

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010924\
Data File : VX039676.D
Acq On : 09 Jan 2024 13:38
Operator : JC/MD
Sample : VX0109WBL01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VX0109WBL01

Integration Parameters: RTEINT3.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\624X010824W.M

Title : METHOD 624 VOLATILE ORGANIC ANALYSIS

Signal : TIC: VX039676.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	70	80	81	rBV4	5490	12917	5.07%	1.218%
2	1.618	81	87	88	rVV6	3218	6234	2.45%	0.588%
3	2.941	298	304	316	rVB	4343	9806	3.85%	0.924%
4	4.897	616	625	640	rBV2	20035	61619	24.18%	5.809%
5	5.952	789	798	815	rBV2	39009	104936	41.17%	9.892%
6	6.757	922	930	952	rBV	64556	160219	62.86%	15.104%
7	8.647	1233	1240	1253	rBV	160065	254865	100.00%	24.026%
8	10.055	1465	1471	1483	rBV	165047	240613	94.41%	22.682%
9	11.079	1634	1639	1655	rVB	153119	206329	80.96%	19.450%
10	13.780	2078	2082	2086	rBV2	2803	3265	1.28%	0.308%

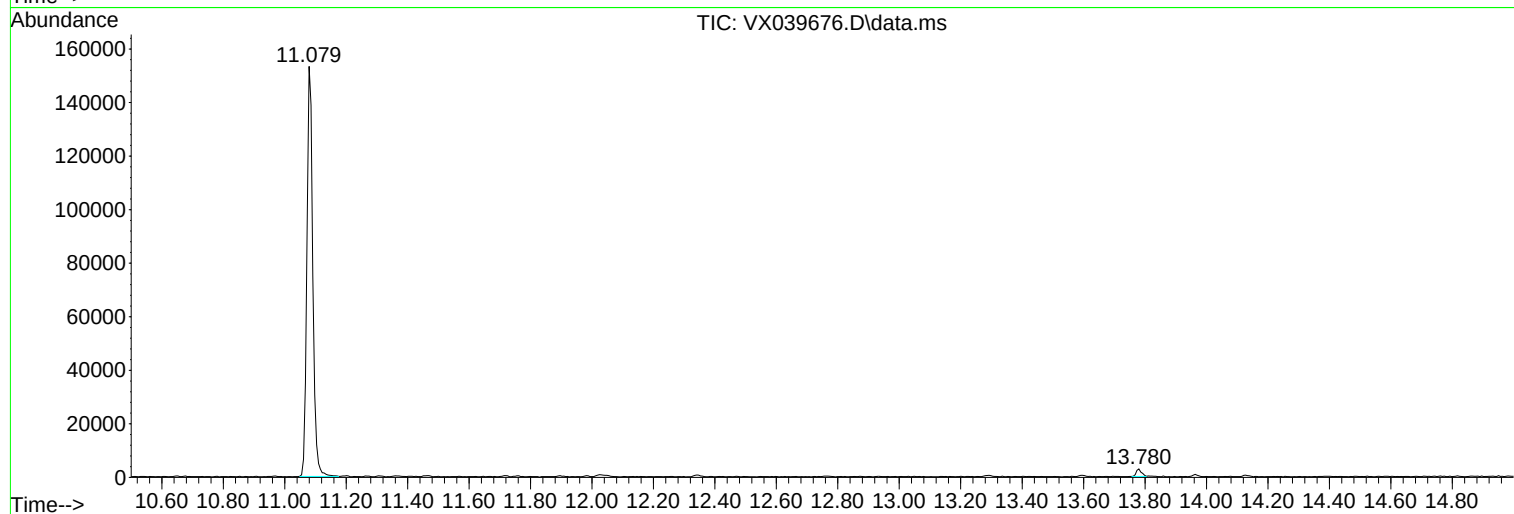
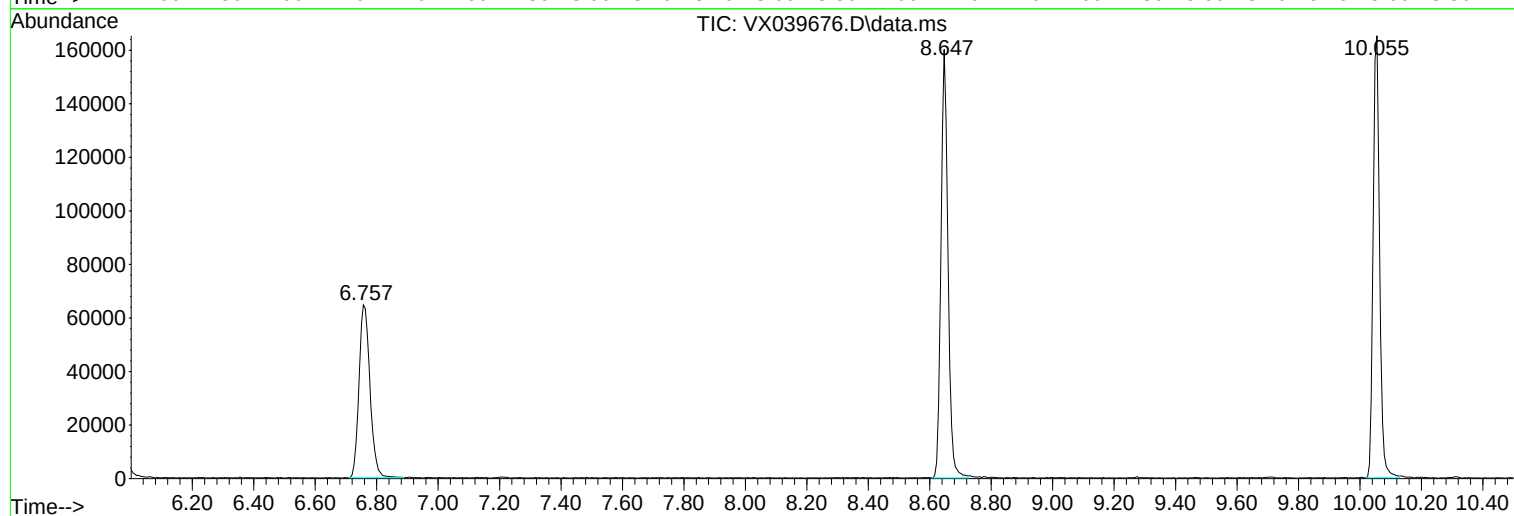
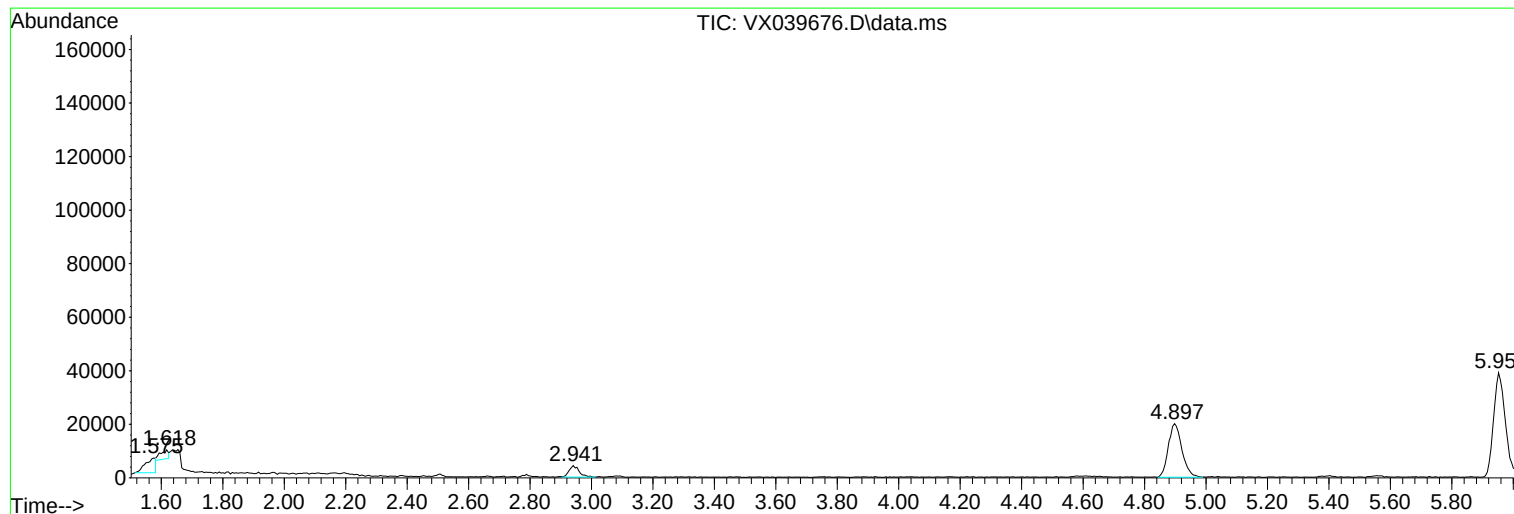
Sum of corrected areas: 1060803

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\624X010824W.M
Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS

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



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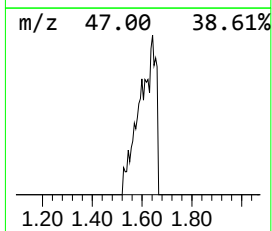
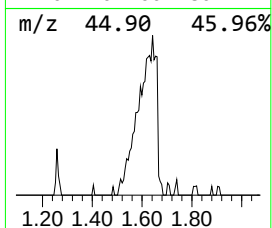
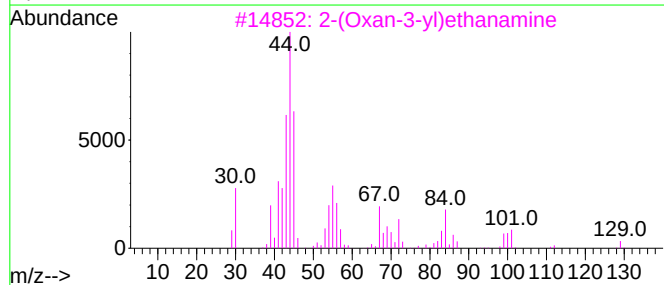
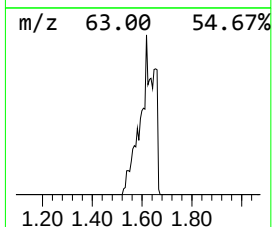
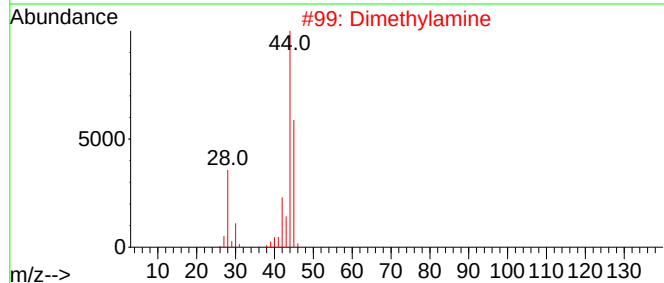
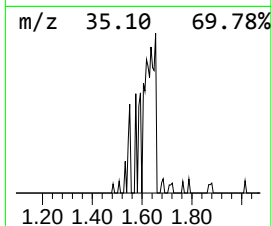
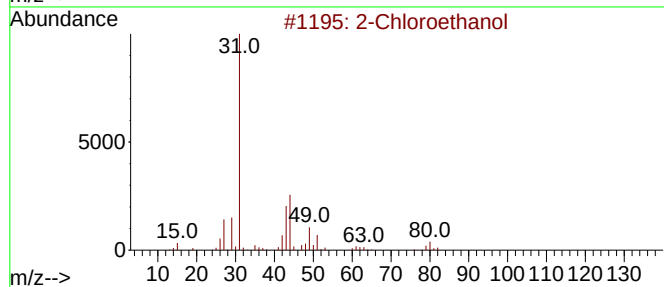
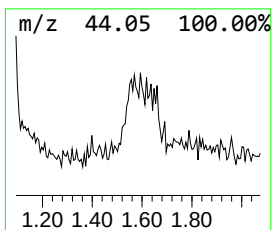
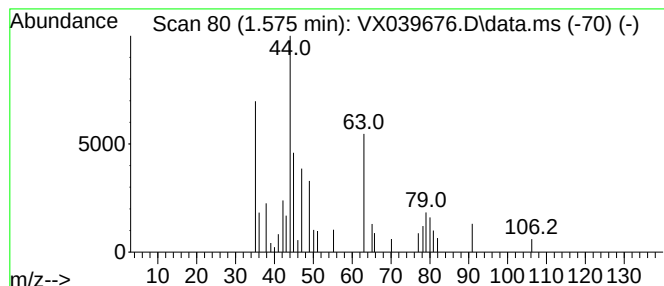
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\624X010824W.M
Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS

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

Peak Number 1 unknown1.575 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.575	6.29 ug/l	12917	Bromochloromethane	4.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Chloroethanol	80	C2H5ClO	000107-07-3	10
2			Dimethylamine	45	C2H7N	000124-40-3	9
3			2-(Oxan-3-yl)ethanamine	129	C7H15NO	1000411-07-5	9
4			Dimethylamine	45	C2H7N	000124-40-3	7
5			Propane	44	C3H8	000074-98-6	5



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.575	1.575	6.3	ug/l	12917	1	4.904	61619	30.0