

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN072718\
 Data File : VN050166.D
 Acq On : 28 Jul 2018 7:07
 Operator : MD\SY
 Sample : J4199-06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-A10

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\82N072418W.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	1.664	3	23	56	rBV	3818734	6037204	100.00%	25.186%
2	4.075	754	773	796	rBV	19782	63304	1.05%	0.264%
3	5.050	1053	1076	1104	rVB4	42591	159172	2.64%	0.664%
4	7.593	1848	1867	1879	rBV	502518	1270974	21.05%	5.302%
5	7.670	1879	1891	1924	rVB	721811	1724660	28.57%	7.195%
6	8.030	1984	2003	2033	rBV	515871	1226822	20.32%	5.118%
7	8.249	2037	2071	2073	rBV3	125932	452311	7.49%	1.887%
8	8.590	2161	2177	2207	rBV	1083673	2277566	37.73%	9.501%
9	10.094	2629	2645	2674	rBV	1889786	3533162	58.52%	14.739%
10	11.409	3040	3054	3075	rBV	1398644	2444953	40.50%	10.200%
11	12.403	3348	3363	3379	rBV	1135412	1948351	32.27%	8.128%
12	13.345	3643	3656	3674	rBV	1112810	1857948	30.77%	7.751%
13	13.544	3700	3718	3730	rBV	549645	910341	15.08%	3.798%
14	13.998	3849	3859	3868	rBV2	42638	64028	1.06%	0.267%

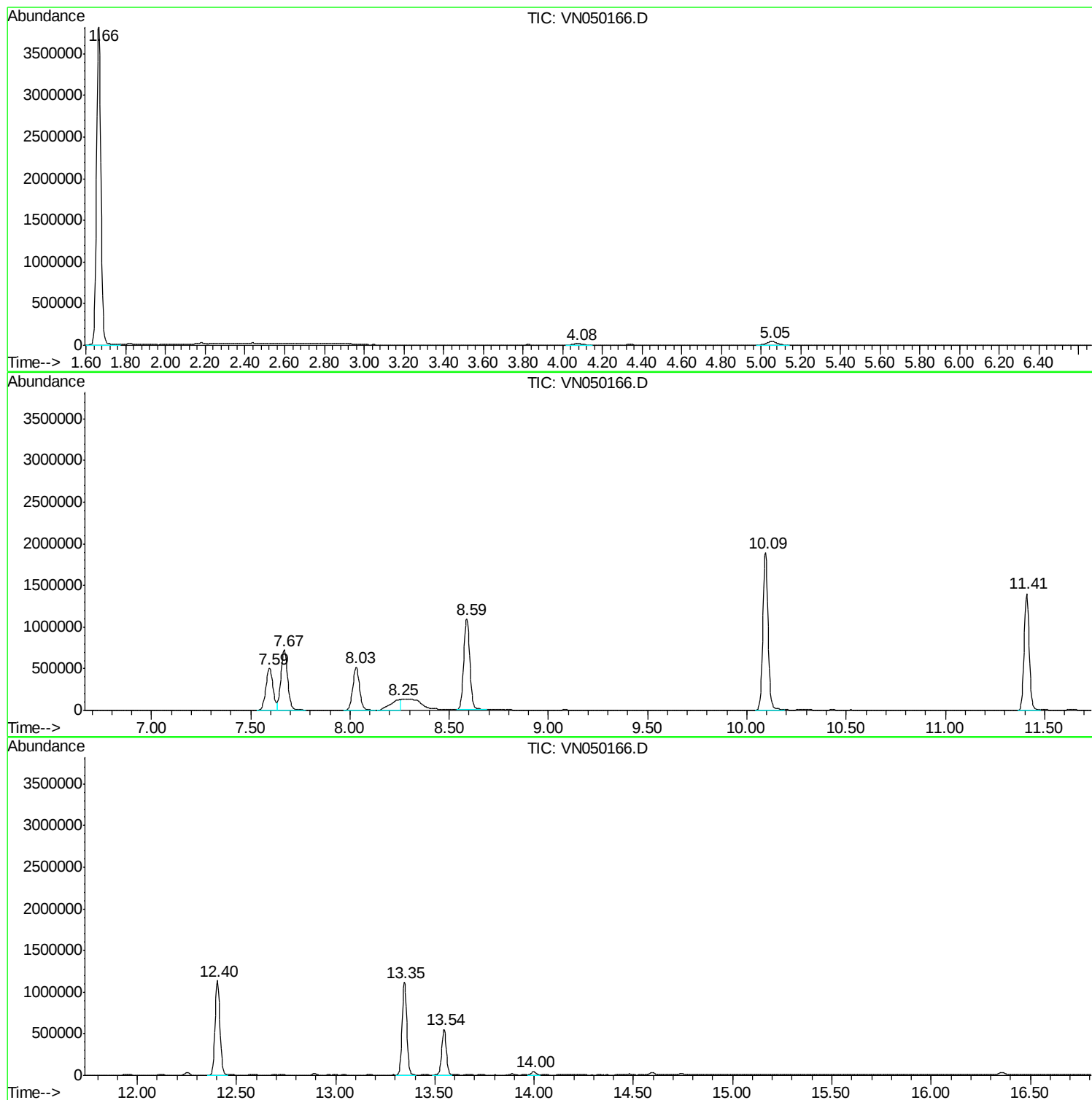
Sum of corrected areas: 23970796

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



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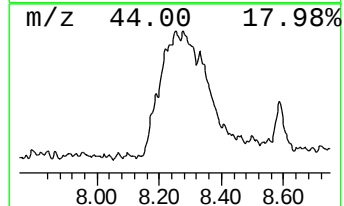
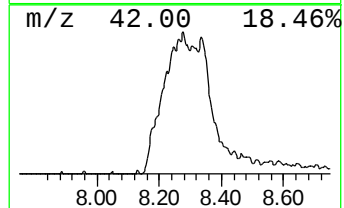
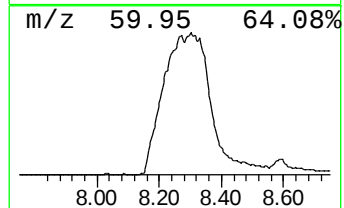
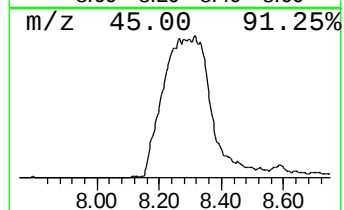
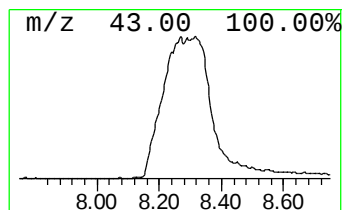
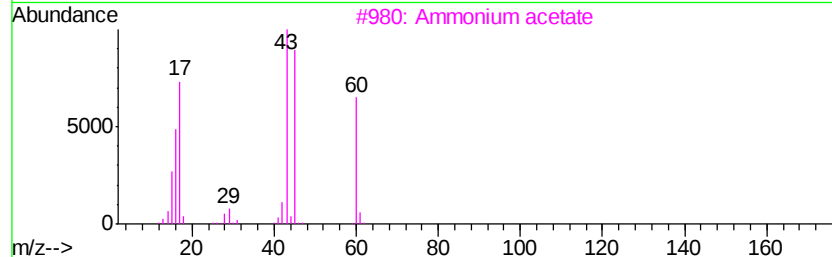
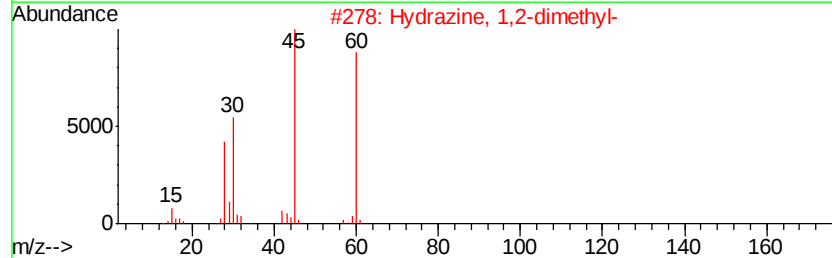
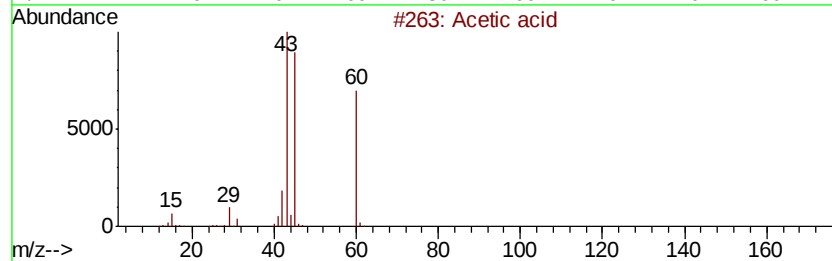
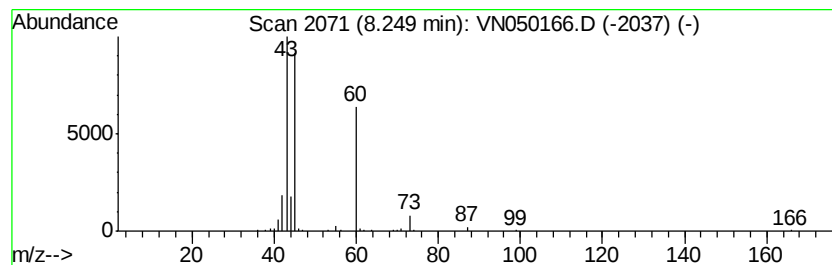
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Acetic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.25	9.93 ug/l	452311	1,4-Difluorobenzene	8.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acetic acid	60	C2H4O2	000064-19-7	86
2		Hydrazine, 1,2-dimethyl-	60	C2H8N2	000540-73-8	59
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		Methanamine, 1-methoxy-N-methyl-...	104	C3H8N2O2	039885-14-8	28



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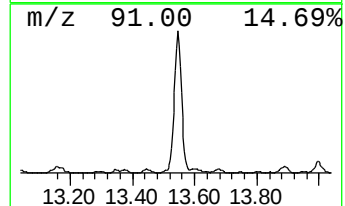
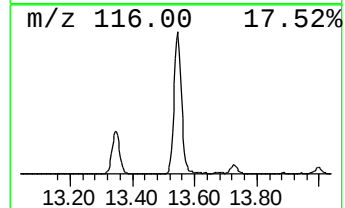
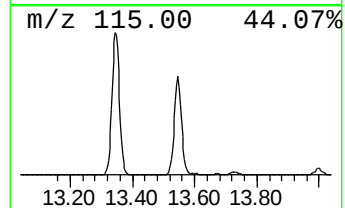
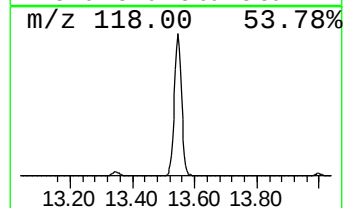
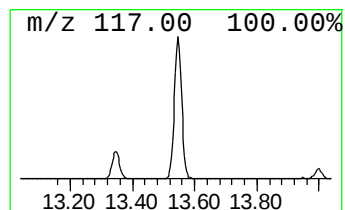
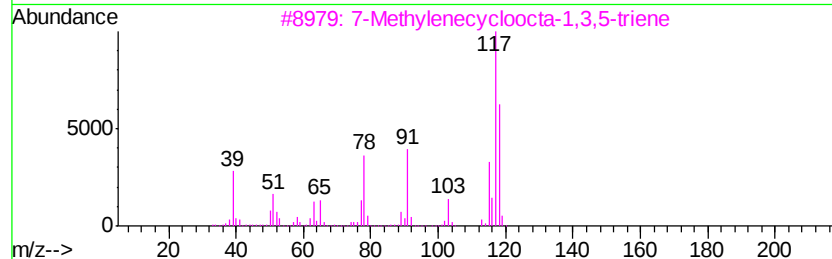
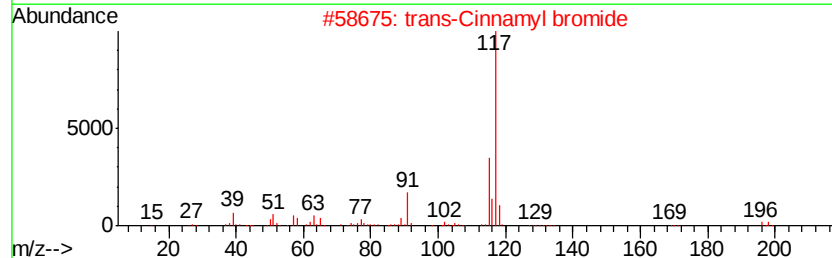
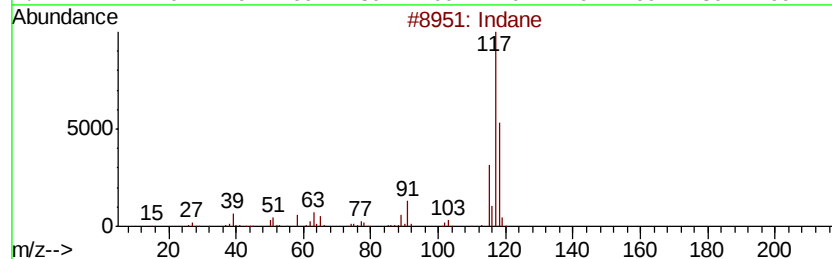
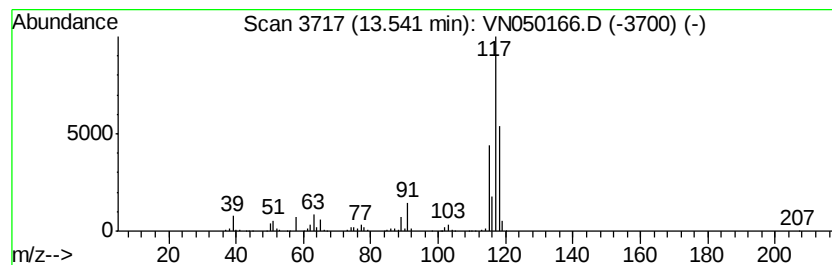
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Peak Number 3 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	24.50 ug/l	910341	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indane	118 C9H10	000496-11-7	95
2	trans-Cinnamyl bromide	196 C9H9Br	026146-77-0	72
3	7-Methylenecycloocta-1,3,5-triene	118 C9H10	002570-13-0	64
4	Benzene, 1-propenyl-	118 C9H10	000637-50-3	62
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118 C9H10	1000191-13-7	53



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Acetic acid	8.25	9.9	ug/l	452311	2	8.59	2277570	50.0
Indane	13.54	24.5	ug/l	910341	4	13.35	1857950	50.0