

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN070618\
 Data File : VN049750.D
 Acq On : 6 Jul 2018 20:05
 Operator : MD\SY
 Sample : J3869-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TW-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_N\METHODS\82N062918W.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	1.654	3	20	61	rBV2	19077643	90578187	100.00%	67.289%
2	7.374	1777	1799	1829	rBV	1100191	2878163	3.18%	2.138%
3	7.593	1846	1867	1878	rBV	1161405	2934485	3.24%	2.180%
4	7.667	1879	1890	1922	rVB	1697872	4097243	4.52%	3.044%
5	8.030	1983	2003	2030	rBV	1195061	2871272	3.17%	2.133%
6	8.287	2042	2083	2085	rBV2	224444	1069006	1.18%	0.794%
7	8.586	2160	2176	2221	rVB	2645704	5700738	6.29%	4.235%
8	10.094	2629	2645	2694	rBV	4398434	8739073	9.65%	6.492%
9	11.409	3037	3054	3085	rBV	3321084	6140620	6.78%	4.562%
10	12.403	3349	3363	3387	rBV	2914139	5016950	5.54%	3.727%
11	13.348	3642	3657	3672	rBV	2778867	4585709	5.06%	3.407%

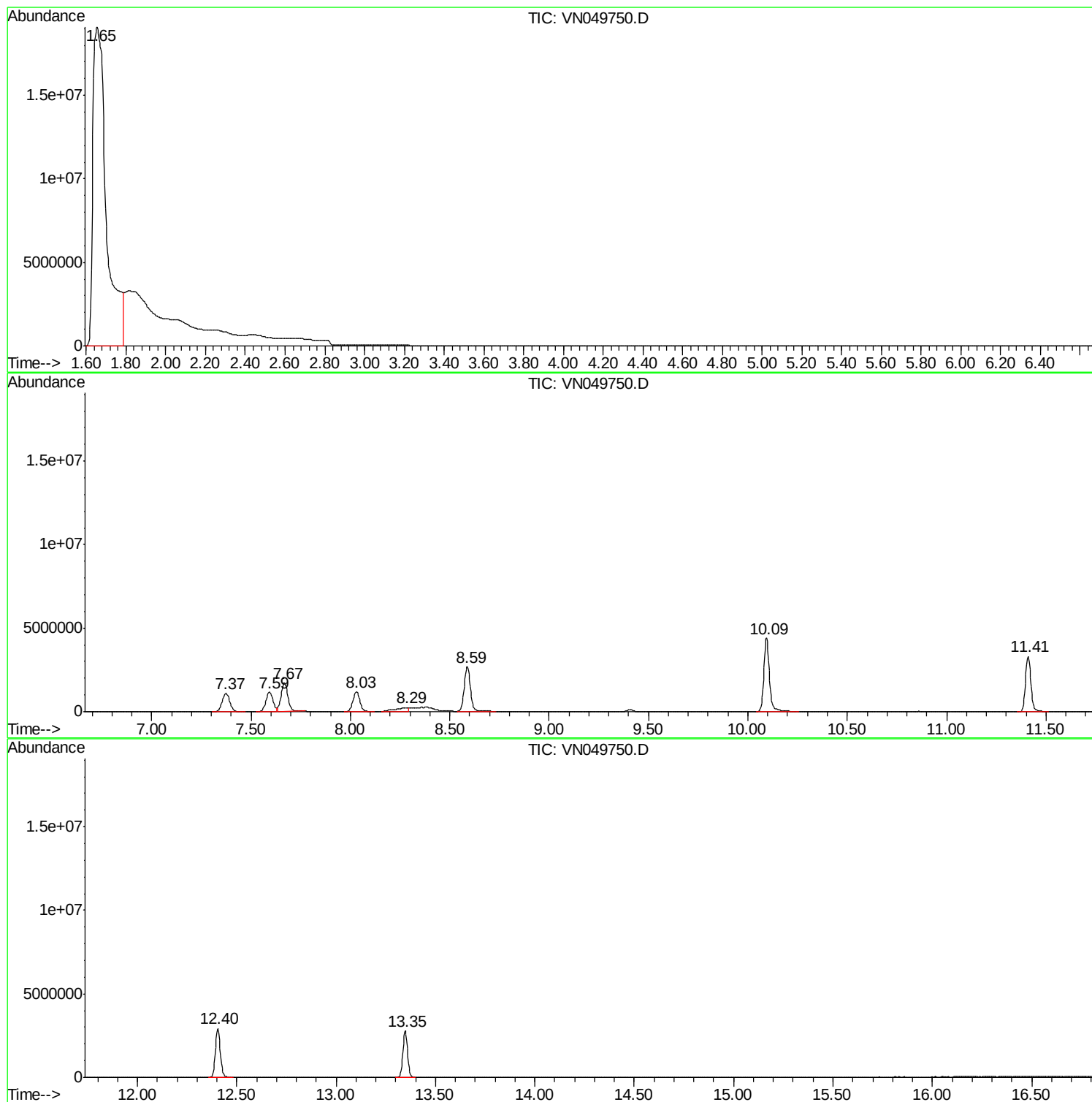
Sum of corrected areas: 134611446

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



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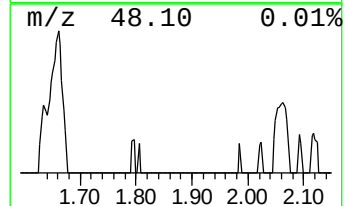
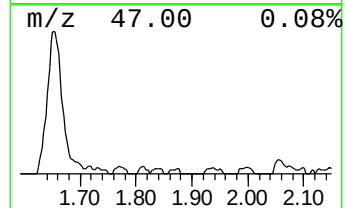
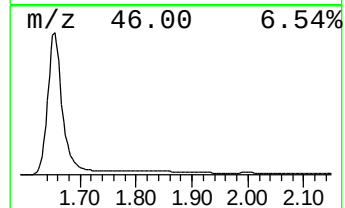
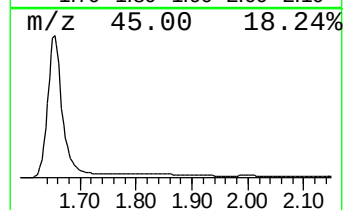
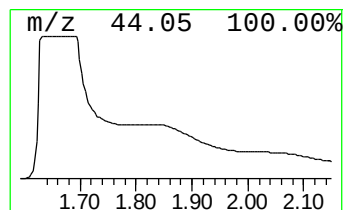
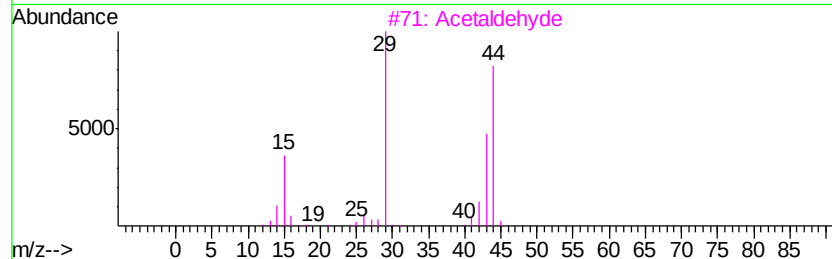
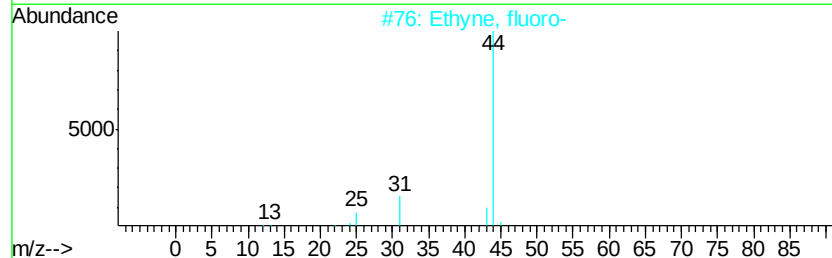
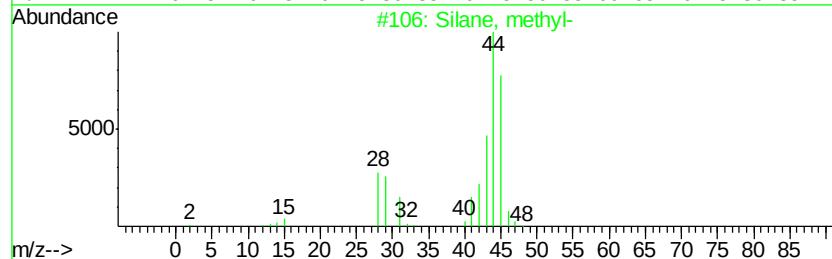
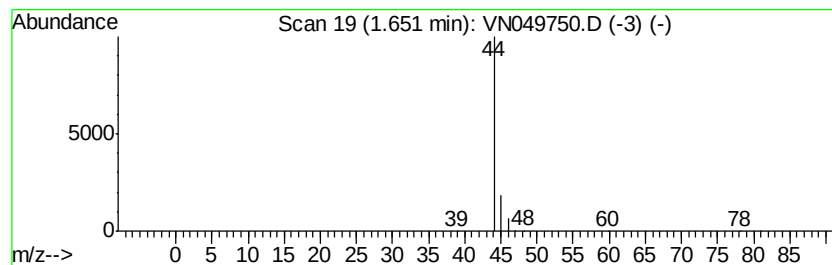
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Peak Number 1 unknown1.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.65	1105.36 ug/l	90578200	Pentafluorobenzene	7.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Silane, methyl-	46	CH6Si	000992-94-9	4
2		Ethyne, fluoro-	44	C2HF	002713-09-9	2
3		Acetaldehyde	44	C2H4O	000075-07-0	2
4		Nitrous oxide	44	N2O	010024-97-2	2
5		Carbon dioxide	44	CO2	000124-38-9	2



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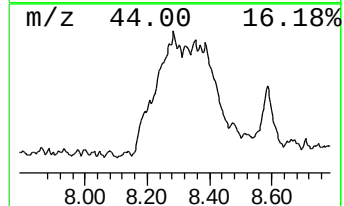
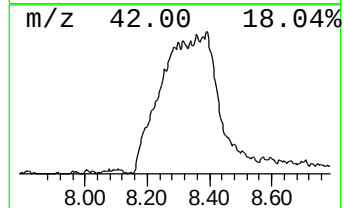
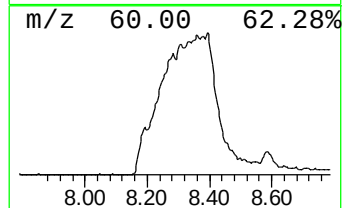
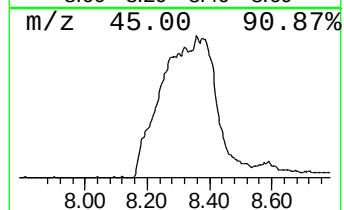
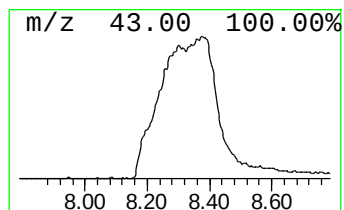
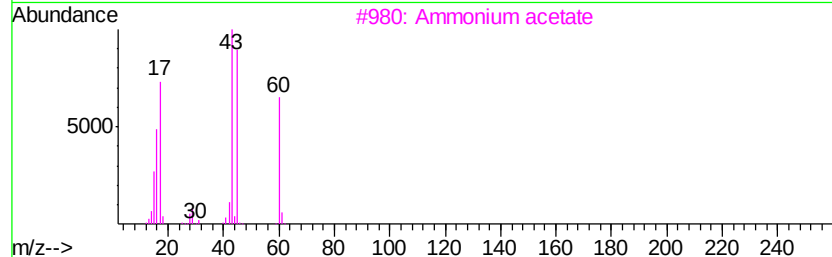
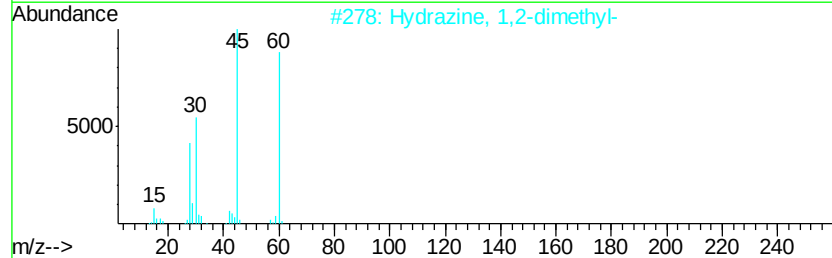
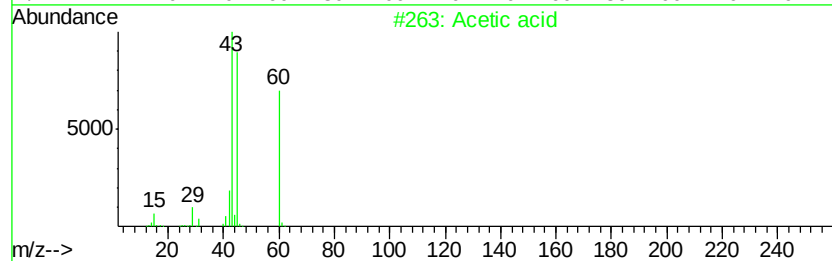
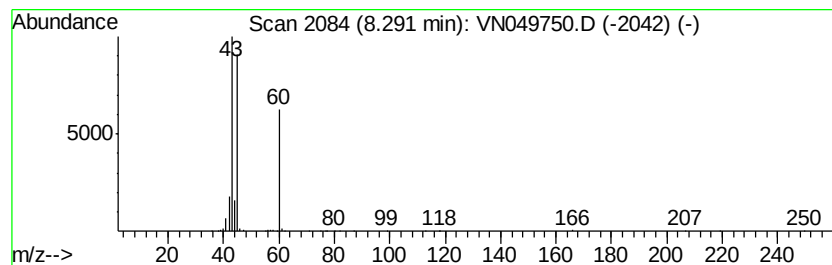
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Peak Number 2 Acetic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.29	9.38 ug/l	1069010	1,4-Difluorobenzene	8.59
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Acetic acid	60 C2H4O2	000064-19-7	90
2	Hydrazine, 1,2-dimethyl-	60 C2H8N2	000540-73-8	64
3	Ammonium acetate	77 C2H7NO2	000631-61-8	43
4	Cyclohexan-1,4,5-triol-3-one-1-c...	190 C7H10O6	1000128-45-5	40
5	Oxirane-2-carboxylic acid, methy...	102 C4H6O3	004538-50-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.65	1.65	1105.4	ug/l	90578200	1	7.67	4097240	50.0
Acetic acid	8.29	9.4	ug/l	1069010	2	8.59	5700740	50.0