

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN071619\
 Data File : VN056757.D
 Acq On : 16 Jul 2019 10:43
 Operator : JC/SP
 Sample : K3814-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 LA119SW005-SW-190712

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_N\METHODS\82N071519W.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	3.822	623	637	653	rBV2	7214	19793	1.31%	0.255%
2	4.066	699	713	738	rBV3	7493	22132	1.47%	0.285%
3	7.590	1791	1809	1821	rBV	182586	456943	30.31%	5.876%
4	7.667	1821	1833	1855	rVB	302776	716921	47.55%	9.219%
5	8.027	1928	1945	1970	rBV	186912	431926	28.65%	5.554%
6	8.307	1994	2032	2034	rBV4	25333	114796	7.61%	1.476%
7	8.584	2103	2118	2155	rVB	503917	1057324	70.12%	13.597%
8	10.091	2571	2587	2618	rBV	795641	1507803	100.00%	19.390%
9	11.407	2982	2996	3025	rBV	782419	1400450	92.88%	18.009%
10	12.400	3287	3305	3322	rBV	610153	1031686	68.42%	13.267%
11	13.342	3585	3598	3616	rVB	612923	1016487	67.42%	13.072%

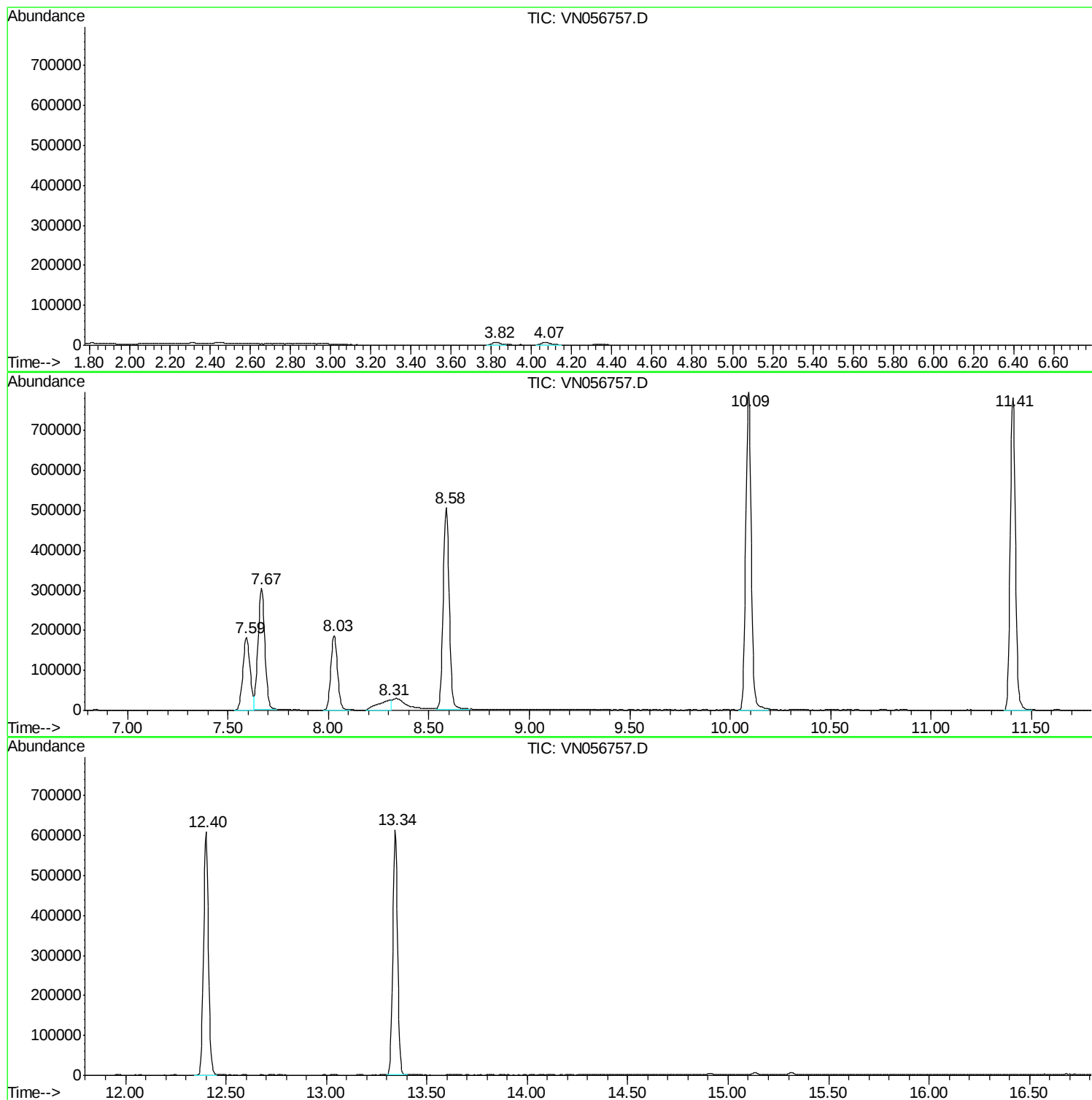
Sum of corrected areas: 7776261

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N071519W.M
Quant Title : SW846 8260

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



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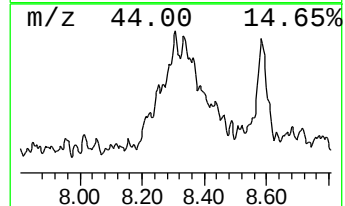
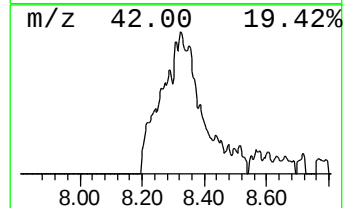
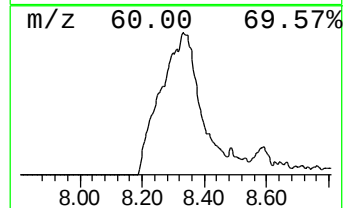
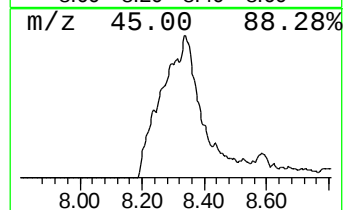
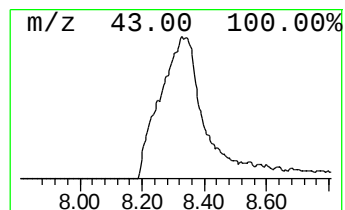
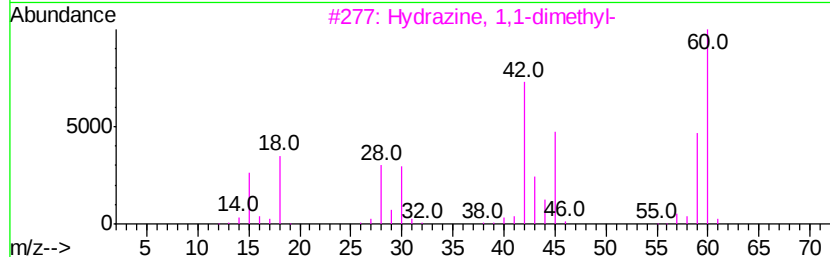
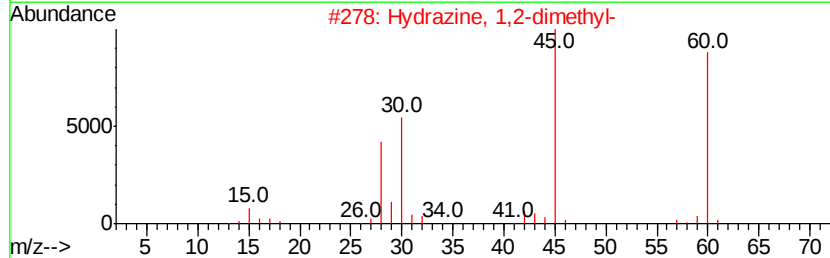
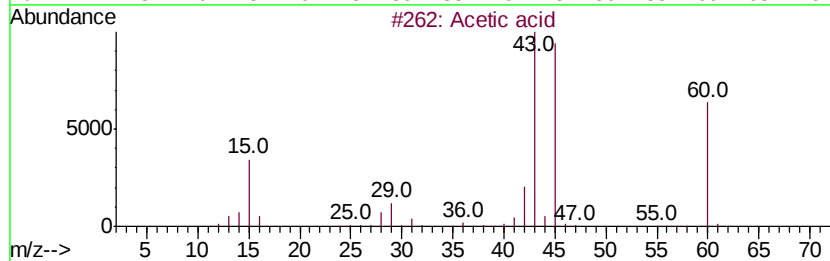
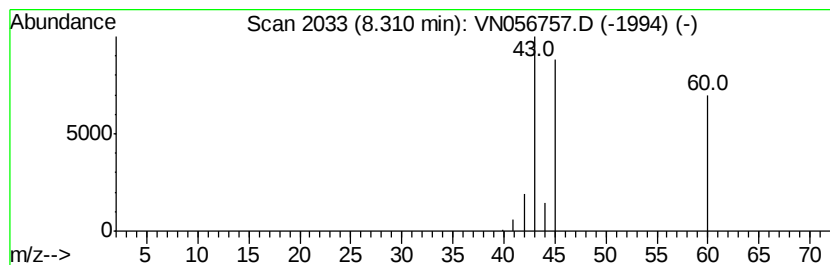
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N071519W.M
Quant Title : SW846 8260

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

Peak Number 1 Acetic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
8.31	5.43 ug/l	114796	1,4-Difluorobenzene	8.59		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid		60	C2H4O2	000064-19-7	90
2	Hvdrazine, 1,2-dimethyl-		60	C2H8N2	000540-73-8	5
3	Hvdrazine, 1,1-dimethyl-		60	C2H8N2	000057-14-7	5
4	Formamide		45	CH3NO	000075-12-7	4
5	Methane, nitroso-		45	CH3NO	000865-40-7	3



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Acetic acid	8.31	5.4	ug/l	114796	2	8.59	1057320	50.0