

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX122218\
 Data File : VX006775.D
 Acq On : 22 Dec 2018 16:36
 Operator : JC/MD
 Sample : J6533-19 10X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 COMP-5

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % 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\82X121818W.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	1.282	18	21	28	rBV	113391	121980	4.09%	0.676%
2	1.489	51	55	67	rBV	913808	1077913	36.18%	5.974%
3	1.611	67	75	80	rVB2	30688	65201	2.19%	0.361%
4	2.117	152	158	168	rBV	275607	446724	15.00%	2.476%
5	4.348	515	524	533	rBV3	22554	64187	2.15%	0.356%
6	4.696	571	581	593	rBV	117825	344281	11.56%	1.908%
7	5.500	702	713	726	rBV	352148	1011559	33.95%	5.606%
8	5.665	728	740	757	rVB	634609	1763904	59.21%	9.776%
9	6.067	795	806	816	rBV	373493	960429	32.24%	5.323%
10	6.860	927	936	953	rBV	919222	2158292	72.45%	11.962%
11	8.713	1233	1240	1249	rBV	1969044	2979139	100.00%	16.511%
12	10.109	1463	1469	1481	rBV	1884564	2571677	86.32%	14.253%
13	11.134	1632	1637	1644	rBV	1521729	1992221	66.87%	11.041%
14	12.079	1786	1792	1800	rBV	1983500	2443239	82.01%	13.541%
15	12.274	1820	1824	1829	rBV	27369	42327	1.42%	0.235%

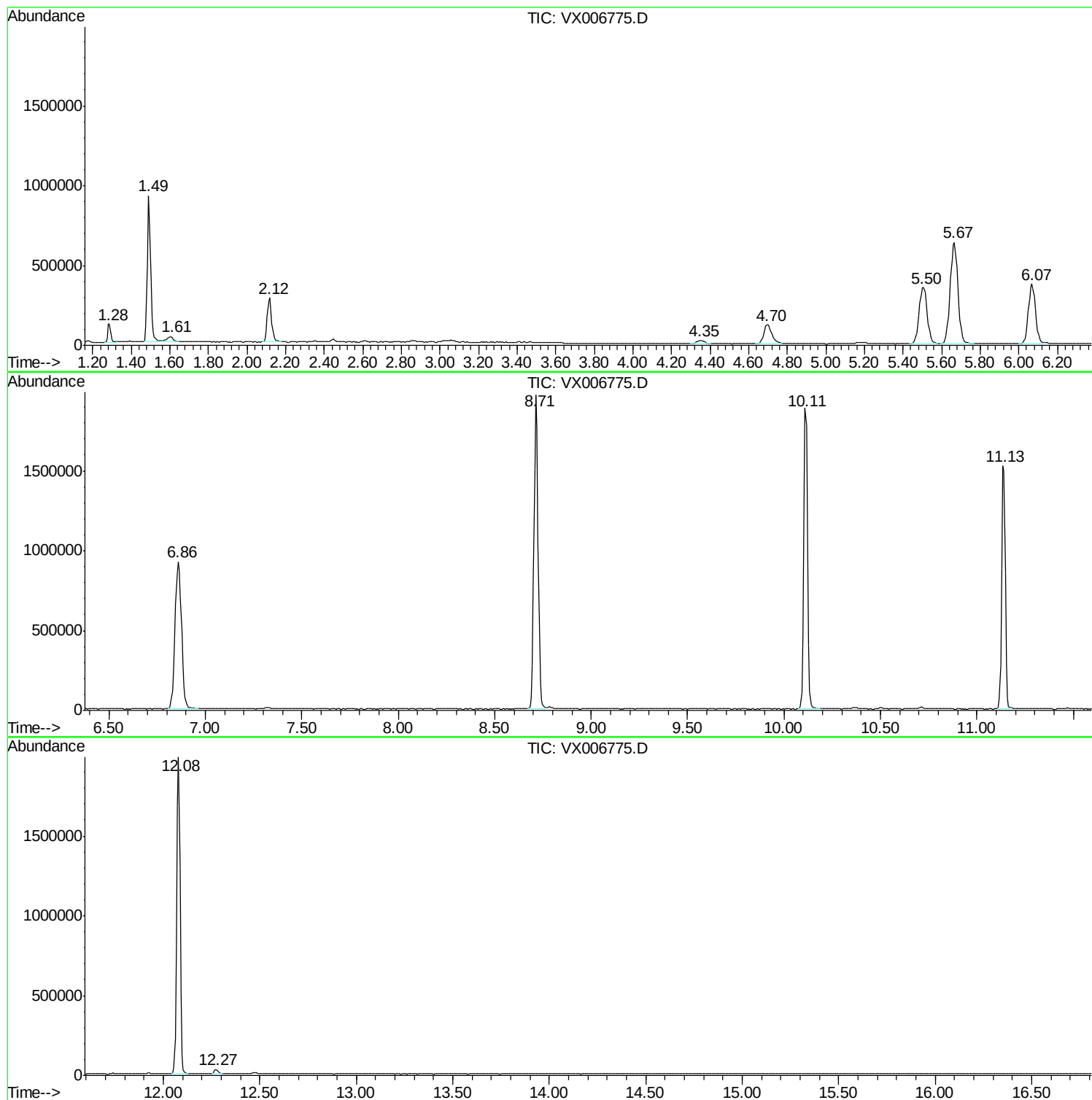
Sum of corrected areas: 18043073

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
Quant Title : SW846 8260

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



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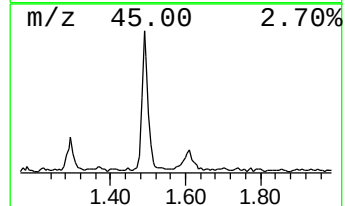
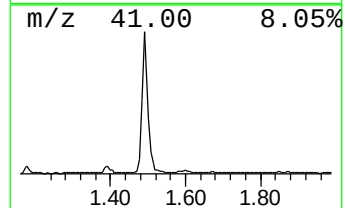
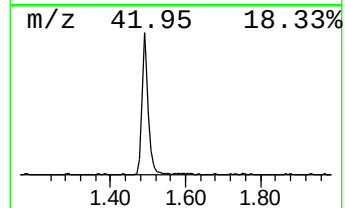
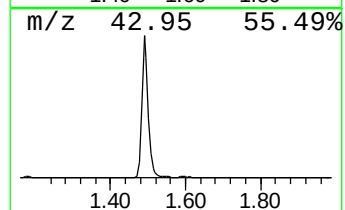
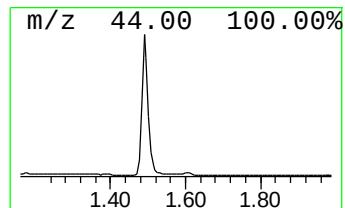
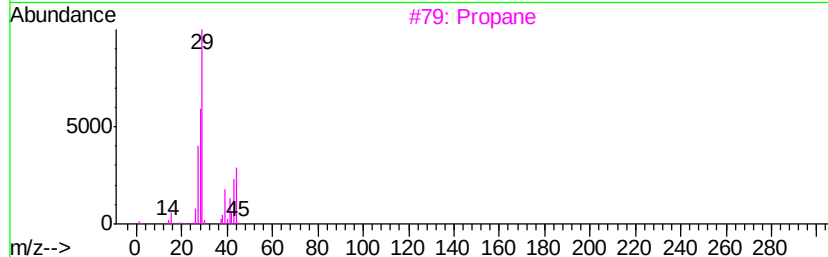
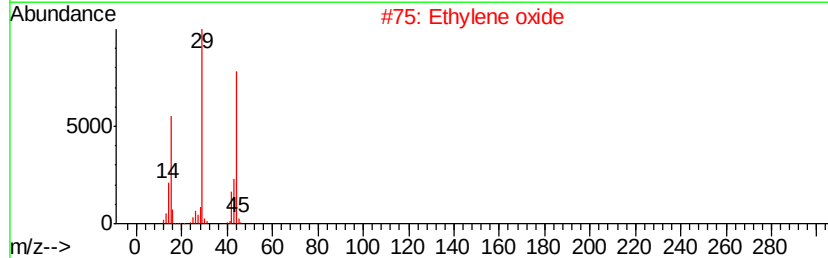
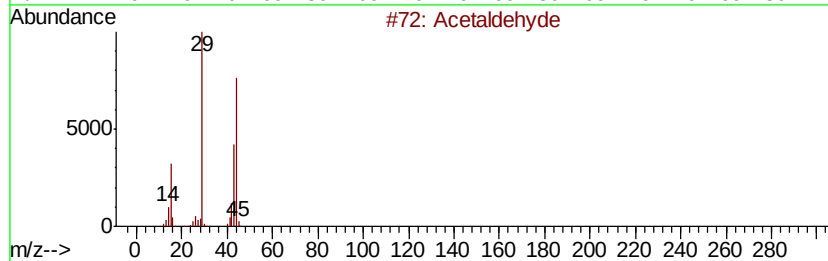
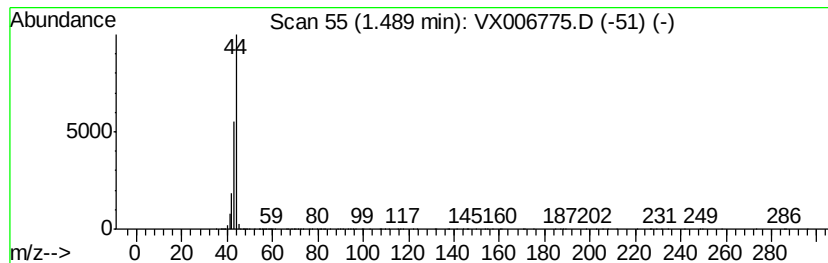
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Peak Number 1 Acetaldehyde Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.49	30.55 ug/l	1077910	Pentafluorobenzene	5.67

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetaldehyd		44	C2H4O	000075-07-0	56
2	Ethylene oxide		44	C2H4O	000075-21-8	5
3	Propane		44	C3H8	000074-98-6	4
4	Alanine		89	C3H7NO2	000056-41-7	4
5	1,2-Propanediamine		74	C3H10N2	000078-90-0	4



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Acetaldehyde	1.49	30.6	ug/l	1077910	1	5.67	1763900	50.0