

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW090219\
 Data File : VW012522.D
 Acq On : 03 Sep 2019 02:23
 Operator : SY/VA
 Sample : K4544-08
 Misc : 5.08G/5ML/MSVOA W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 T16-17-B3-(2.25)

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_W\METHOD\82W082319S.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	2.147	34	40	47	rBV	13295	25157	2.44%	0.543%
2	4.903	480	492	504	rBV2	14757	50950	4.94%	1.099%
3	5.933	651	661	671	rBV4	5518	16100	1.56%	0.347%
4	7.531	914	923	935	rBV5	3706	11229	1.09%	0.242%
5	7.878	969	980	985	rBV	154447	366584	35.55%	7.906%
6	7.945	985	991	1001	rVB	299011	665287	64.52%	14.348%
7	8.305	1041	1050	1062	rBV	163044	363504	35.25%	7.840%
8	8.841	1127	1138	1150	rBV	426294	848690	82.30%	18.304%
9	10.323	1371	1381	1389	rBV	591862	1031186	100.00%	22.240%
10	10.701	1434	1443	1452	rBV2	9990	19512	1.89%	0.421%
11	11.627	1588	1595	1606	rBV	335966	592122	57.42%	12.770%
12	12.615	1750	1757	1769	rVB2	185493	316646	30.71%	6.829%
13	13.554	1905	1911	1920	rBV	132736	221927	21.52%	4.786%
14	13.645	1921	1926	1932	rVB3	6446	10818	1.05%	0.233%
15	14.072	1991	1996	2002	rVB	17704	27270	2.64%	0.588%
16	14.523	2063	2070	2077	rVB3	7446	13906	1.35%	0.300%
17	14.724	2098	2103	2107	rVB	14071	19744	1.91%	0.426%
18	15.493	2224	2229	2235	rBV	8364	13822	1.34%	0.298%
19	15.816	2271	2282	2290	rBV7	9624	22250	2.16%	0.480%

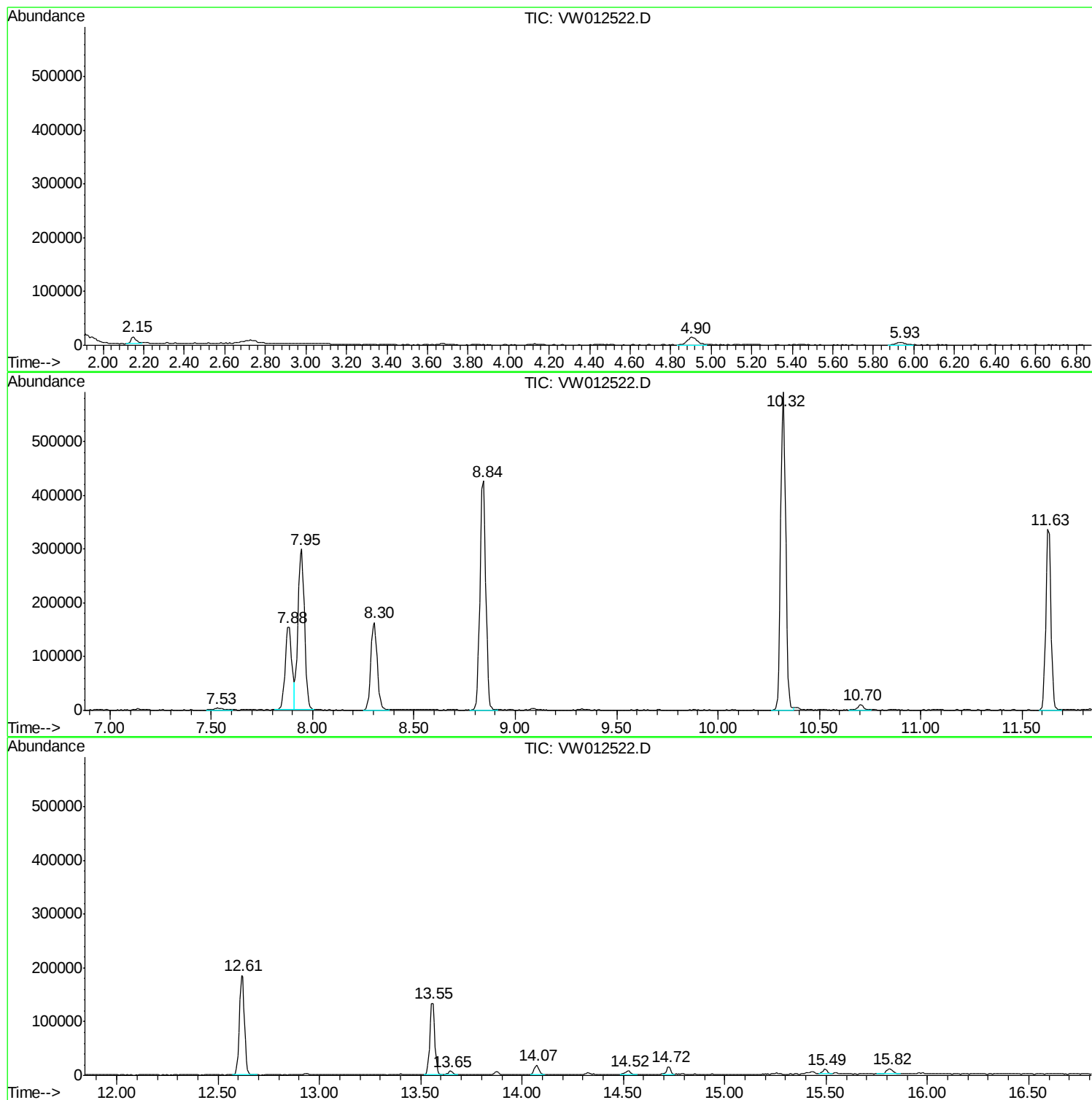
Sum of corrected areas: 4636704

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W082319S.M
Quant Title : SW846 8260

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



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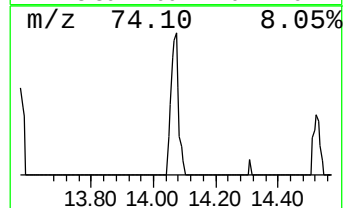
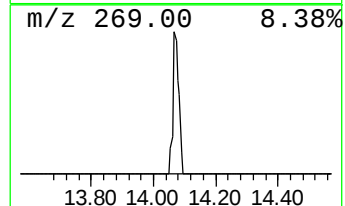
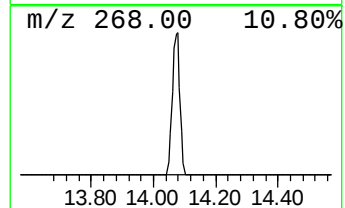
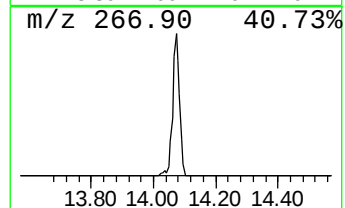
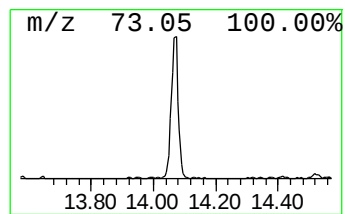
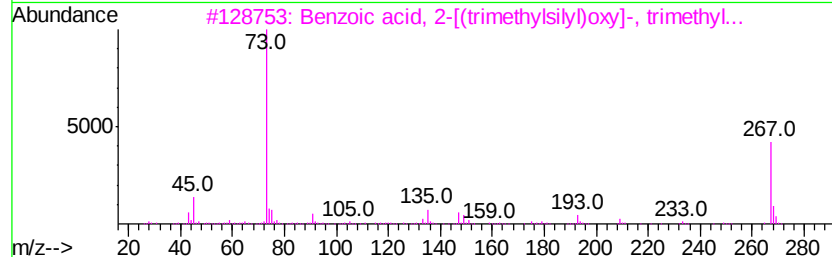
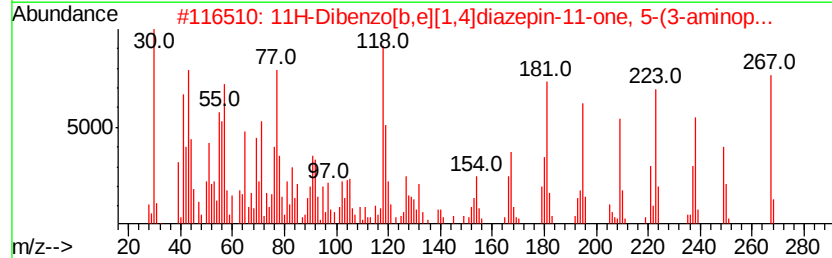
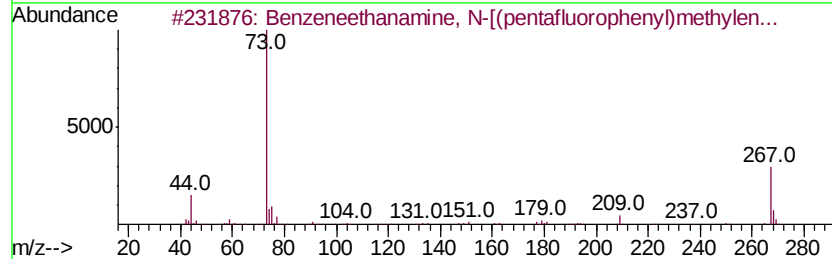
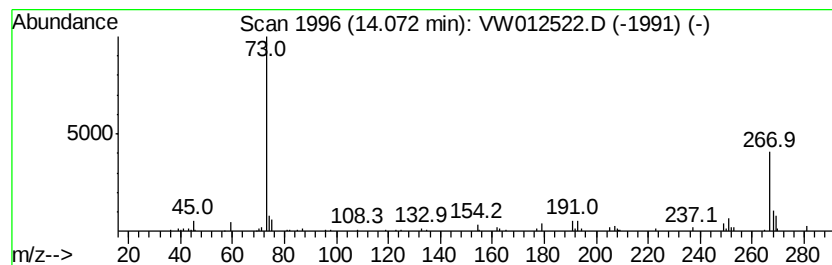
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzeneethanamine, N-[(pent... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.07	6.14 ug/l	27270	1,4-Dichlorobenzene-d4	13.56		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzeneethanamine, N-[(pentafluoro...	475	C21H26F5N02Si2	055429-85-1	50
2		11H-Dibenzo[b,e][1,4]diazepin-11...	267	C16H17N3O	013450-73-2	46
3		Benzoic acid, 2-[(trimethylsilyl)...	282	C13H22O3Si2	003789-85-3	38
4		4-Methylcatechol, bis(trimethyls...	268	C13H24O2Si2	1000332-79-9	22
5		3,5-Dimethoxyphenylacetic acid, ...	268	C13H20O4Si	1000071-82-4	14



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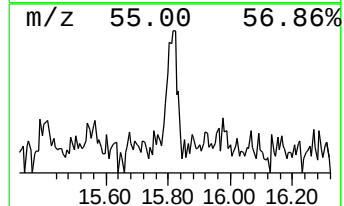
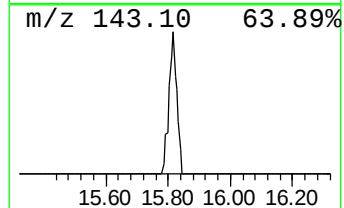
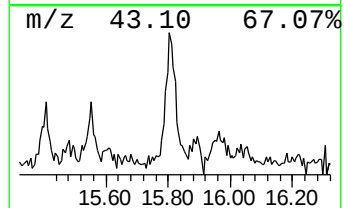
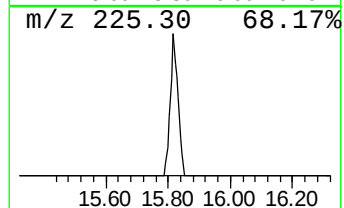
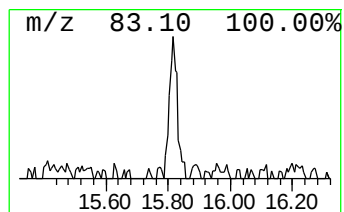
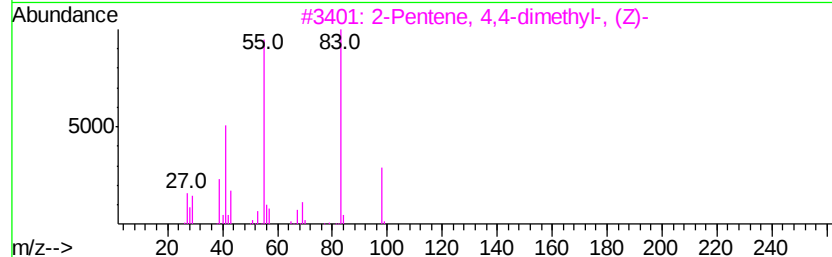
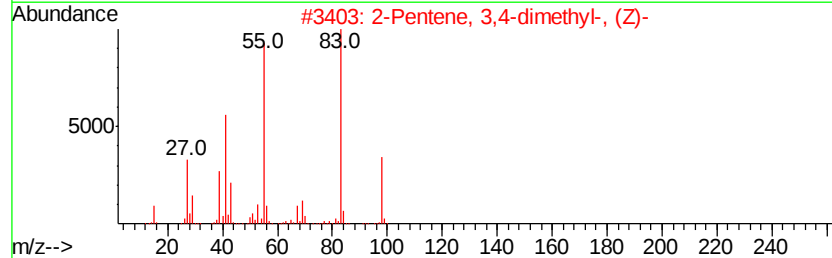
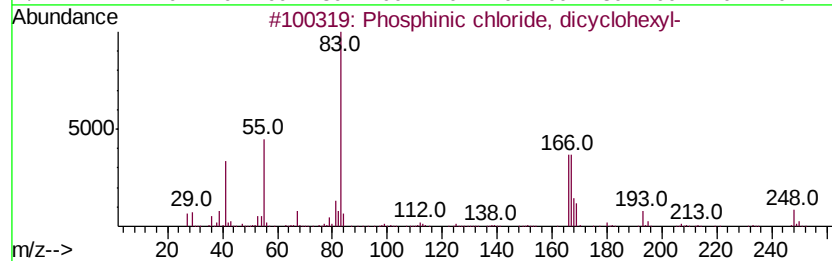
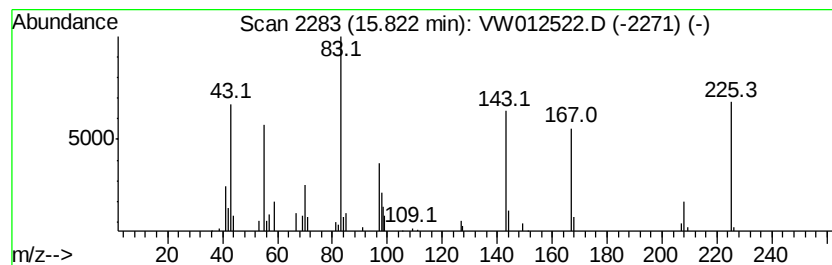
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Peak Number 2 unknown15.82 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.82	5.01 ug/l	22250	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phosphinic chloride, dicyclohexyl-	248	C12H22ClOP	015873-72-0	25
2		2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	22
3		2-Pentene, 4,4-dimethyl-, (Z)-	98	C7H14	000762-63-0	22
4		2-Pentene, 2,3-dimethyl-	98	C7H14	010574-37-5	22
5		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	22



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzeneethanamine...	14.07	6.1	ug/l	27270	4	13.56	221927	50.0
unknown15.82	15.82	5.0	ug/l	22250	4	13.56	221927	50.0