

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100419\
 Data File : VN058547.D
 Acq On : 4 Oct 2019 17:37
 Operator : JC/SP
 Sample : K5169-09
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-4R_100219

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	4.040	682	705	719	rBV	30713	90677	3.72%	0.615%
2	7.574	1785	1804	1815	rBV	275215	687500	28.19%	4.666%
3	7.648	1815	1827	1847	rVB	541565	1277950	52.40%	8.672%
4	8.011	1922	1940	1962	rBV	339911	796326	32.65%	5.404%
5	8.574	2098	2115	2148	rBV	781099	1657544	67.97%	11.248%
6	10.085	2571	2585	2605	rBV	1312426	2438697	100.00%	16.550%
7	10.291	2631	2649	2673	rBV2	104569	409222	16.78%	2.777%
8	11.407	2984	2996	3019	rVB	1078727	1868387	76.61%	12.679%
9	12.406	3295	3307	3318	rBV	833549	1479674	60.67%	10.041%
10	13.349	3588	3600	3615	rBV	938484	1543022	63.27%	10.471%
11	13.892	3756	3769	3788	rVB2	61961	156608	6.42%	1.063%
12	15.275	4190	4199	4211	rBV	71268	124864	5.12%	0.847%
13	16.027	4420	4433	4463	rBV	982319	2022629	82.94%	13.726%
14	16.686	4629	4638	4651	rVB2	91914	182603	7.49%	1.239%

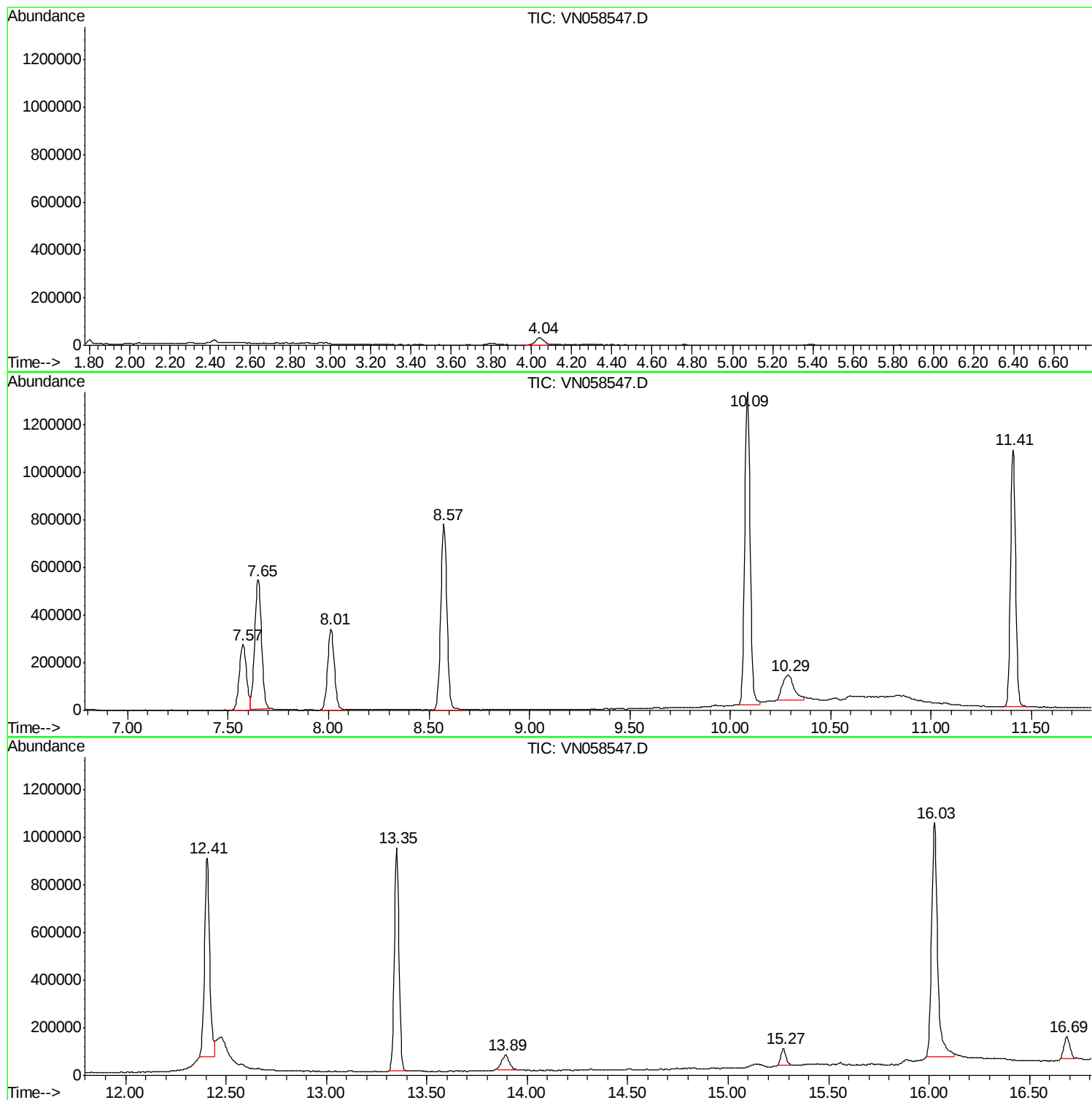
Sum of corrected areas: 14735703

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



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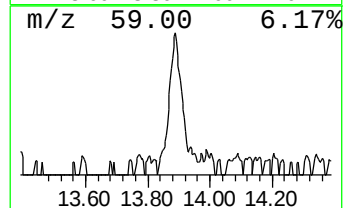
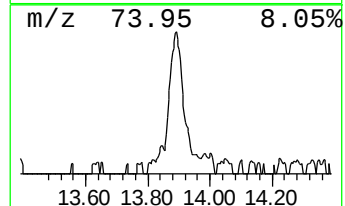
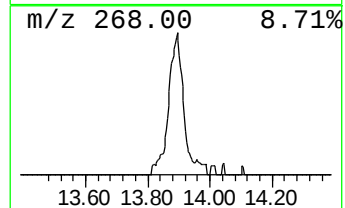
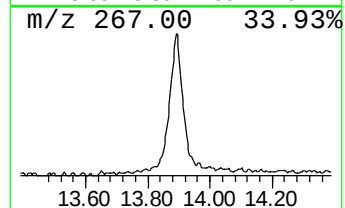
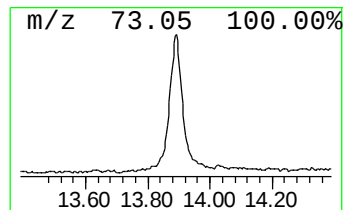
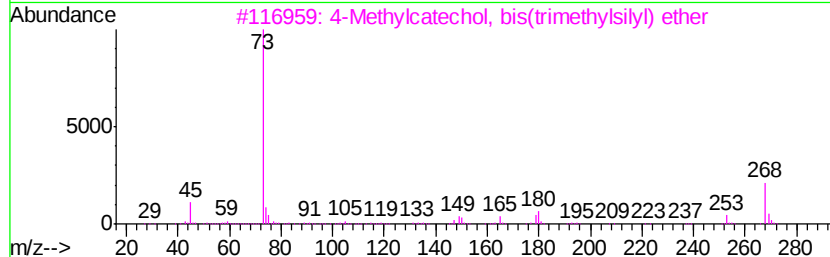
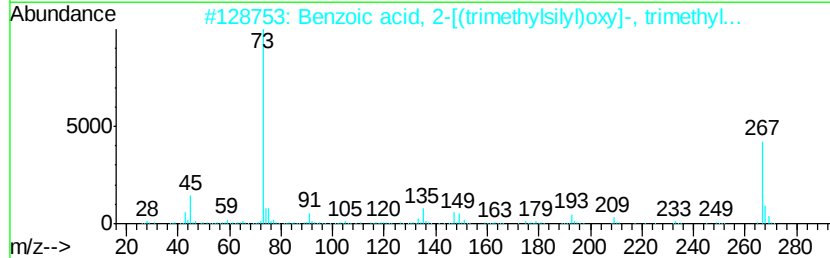
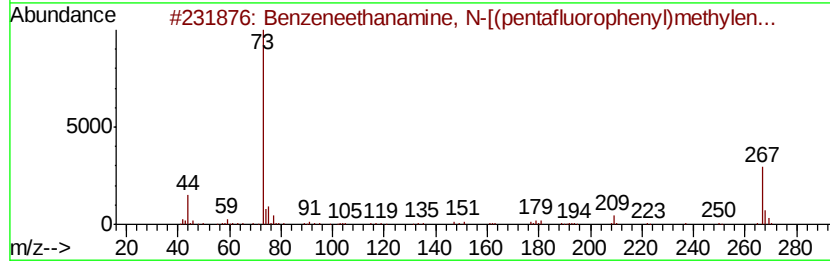
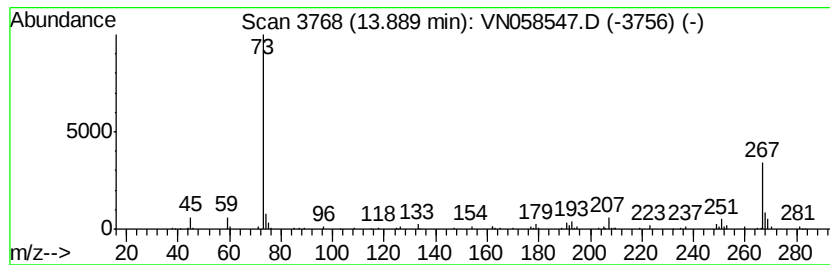
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzeneethanamine, N-[(pent... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.89	5.07 ug/l	156608	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		4-Methylcatechol, bis(trimethyls...	268	C13H24O2Si2	1000332-79-9	25
4		1-(3-Chlorophenyl)piperazine, 4-...	268	C13H21ClN2Si	1000373-99-3	25
5		11H-Dibenzo[b,e][1,4]diazepin-11...	267	C16H17N3O	013450-73-2	9



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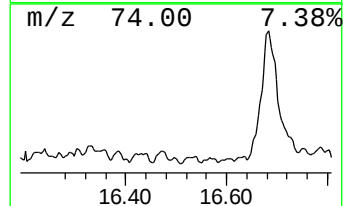
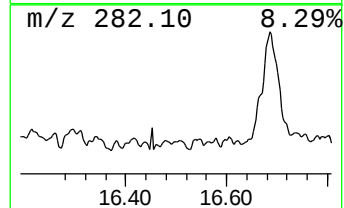
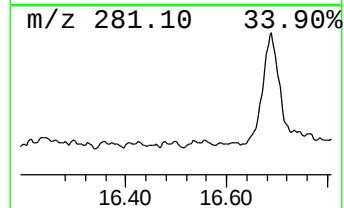
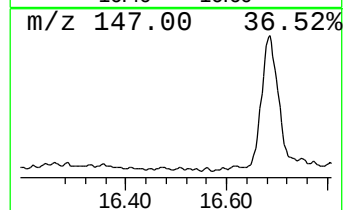
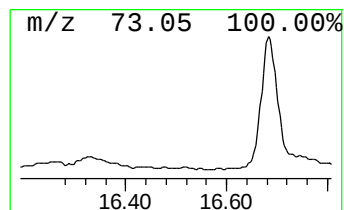
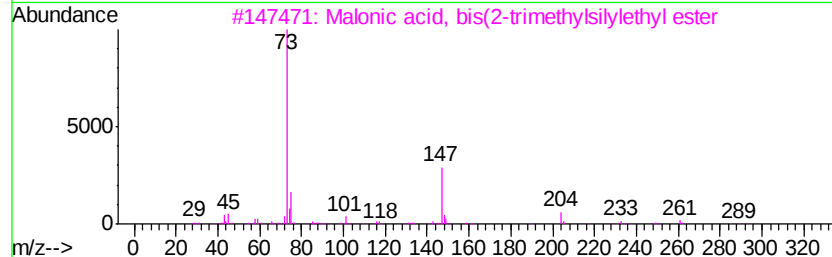
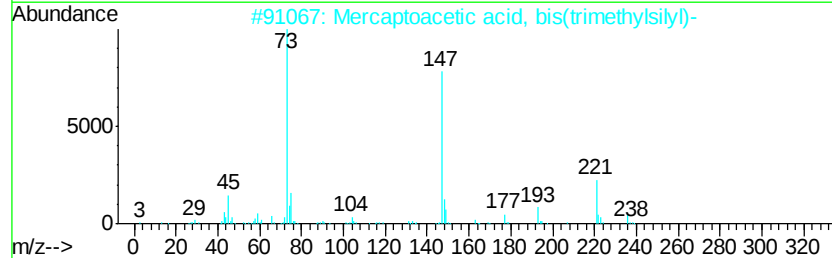
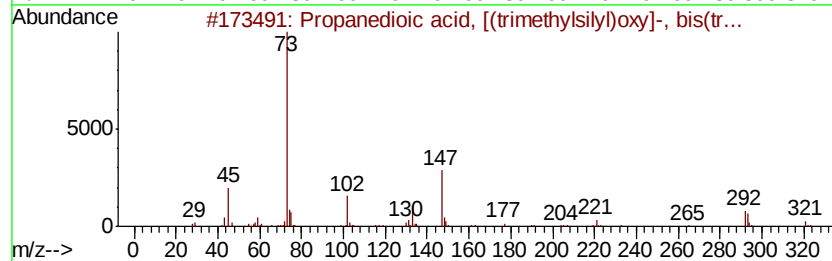
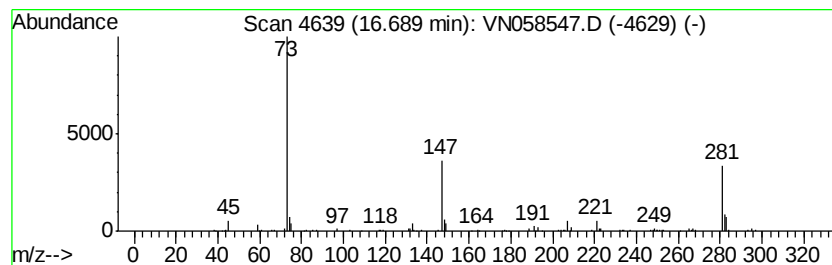
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Peak Number 4 unknown16.69 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.69	5.92 ug/l	182603	1,4-Dichlorobenzene-d4	13.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanedioic acid, [(trimethylsilyl)oxy]-, bis(tri...	336	C12H28O5Si3	038165-93-4	23
2		Mercaptoacetic acid, bis(trimethylsilyl)oxy]-, bis(tri...	236	C8H20O2SSi2	006398-62-5	17
3		Malonic acid, bis(2-trimethylsilyl)oxy]-, bis(tri...	304	C13H28O4Si2	090744-45-9	17
4		2-Hexenedioic acid, bis(trimethylsilyl)oxy]-, bis(tri...	288	C12H24O4Si2	055494-10-5	17
5		Ethanedioic acid, bis(trimethylsilyl)oxy]-, bis(tri...	234	C8H18O4Si2	018294-04-7	17



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzeneethanamine...	13.89	5.1	ug/l	156608	4	13.35	1543020	50.0
unknown16.69	16.69	5.9	ug/l	182603	4	13.35	1543020	50.0