

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU121718\
Data File : VU028731.D
Acq On : 17 Dec 2018 19:32
Operator : MD/SY
Sample : J6440-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
COMP-3

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_U\METHOD\82U120618W.M
Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.453	81	87	163	rBV2	28208702	45533490	100.00%	81.061%
2	3.620	740	761	810	rBV	839890	2155446	4.73%	3.837%
3	5.955	1475	1487	1530	rVB	1779754	3437650	7.55%	6.120%
4	6.215	1546	1568	1585	rBV	1086339	2054098	4.51%	3.657%
5	7.562	1974	1987	2000	rBV	315866	550128	1.21%	0.979%
6	9.086	2449	2461	2479	rBV	273465	465506	1.02%	0.829%
7	9.372	2537	2550	2572	rVB	904186	1472199	3.23%	2.621%
8	13.739	3897	3908	3929	rVB	318667	503607	1.11%	0.897%

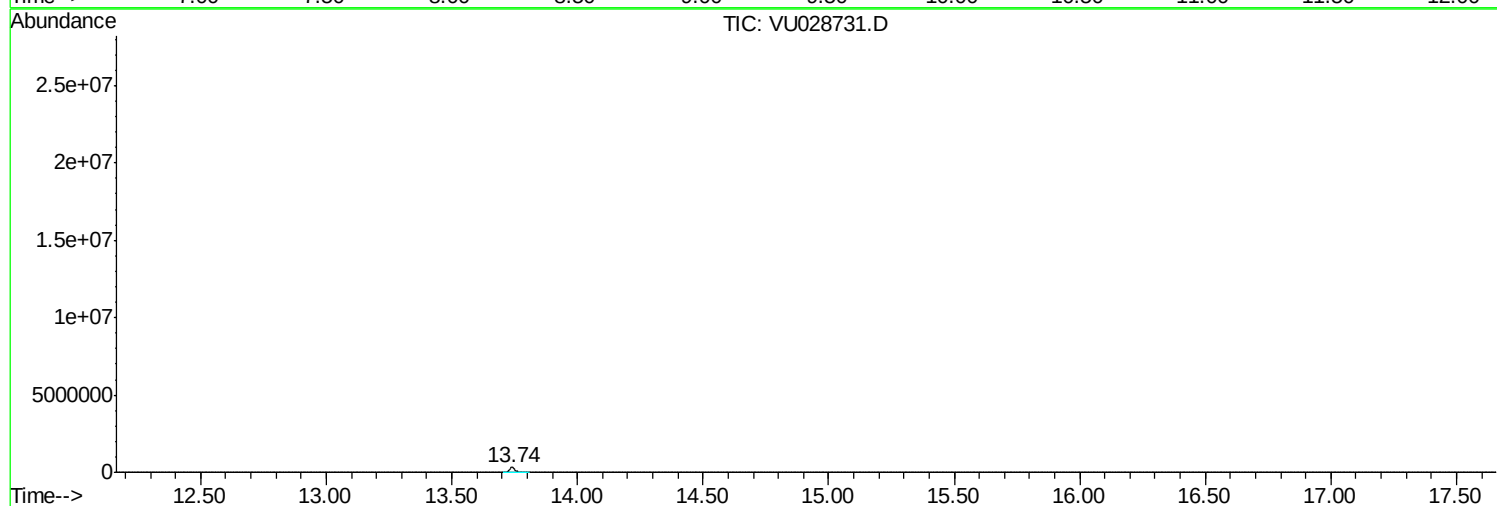
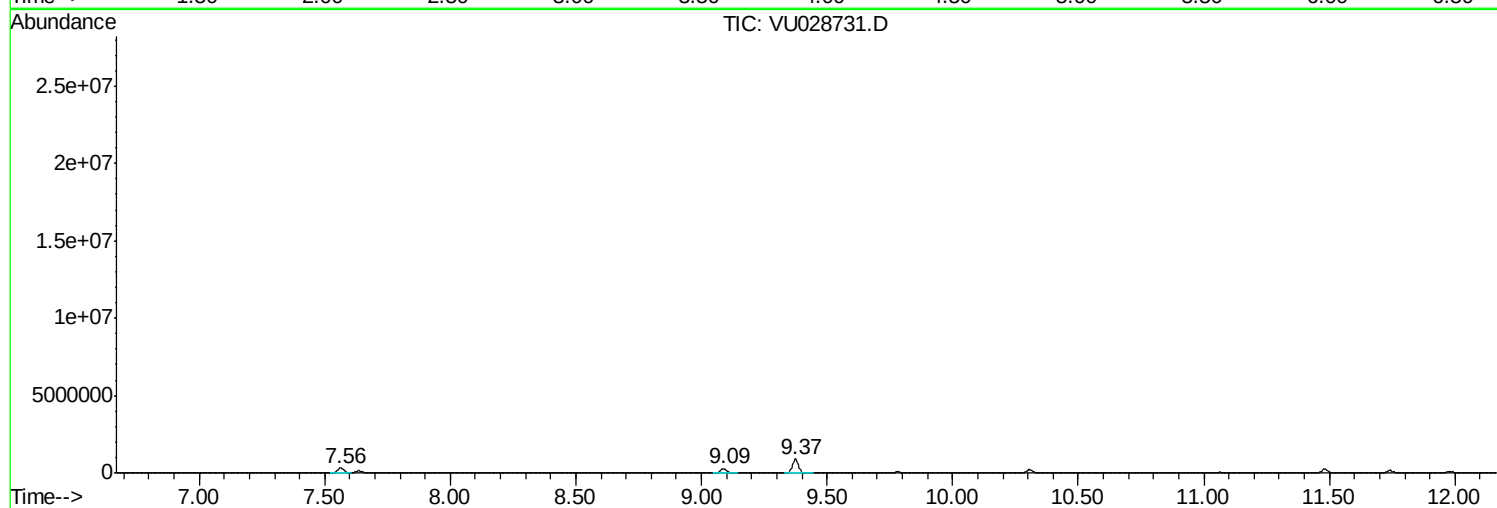
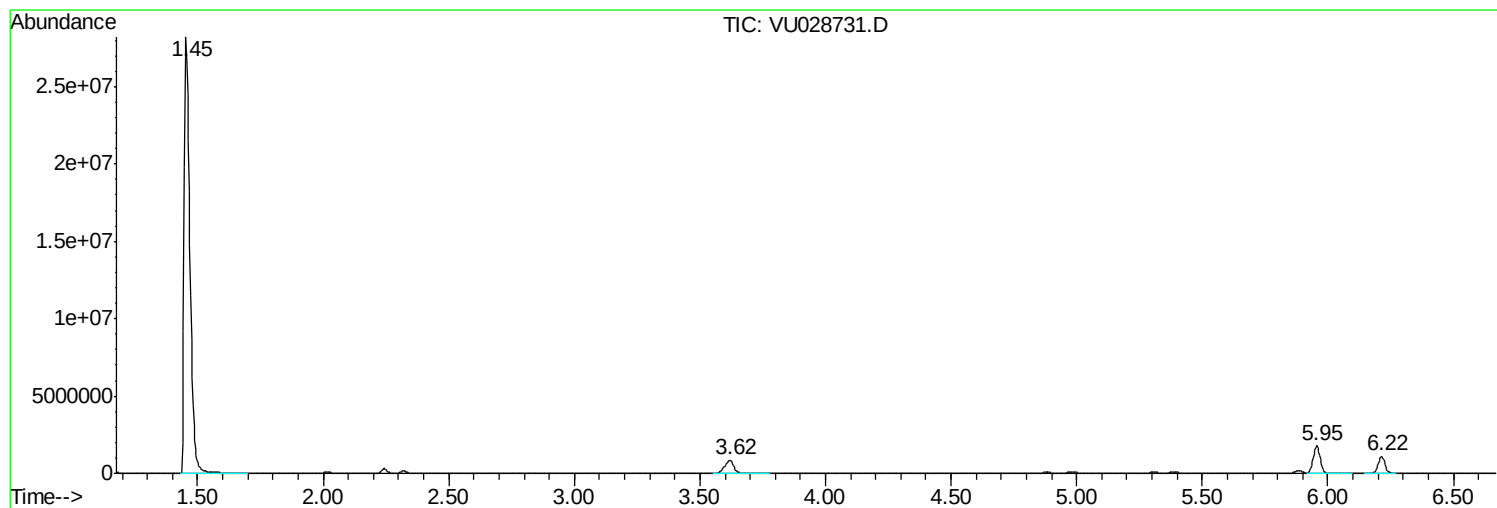
Sum of corrected areas: 56172124

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Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_U\METHOD\82U120618W.M
Quant Title : SW846 8260

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



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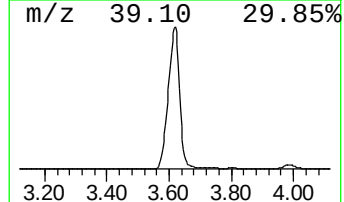
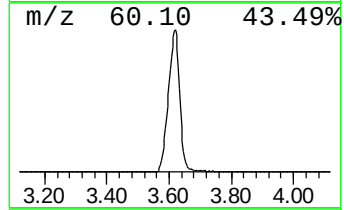
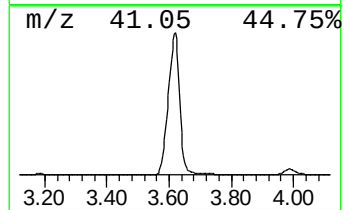
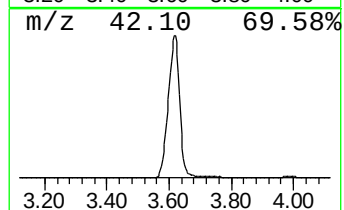
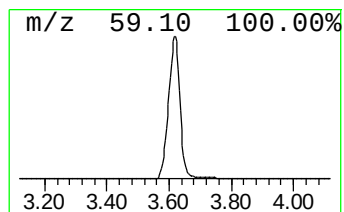
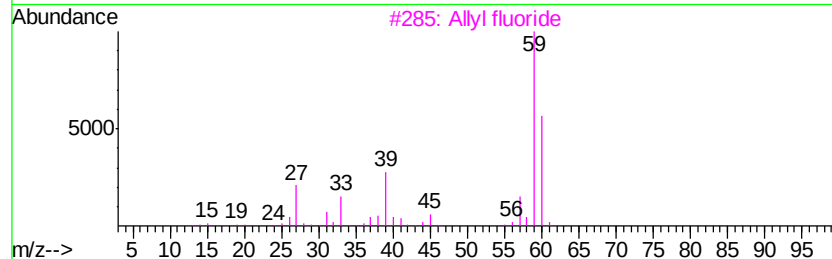
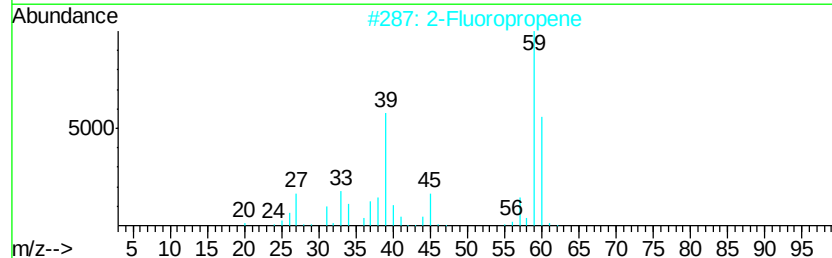
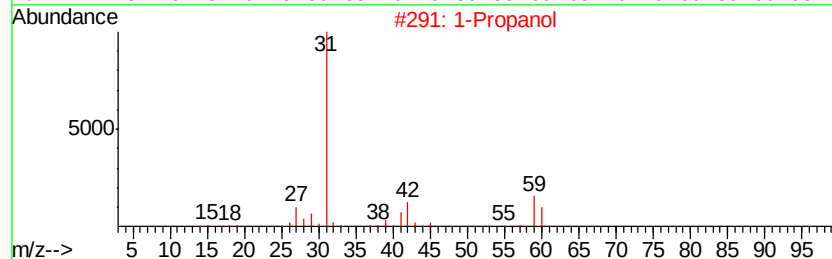
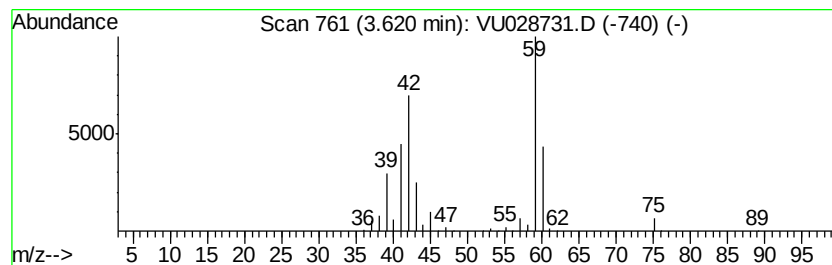
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 1-Propanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.62	31.35 ug/l	2155450	Pentafluorobenzene	4.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Propanol	60	C3H8O	000071-23-8	78
2		2-Fluoropropene	60	C3H5F	001184-60-7	53
3		Allyl fluoride	60	C3H5F	000818-92-8	42
4		Propanedioic acid	104	C3H4O4	000141-82-2	9
5		1-Propanamine, N,2-dimethyl-N-ni...	116	C5H12N2O	034419-76-6	9



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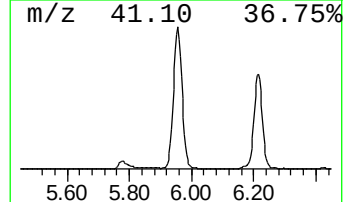
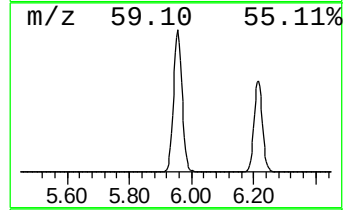
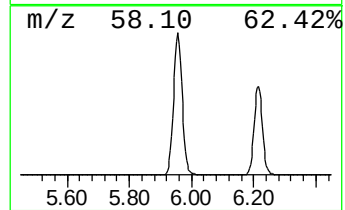
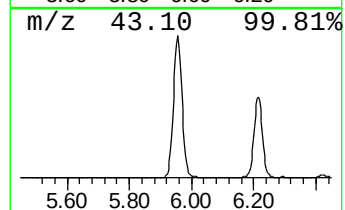
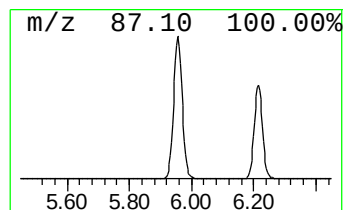
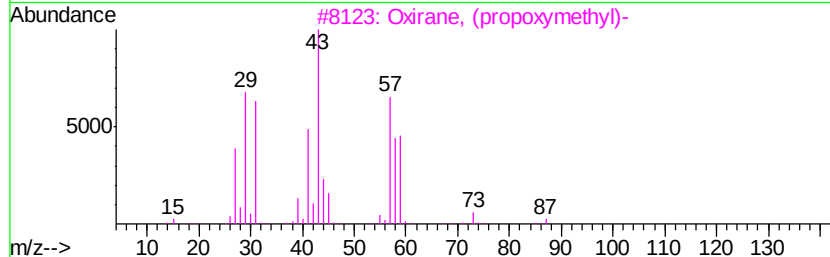
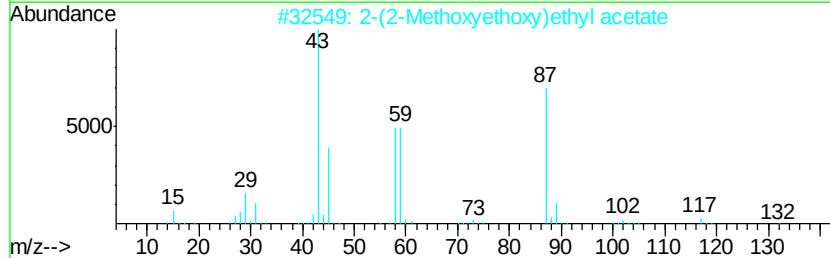
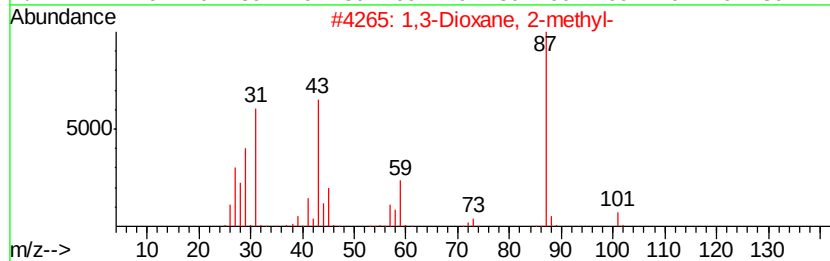
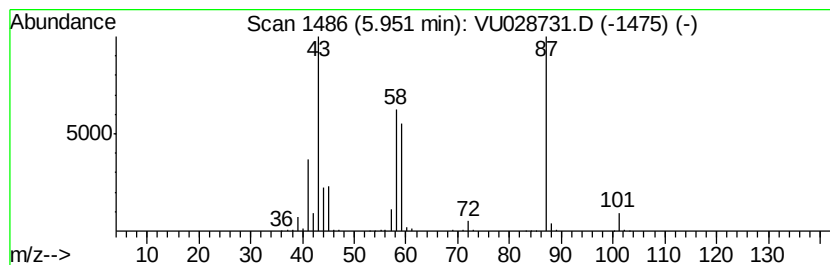
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Peak Number 3 1,3-Dioxane, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.95	50.00 ug/l	3437650	1,4-Difluorobenzene	5.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Dioxane, 2-methyl-	102	C5H10O2	000626-68-6	59
2		2-(2-Methoxyethoxy)ethyl acetate	162	C7H14O4	1000351-95-4	45
3		Oxirane, (propoxymethyl)-	116	C6H12O2	003126-95-2	37
4		Monomethyl malonate	118	C4H6O4	016695-14-0	33
5		Oxazolidin-2-one	87	C3H5NO2	000497-25-6	32



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Propanol	3.62	31.4	ug/l	2155450	1	4.98	3437650	50.0
1,3-Dioxane, 2-me...	5.95	50.0	ug/l	3437650	2	5.88	3437650	50.0