

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032423\
Data File : VX034726.D
Acq On : 24 Mar 2023 17:15
Operator : JC/MD
Sample : 02073-04 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1660

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\82X031023W.M

Title : SW846 8260

Signal : TIC: VX034726.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.392	206	214	247	rBV	10144875	22174886	100.00%	60.690%
2	3.946	449	469	471	rBV3	77192	304595	1.37%	0.834%
3	5.385	693	705	717	rBV2	185333	544960	2.46%	1.491%
4	5.550	718	732	754	rVB	342420	1016249	4.58%	2.781%
5	5.836	768	779	790	rBV	96051	287483	1.30%	0.787%
6	5.952	790	798	819	rVB	225671	612795	2.76%	1.677%
7	6.763	920	931	948	rBV	520543	1256253	5.67%	3.438%
8	8.647	1234	1240	1246	rBV	1096967	1686640	7.61%	4.616%
9	8.720	1246	1252	1267	rVB	2991094	4665943	21.04%	12.770%
10	10.055	1465	1471	1485	rBV	1064773	1462519	6.60%	4.003%
11	11.079	1634	1639	1650	rBV	912417	1146357	5.17%	3.137%
12	12.018	1788	1793	1801	rBV	1075529	1379563	6.22%	3.776%

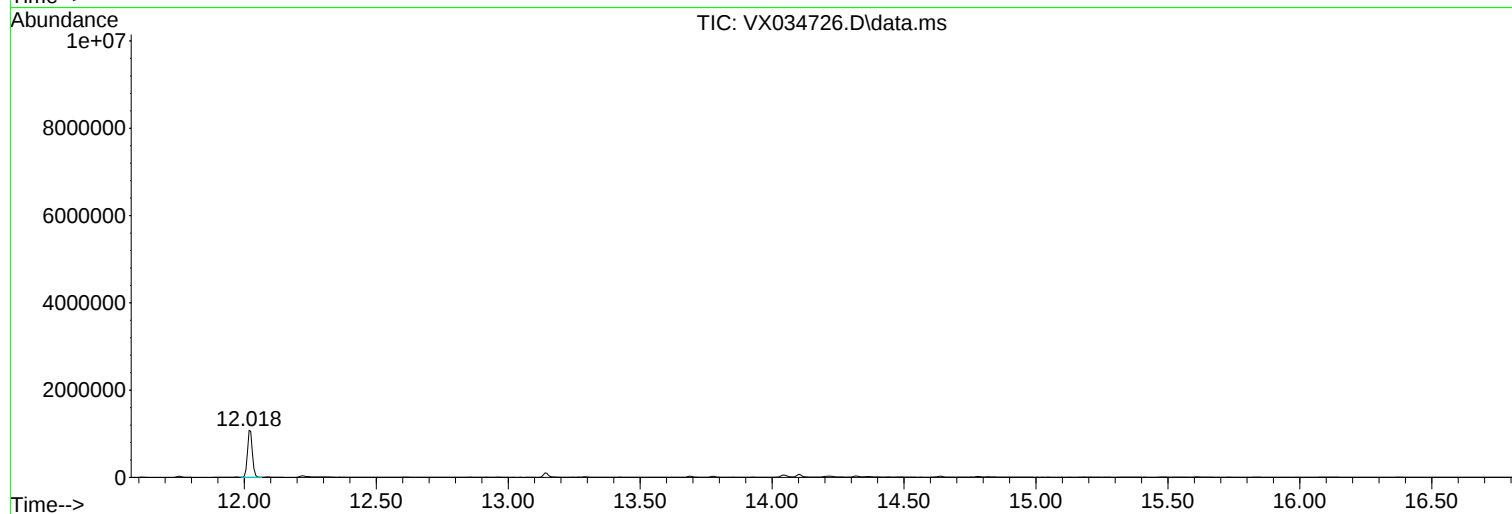
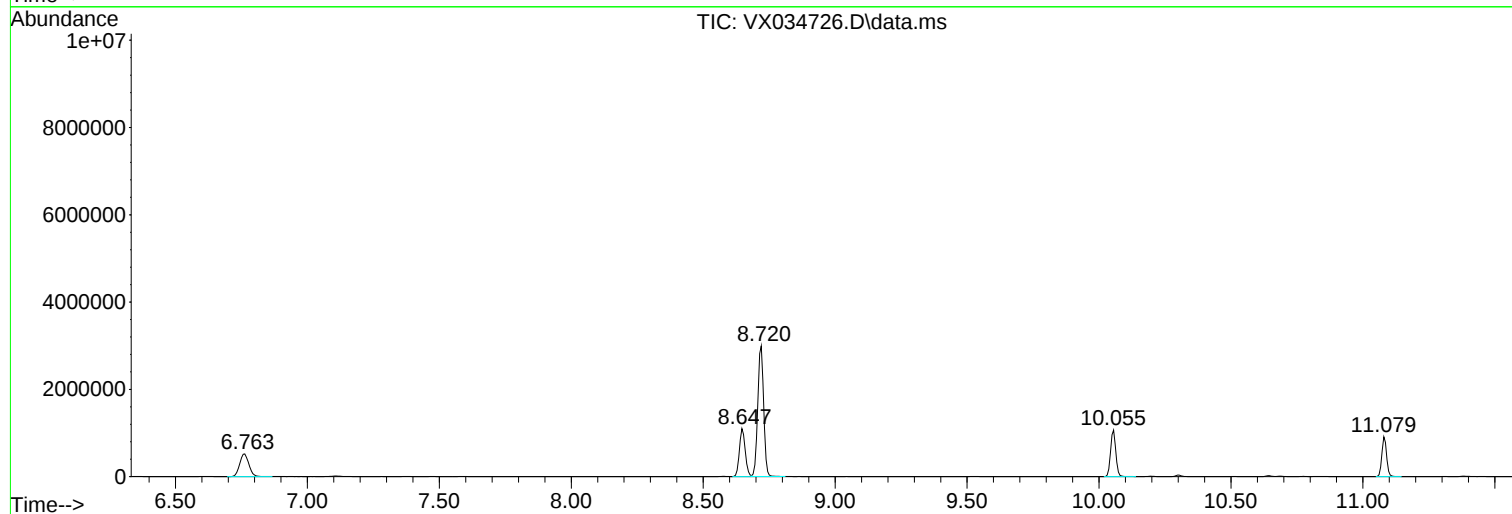
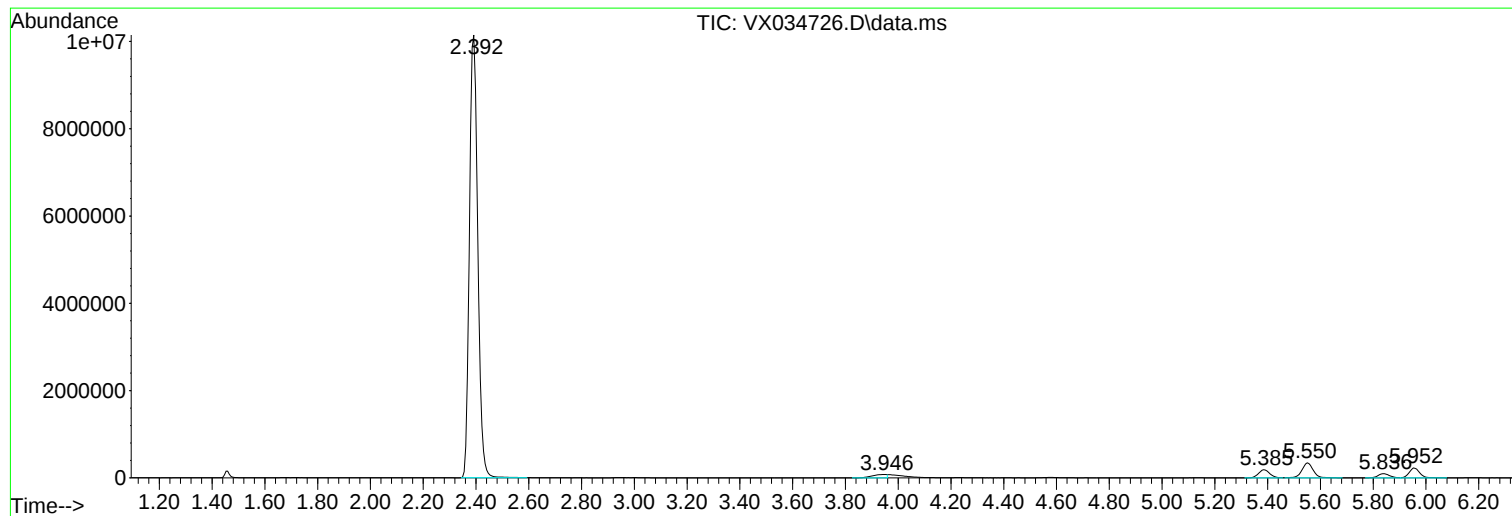
Sum of corrected areas: 36538243

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
Quant Title : SW846 8260

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



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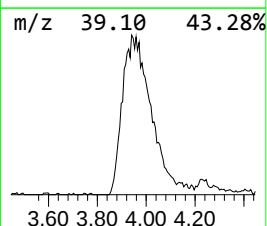
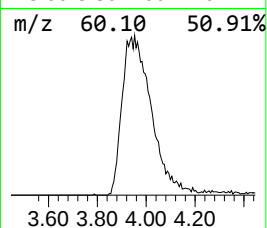
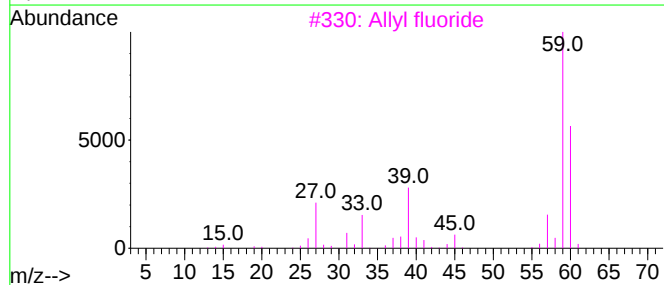
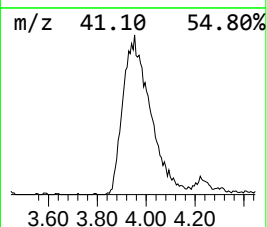
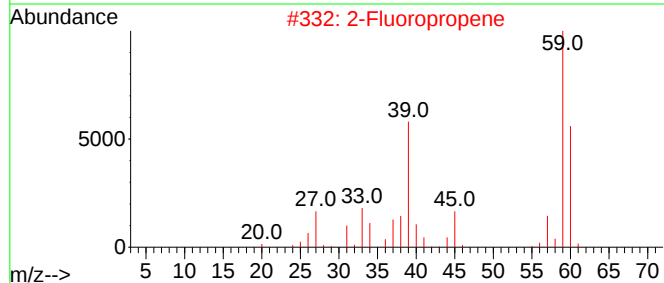
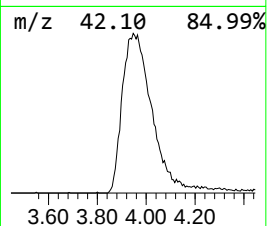
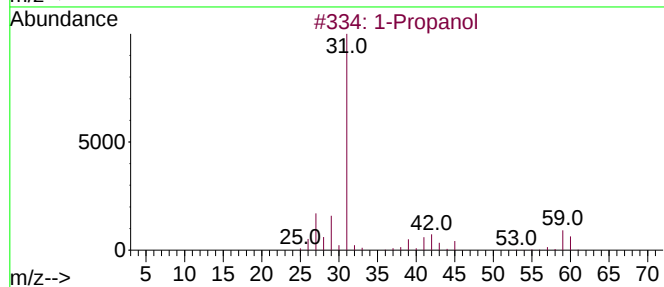
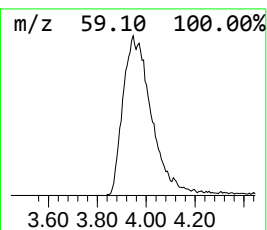
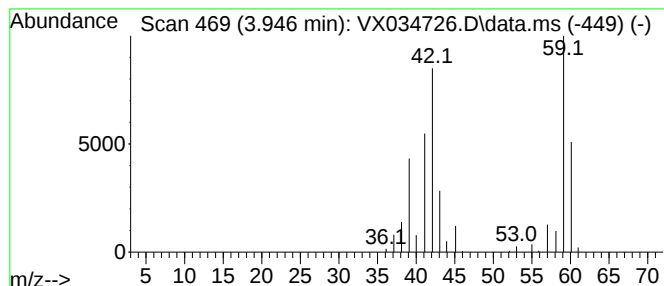
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
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Peak Number 1 1-Propanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.946	14.99 ug/l	304595	Pentafluorobenzene	5.550

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propanol			60	C3H8O	000071-23-8	59
2	2-Fluoropropene			60	C3H5F	001184-60-7	52
3	Allyl fluoride			60	C3H5F	000818-92-8	38
4	Cyclopropane			42	C3H6	000075-19-4	35
5	(3,3-Dimethyloxiranyl)methanol			102	C5H10O2	1010306-71-7	9



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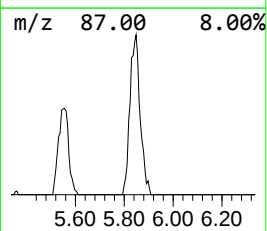
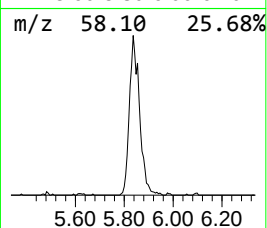
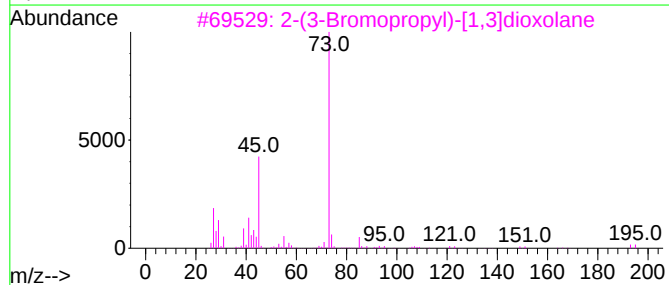
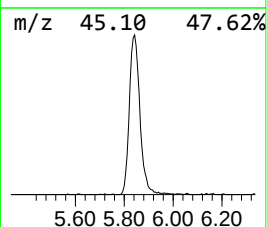
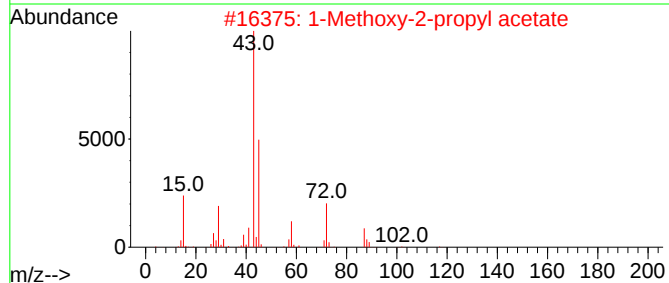
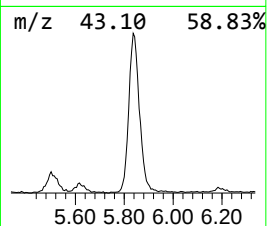
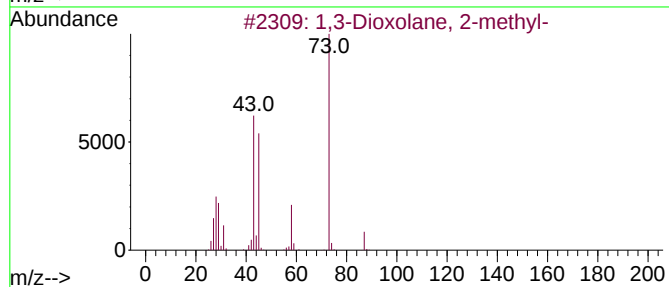
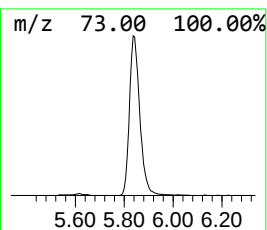
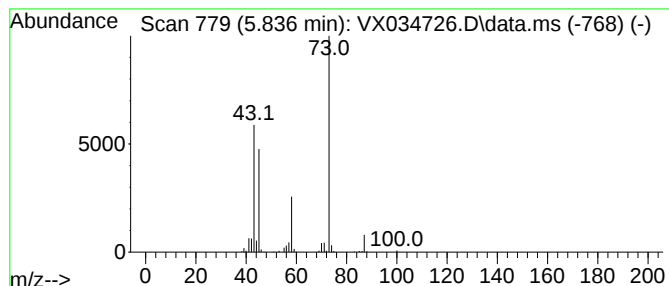
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Peak Number 2 1,3-Dioxolane, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.836	14.14 ug/l	287483	Pentafluorobenzene	5.550

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	64		
2	1-Methoxy-2-propyl acetate	132	C6H12O3	000108-65-6	33		
3	2-(3-Bromopropyl)-[1,3]dioxolane	194	C6H11BrO2	062563-07-9	33		
4	2-Hydroxy-3-hexanone	116	C6H12O2	054073-43-7	28		
5	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	17		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Propanol	3.946	15.0	ug/l	304595	1	5.550	1016250	50.0
1,3-Dioxolane, ...	5.836	14.1	ug/l	287483	1	5.550	1016250	50.0