

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100419\
 Data File : VN058539.D
 Acq On : 4 Oct 2019 14:41
 Operator : JC/SP
 Sample : K5176-19 100X
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 30459-61

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\82N092719W.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	2.423	192	202	212	rBV6	17529	32794	1.40%	0.213%
2	6.799	1546	1563	1585	rBV2	25177	83159	3.56%	0.540%
3	7.574	1783	1804	1815	rBV	283062	699564	29.97%	4.544%
4	7.651	1815	1828	1849	rVB	559659	1331057	57.01%	8.646%
5	8.011	1920	1940	1961	rBV2	356295	827795	35.46%	5.377%
6	8.574	2099	2115	2141	rBV	805154	1702616	72.93%	11.059%
7	10.085	2571	2585	2603	rBV	1257401	2334586	100.00%	15.164%
8	10.291	2630	2649	2670	rBV2	127682	500290	21.43%	3.250%
9	11.407	2983	2996	3022	rVB	1118144	1978334	84.74%	12.850%
10	12.407	3295	3307	3317	rBV	892104	1613523	69.11%	10.481%
11	13.349	3588	3600	3619	rVB	1031891	1705594	73.06%	11.079%
12	13.895	3757	3770	3784	rVB	74420	181090	7.76%	1.176%
13	15.275	4191	4199	4213	rVB	80913	138125	5.92%	0.897%
14	16.027	4420	4433	4463	rBV	1054317	2103355	90.10%	13.662%
15	16.686	4630	4638	4651	rVB2	87471	163440	7.00%	1.062%

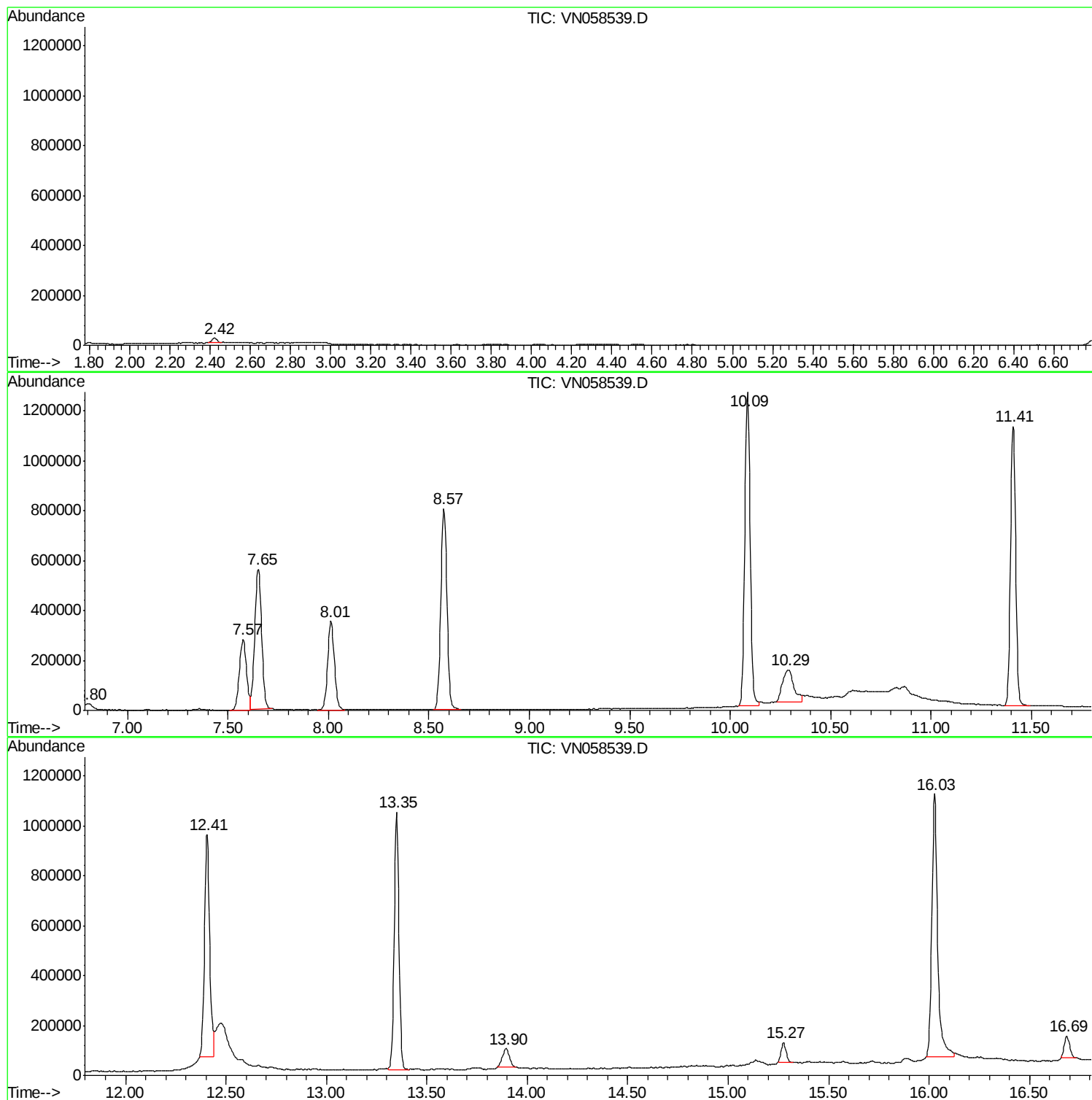
Sum of corrected areas: 15395322

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N092719W.M
Quant Title : SW846 8260

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



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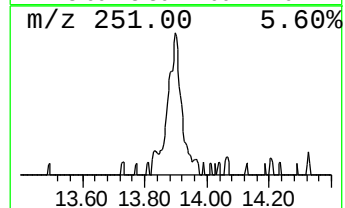
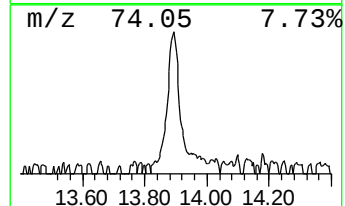
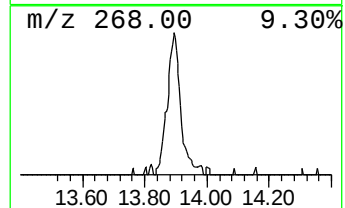
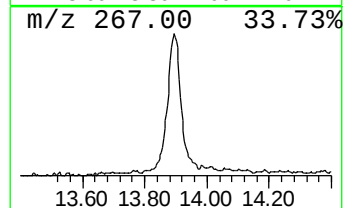
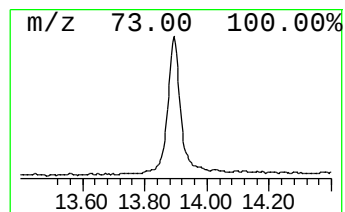
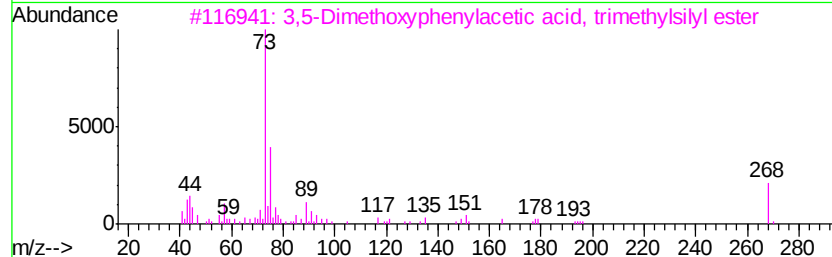
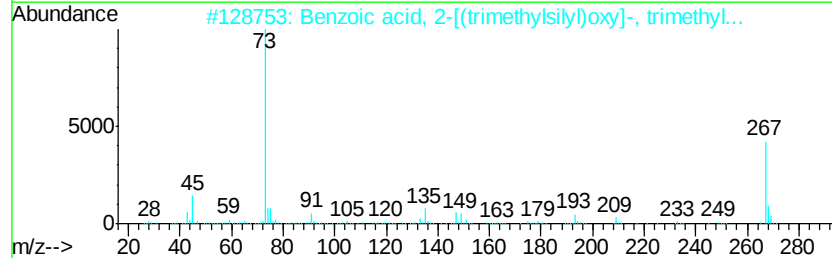
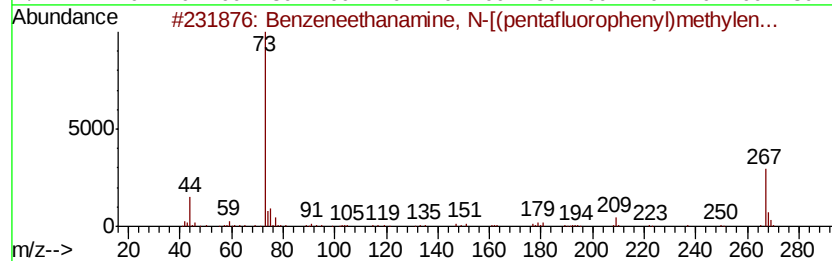
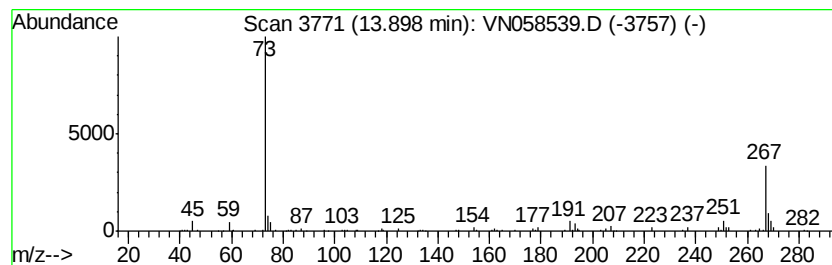
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Peak Number 2 Benzeneethanamine, N-[(pent... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.90	5.31 ug/l	181090	1,4-Dichlorobenzene-d4	13.35		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzeneethanamine, N-[(pentafluoro...	475	C21H26F5N02Si2	055429-85-1	50
2		Benzoic acid, 2-[(trimethylsilyl)...	282	C13H22O3Si2	003789-85-3	45
3		3,5-Dimethoxyphenylacetic acid, ...	268	C13H20O4Si	1000071-82-4	27
4		4-Methylcatechol, bis(trimethyls...	268	C13H24O2Si2	1000332-79-9	25
5		1-Boraindane, 3-methyl-1-[1-(tri...	266	C17H23BSi	190316-36-0	16



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzeneethanamine...	13.90	5.3	ug/l	181090	4	13.35	1705590	50.0