

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW082220\
 Data File : VW016309.D
 Acq On : 21 Aug 2020 16:37
 Operator : SY/VA
 Sample : L3734-04
 Misc : 6.13G/5ML/MSVOA W/SOIL
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 EROC-BF-1C

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\82W081420S.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	3.684	286	292	303	rVB4	4335	10776	1.17%	0.217%
2	4.123	354	364	373	rBV2	4290	11550	1.26%	0.232%
3	4.916	481	494	510	rBV2	52181	173387	18.90%	3.485%
4	7.141	849	859	867	rBV2	4862	14272	1.56%	0.287%
5	7.885	972	981	985	rBV	45062	99311	10.83%	1.996%
6	7.945	985	991	1003	rVB	234366	531732	57.97%	10.688%
7	8.305	1042	1050	1062	rBV	89877	196281	21.40%	3.945%
8	8.842	1129	1138	1149	rBV	347596	683516	74.51%	13.739%
9	10.323	1373	1381	1389	rBV	511717	917298	100.00%	18.438%
10	10.701	1438	1443	1449	rVB3	5970	10384	1.13%	0.209%
11	11.634	1588	1596	1609	rBV	414048	703916	76.74%	14.149%
12	11.841	1624	1630	1638	rBV4	5377	9894	1.08%	0.199%
13	12.183	1680	1686	1691	rVB4	6532	14038	1.53%	0.282%
14	12.621	1750	1758	1767	rVB	301700	498016	54.29%	10.010%
15	12.871	1794	1799	1802	rBV	7275	11728	1.28%	0.236%
16	12.945	1806	1811	1817	rVV2	12053	21653	2.36%	0.435%
17	13.256	1856	1862	1871	rBV	16310	30236	3.30%	0.608%
18	13.560	1905	1912	1921	rBV	330689	538479	58.70%	10.824%
19	13.652	1921	1927	1933	rVV2	12017	18923	2.06%	0.380%
20	13.792	1947	1950	1957	rVB3	6611	10186	1.11%	0.205%
21	13.877	1959	1964	1971	rBV5	9305	17784	1.94%	0.357%
22	14.072	1989	1996	2007	rVB2	57946	111739	12.18%	2.246%
23	14.207	2012	2018	2023	rBV5	8608	16748	1.83%	0.337%
24	14.335	2035	2039	2044	rBV7	4404	10108	1.10%	0.203%
25	14.457	2056	2059	2063	rVB3	6811	9603	1.05%	0.193%
26	14.524	2068	2070	2079	rVB8	7743	17570	1.92%	0.353%
27	14.725	2100	2103	2109	rVB5	6129	10827	1.18%	0.218%
28	14.798	2109	2115	2122	rBV5	12921	34076	3.71%	0.685%
29	14.963	2137	2142	2147	rBV3	7023	13001	1.42%	0.261%
30	15.206	2171	2182	2187	rBV2	23508	50974	5.56%	1.025%
31	15.371	2204	2209	2213	rBV	7850	11868	1.29%	0.239%
32	15.499	2221	2230	2238	rBV	55027	106726	11.63%	2.145%
33	15.865	2286	2290	2297	rVB7	7036	13704	1.49%	0.275%
34	16.182	2332	2342	2347	rBV7	5006	10809	1.18%	0.217%

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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

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Title : SW846 8260

35	16.438	2379	2384	2390	rBV2	7214	14663	1.60%	0.295%
36	16.645	2410	2418	2425	rBV4	8001	19287	2.10%	0.388%

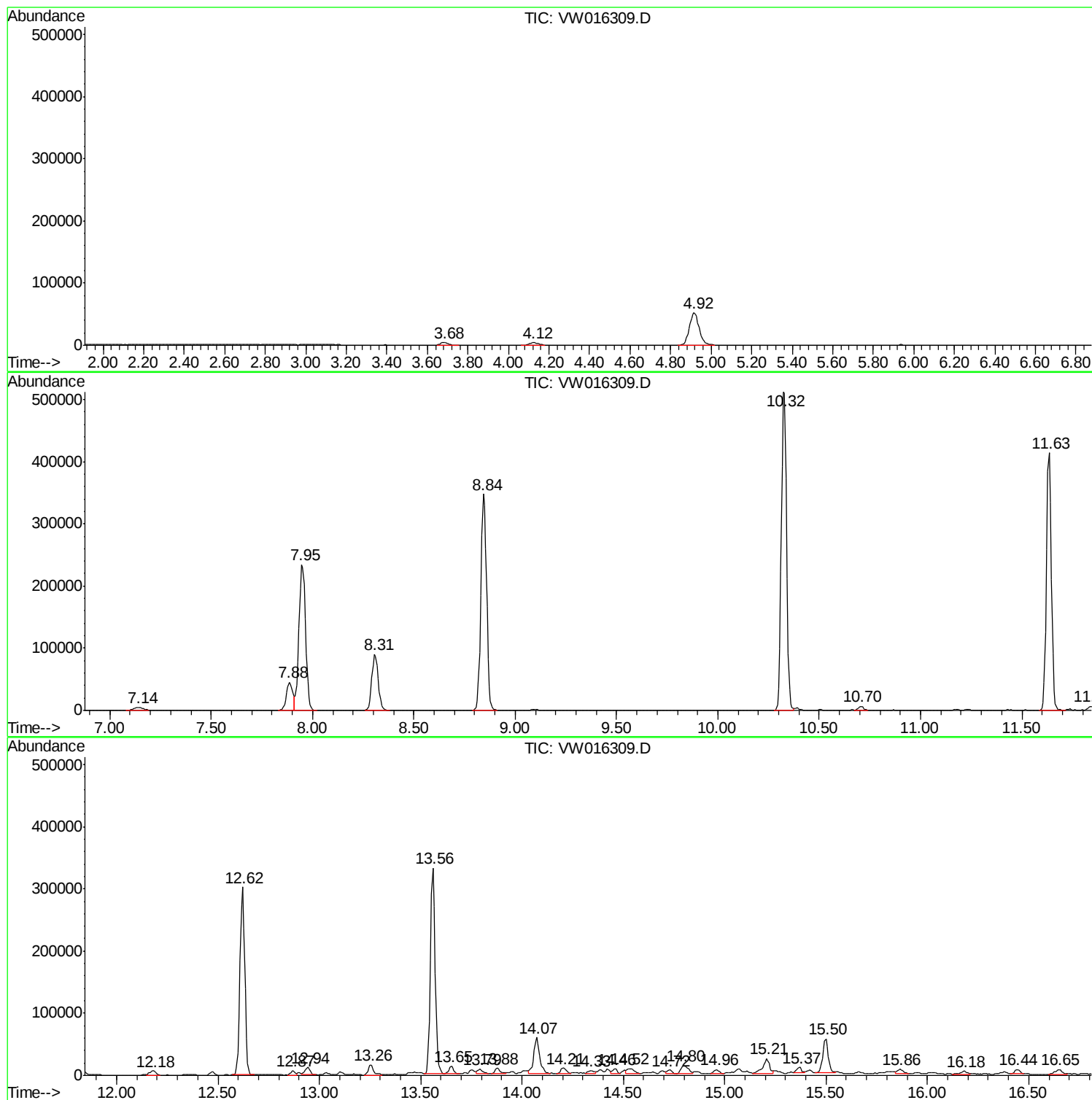
Sum of corrected areas: 4975063

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



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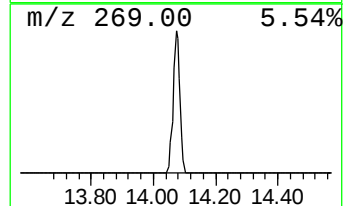
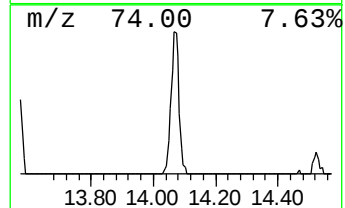
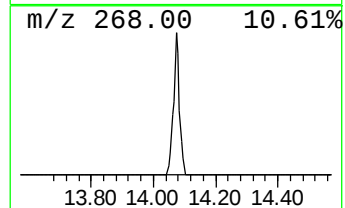
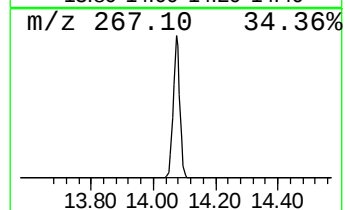
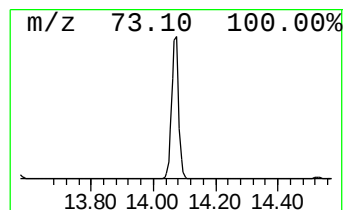
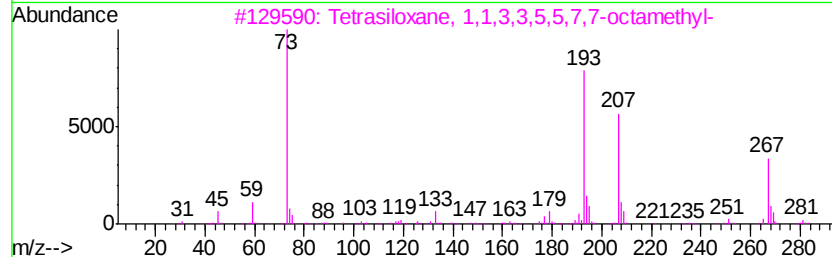
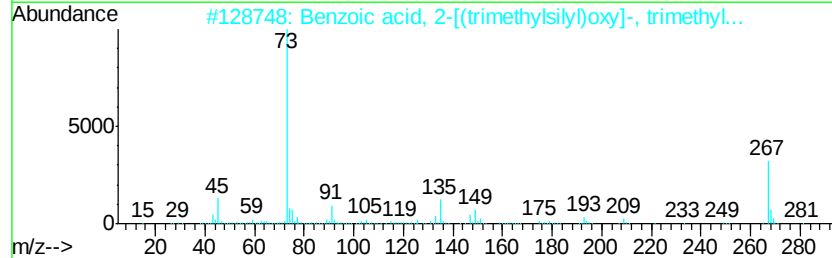
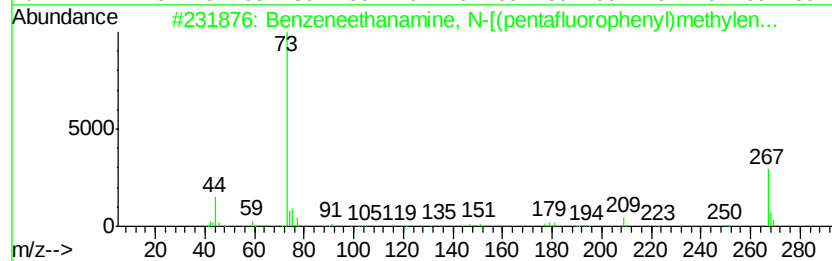
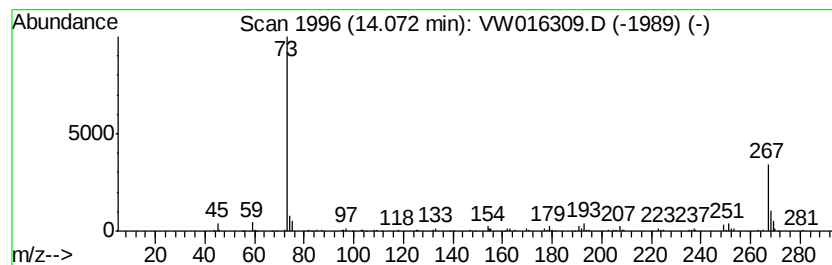
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.07	10.38 ug/l	111739	1,4-Dichlorobenzene-d4	13.56		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzeneethanamine, N-[(pentafluoro...	475	C21H26F5N02Si2	055429-85-1	53
2		Benzoic acid, 2-[(trimethylsilyl)oxy]-, trimethyl...	282	C13H22O3Si2	003789-85-3	39
3		Tetrasiloxane, 1,1,3,3,5,5,7,7-octamethyl-	282	C8H26O3Si4	001000-05-1	39
4		Silane, trimethyl[[(p-(trimethylsilyl)oxy)-, trimethyl...	268	C13H24O2Si2	018401-58-6	9
5		Silane, (2-ethyl-3,3-dimethyl-4-oxo-4-phenylbut-1-en-1-yl)dimethyl-	208	C13H24Si	095798-13-3	9



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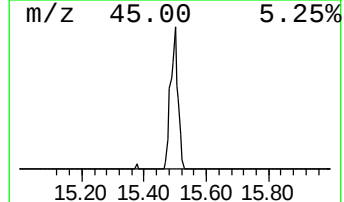
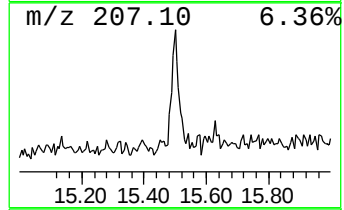
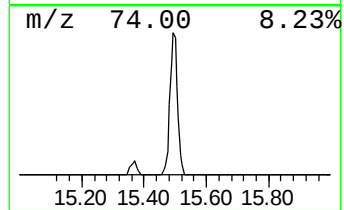
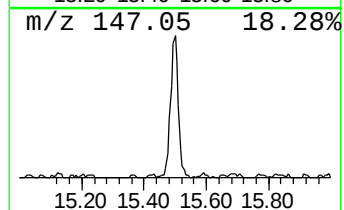
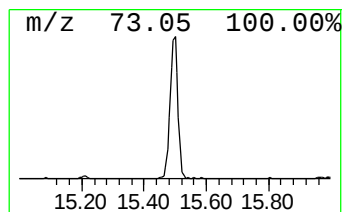
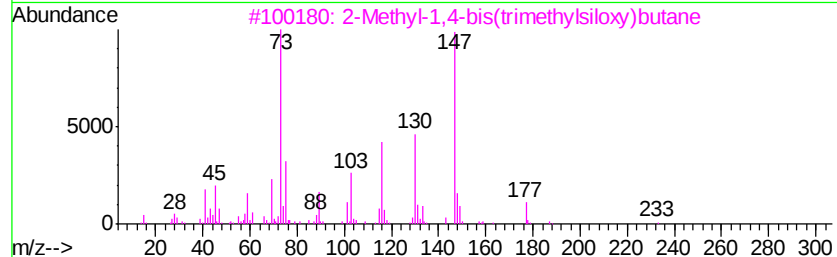
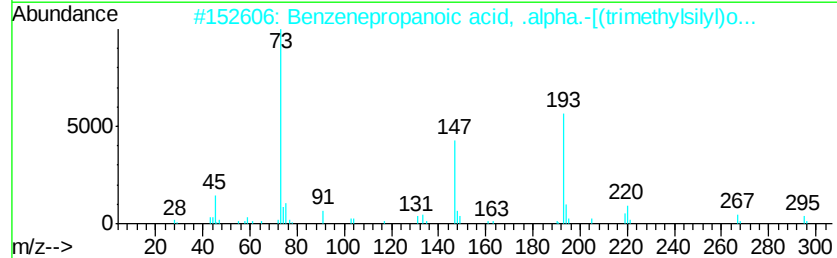
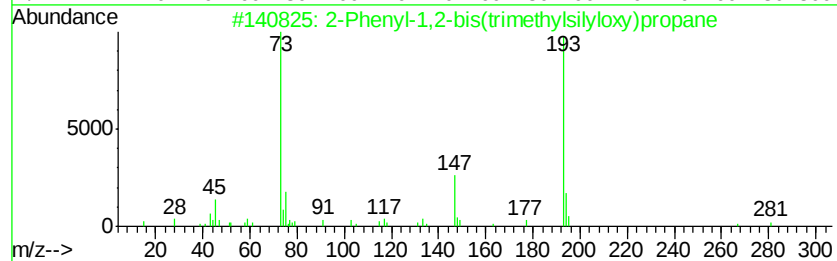
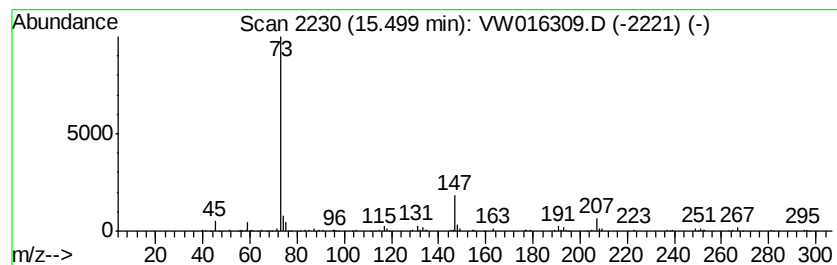
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Peak Number 2 2-Phenyl-1,2-bis(trimethylsiloxy)propane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.50	9.91 ug/l	106726	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Phenyl-1,2-bis(trimethylsilyloxy)propane	296	C15H28O2Si2	294847-15-7	53
2		Benzenepropanoic acid, .alpha.-[trimethylsilyloxy]propane	310	C15H26O3Si2	027750-45-4	42
3		2-Methyl-1,4-bis(trimethylsiloxy)butane	248	C11H28O2Si2	105747-05-5	38
4		Ribo-hexos-3-ulose, 2,4,5,6-tetra-O-(trimethylsilyl)-	524	C20H48N2O6Si4	062108-14-9	36
5		Silane, trimethyl[4-(trimethylsilyloxy)phenyl]-	218	C10H26O2Si2	007140-91-2	36



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzeneethanamine...	14.07	10.4	ug/l	111739	4	13.56	538479	50.0
2-Phenyl-1,2-bis(...	15.50	9.9	ug/l	106726	4	13.56	538479	50.0