

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX012721\
Data File : VX020777.D
Acq On : 27 Jan 2021 18:01
Operator : JC/SP
Sample : M1217-18
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
1726-1727

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\82X012521W.M

Title : SW846 8260

Signal : TIC: VX020777.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.611	238	250	355	rBV	8372153	59390349	10.89%	9.244%
2	7.306	1004	1020	1084	rBV	6451346	37561470	6.89%	5.847%
3	11.177	1625	1655	1658	rBV6	80172400	545491955	100.00%	84.909%

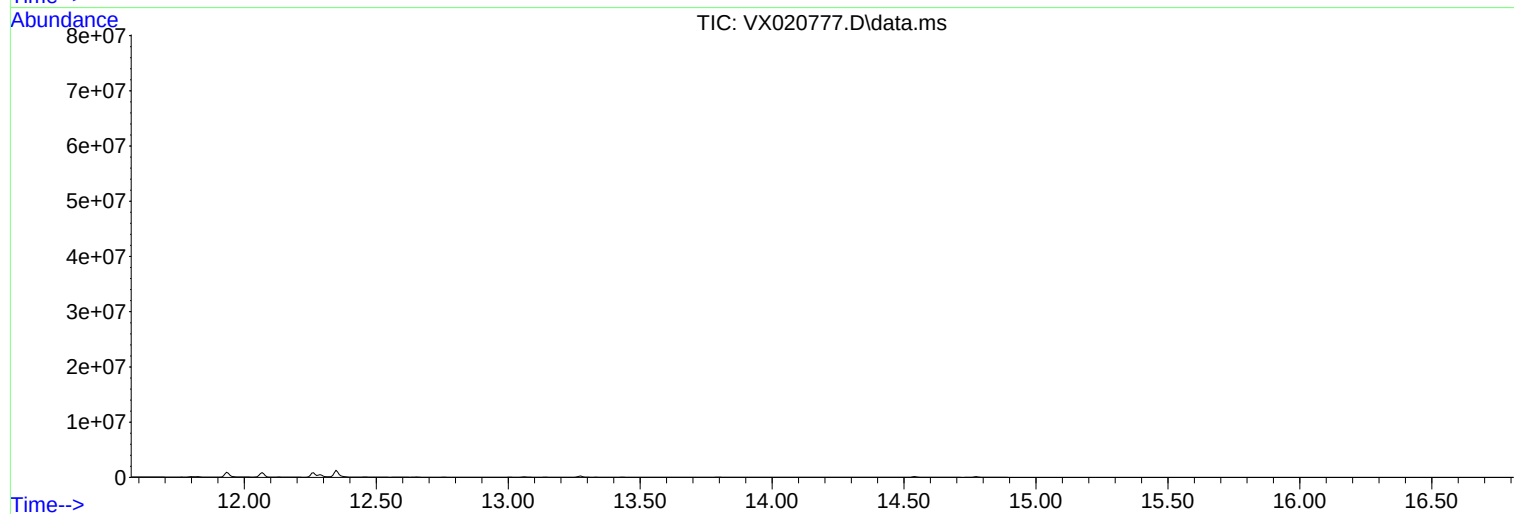
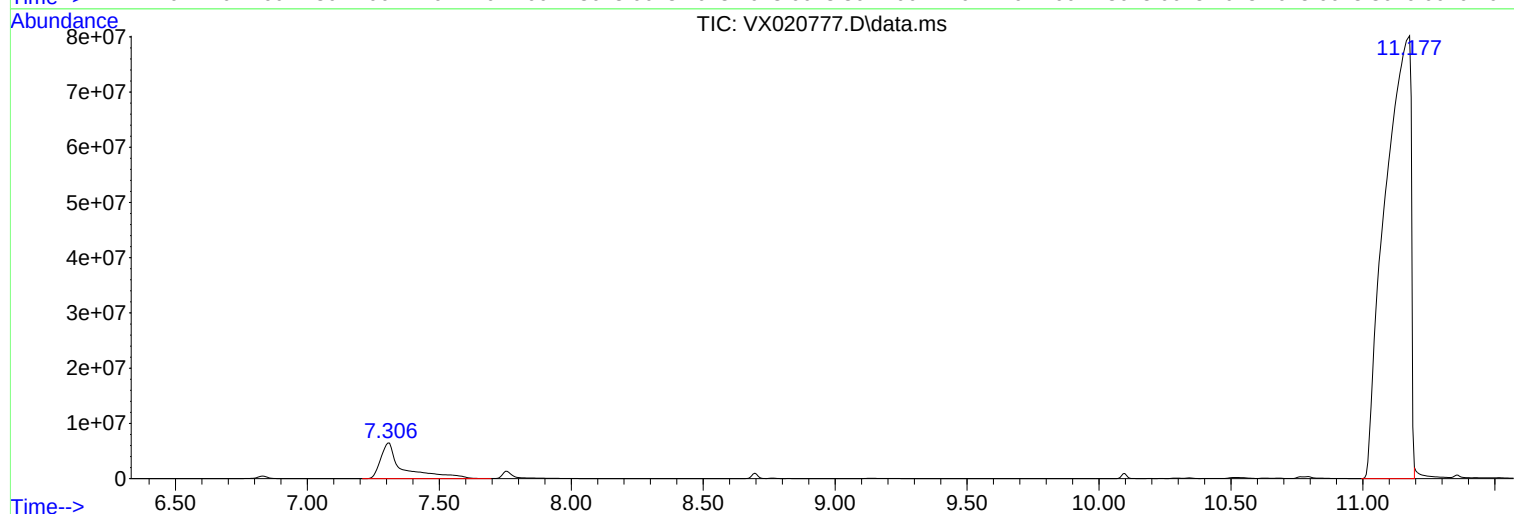
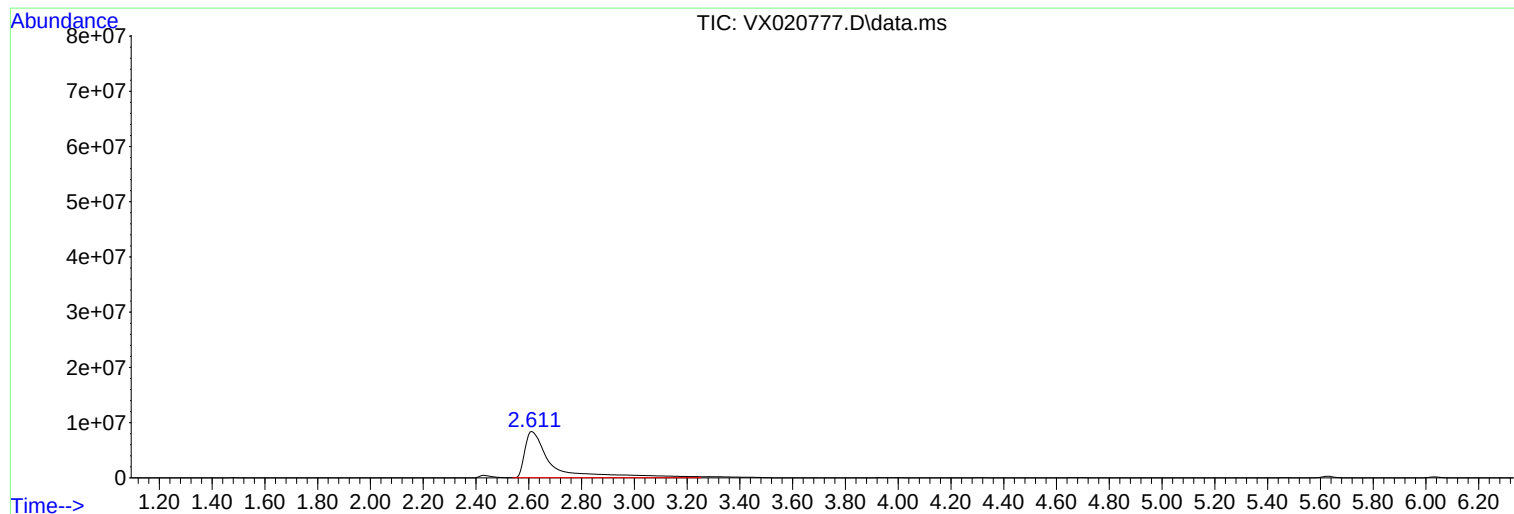
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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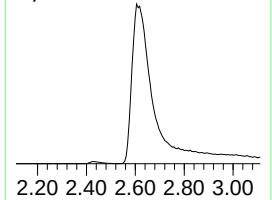
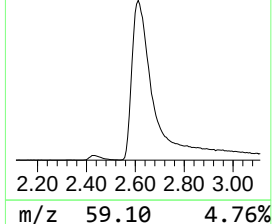
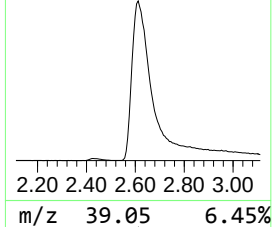
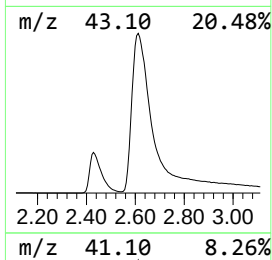
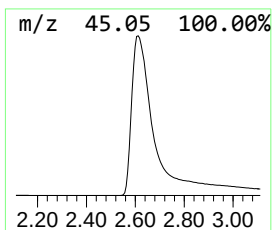
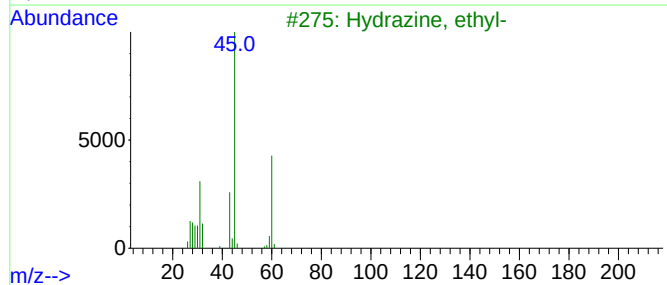
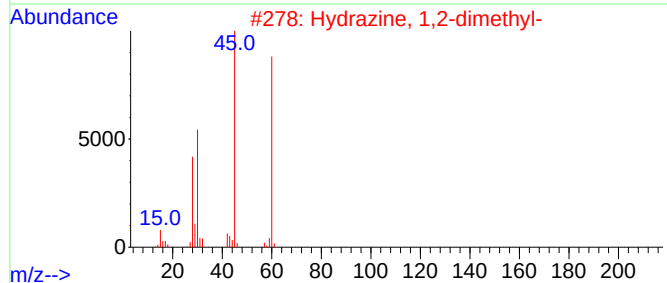
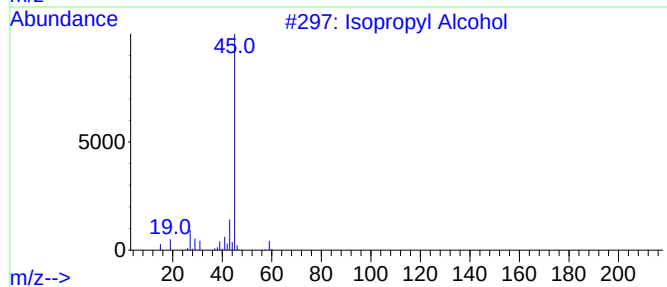
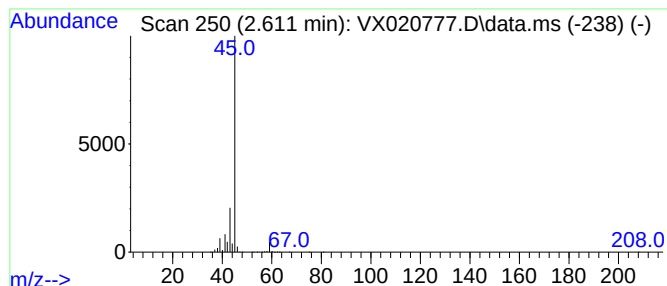
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Isopropyl Alcohol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.611	79.06 ug/l	59390300	Pentafluorobenzene	5.629

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl Alcohol	60	C3H8O	000067-63-0	86
2			Hydrazine, 1,2-dimethyl-	60	C2H8N2	000540-73-8	9
3			Hydrazine, ethyl-	60	C2H8N2	000624-80-6	9
4			Formamide	45	CH3NO	000075-12-7	5
5			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	002155-30-8	4



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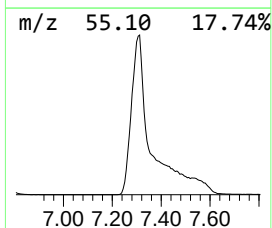
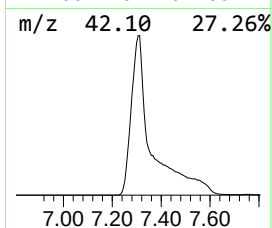
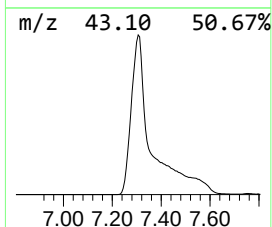
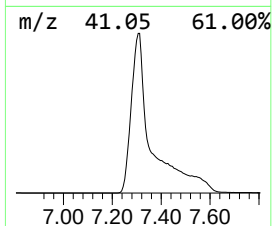
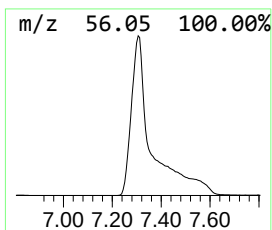
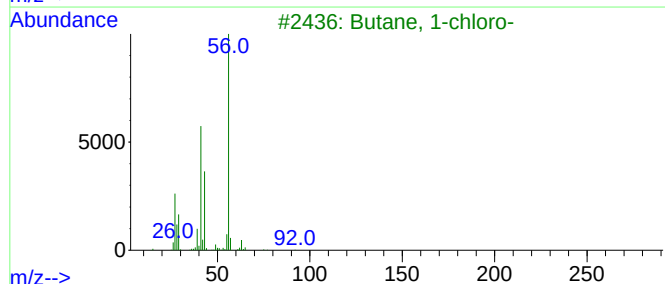
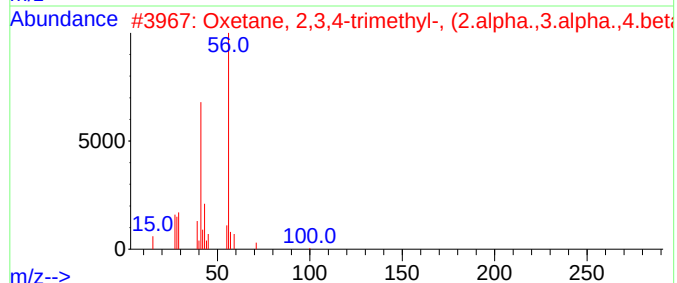
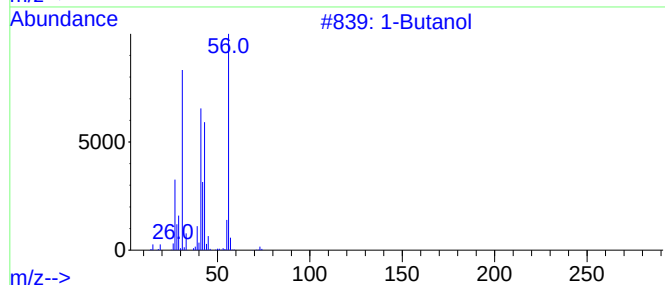
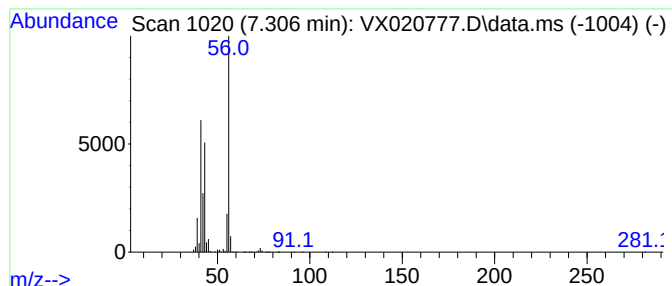
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Peak Number 2 1-Butanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.306	50.00 ug/l	37561500	1,4-Difluorobenzene	6.830

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Butanol	74	C4H10O	000071-36-3	91
2			Oxetane, 2,3,4-trimethyl-, (2.alpha...	100	C6H12O	032347-12-9	38
3			Butane, 1-chloro-	92	C4H9Cl	000109-69-3	38
4			3-Butenoic acid, 2-amino-	101	C4H7NO2	056512-51-7	9
5			Formic acid, butyl ester	102	C5H10O2	000592-84-7	9



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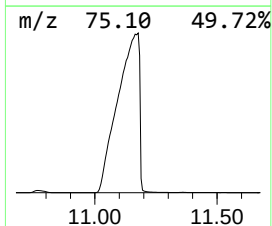
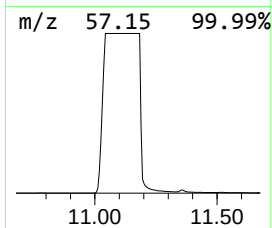
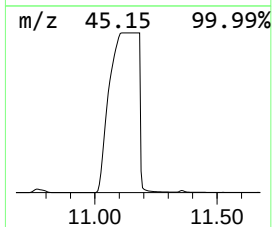
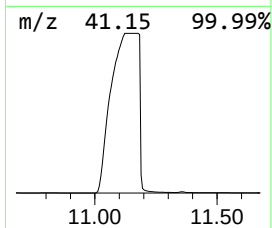
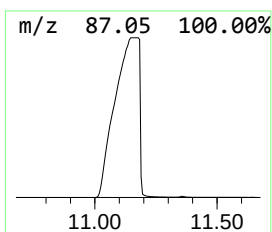
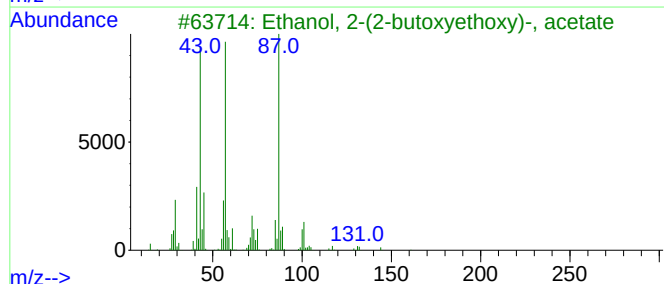
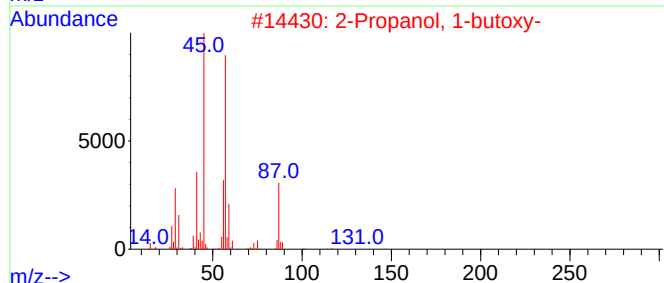
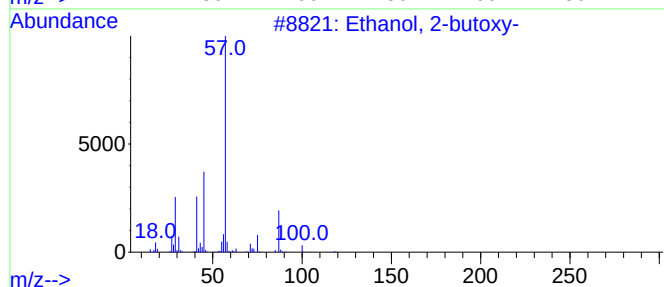
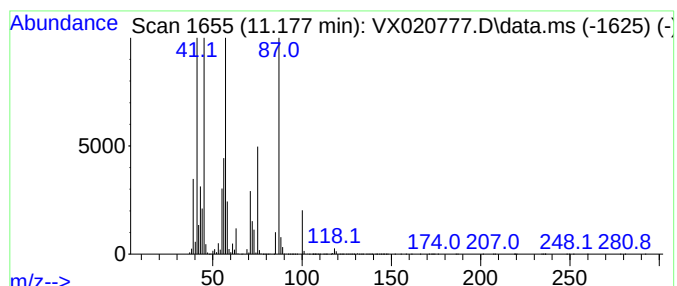
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Peak Number 3 Ethanol, 2-butoxy- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.177	50.00 ug/l	545492000	1,4-Dichlorobenzene-d4	12.067

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	58
2			2-Propanol, 1-butoxy-	132	C7H16O2	005131-66-8	50
3			Ethanol, 2-(2-butoxyethoxy)-, ac...	204	C10H20O4	000124-17-4	47
4			2-Ethyl-2-hydroxybutyric acid	132	C6H12O3	003639-21-2	38
5			Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	37



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Isopropyl Alcohol	2.611	79.1	ug/l	59390300	1	5.629	37561500	50.0
1-Butanol	7.306	50.0	ug/l	37561500	2	6.830	37561500	50.0
Ethanol, 2-butoxy-	11.177	50.0	ug/l	545492000	4	12.067	545492000	50.0