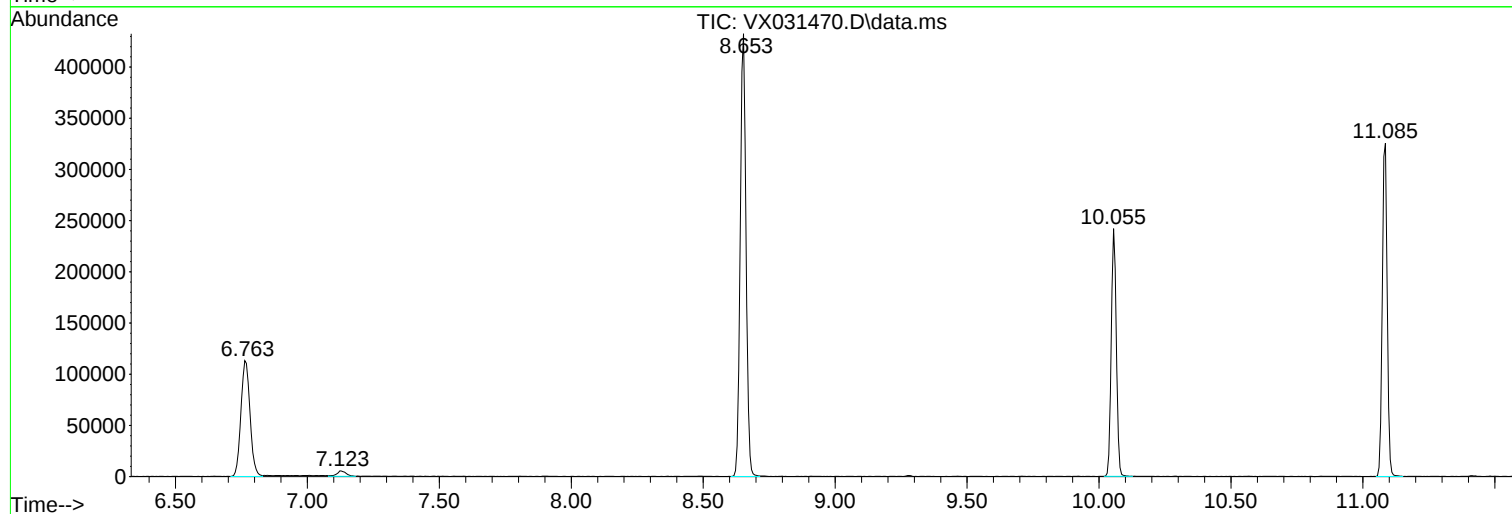
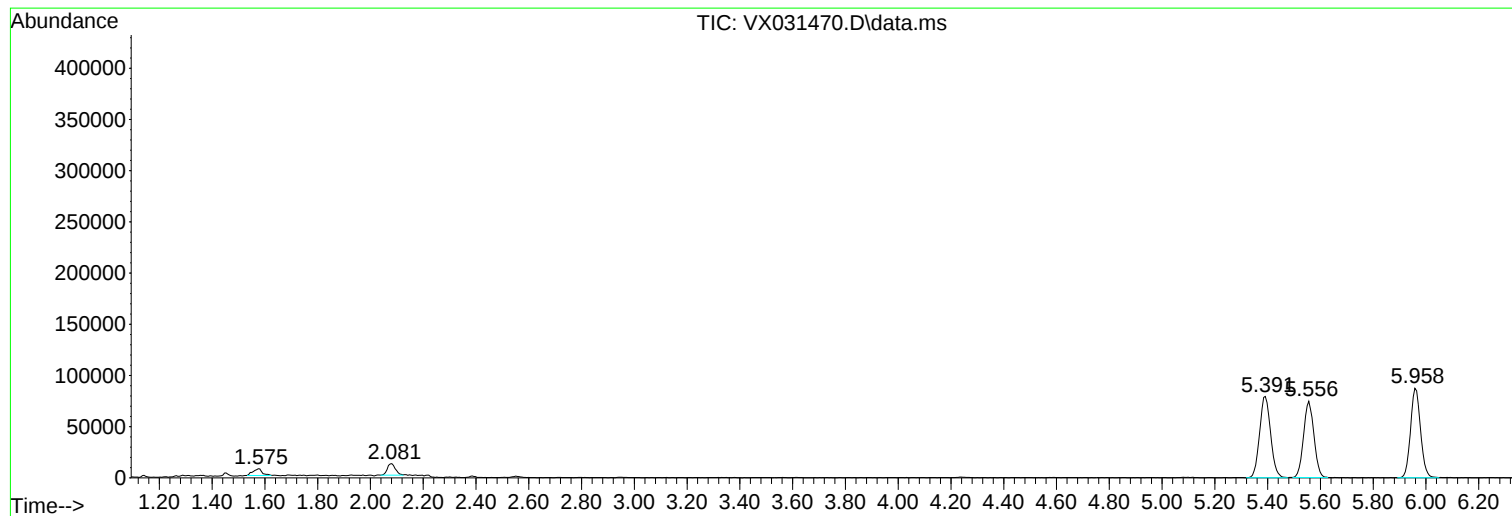
Sum of corrected areas: 2687726

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

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



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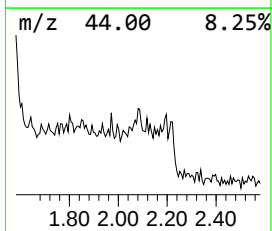
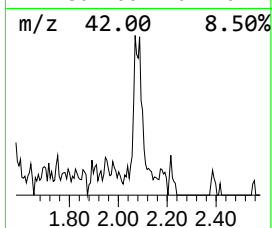
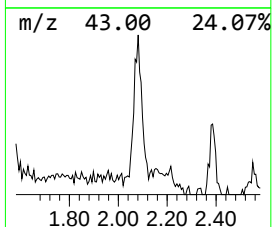
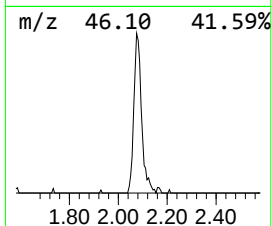
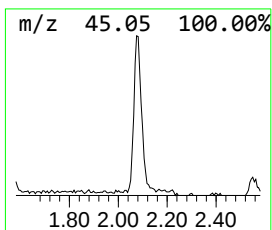
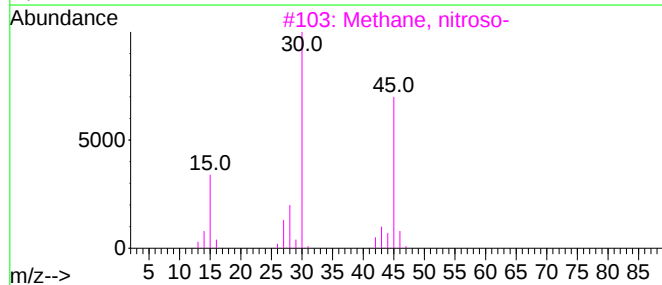
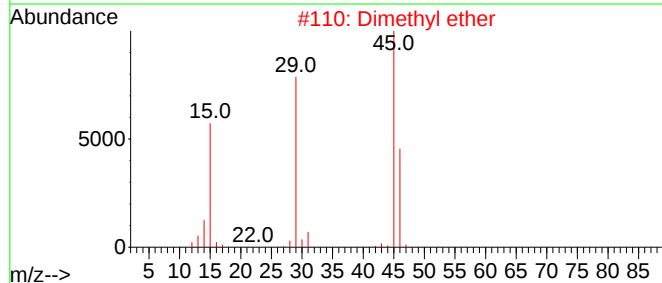
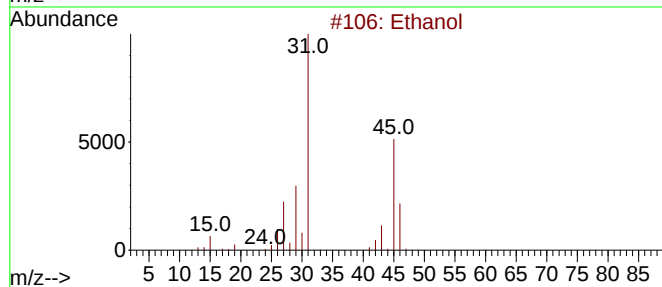
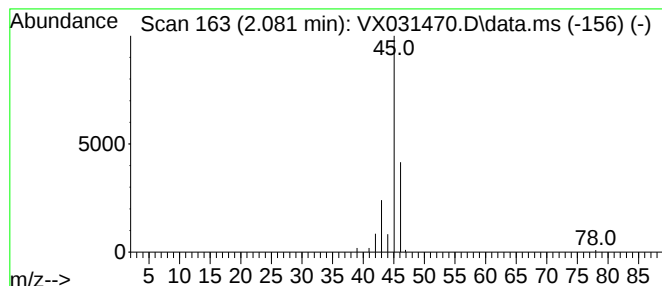
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

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

Peak Number 1 Ethanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.081	5.97 ug/l	24367	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol	46	C2H6O	000064-17-5	74
2			Dimethyl ether	46	C2H6O	000115-10-6	9
3			Methane, nitroso-	45	CH3NO	000865-40-7	7
4			Formic acid	46	CH2O2	000064-18-6	5
5			2-Hexanol, 3-methyl-	116	C7H16O	002313-65-7	4



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol	2.081	6.0	ug/l	24367	1	5.556	203972	50.0