

Data Path : W:\HPCHEM1\MSVOA W\DATA\VW050818\
 Data File : VW002298.D
 Acq On : 08 May 2018 16:52
 Operator : JC/SY
 Sample : J2722-02
 Misc : 5.12G/5ML/MSVOA W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 OILY-STONES

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 : W:\HPCHEM1\MSVOA_W\METHOD\82W050818S.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	1.752	19	27	28	rBV	6755	13164	1.77%	0.294%
2	1.807	28	36	55	rVB	285286	744883	100.00%	16.647%
3	2.630	165	171	174	rBV3	4613	10807	1.45%	0.242%
4	4.129	409	417	425	rVB3	2949	8195	1.10%	0.183%
5	4.410	455	463	473	rBV2	7425	23484	3.15%	0.525%
6	4.910	536	545	548	rBV6	3930	10947	1.47%	0.245%
7	7.153	904	913	926	rBV	28043	80790	10.85%	1.806%
8	7.885	1023	1033	1037	rBV	96527	218602	29.35%	4.885%
9	7.946	1037	1043	1053	rVB	205780	460917	61.88%	10.301%
10	8.306	1087	1102	1112	rBV	89642	208053	27.93%	4.650%
11	8.842	1182	1190	1199	rBV	296475	598639	80.37%	13.379%
12	10.330	1427	1434	1442	rBV	273415	480228	64.47%	10.733%
13	10.714	1490	1497	1504	rVB3	26703	45409	6.10%	1.015%
14	11.073	1550	1556	1564	rBV	20148	33998	4.56%	0.760%
15	11.445	1614	1617	1624	rVB4	5206	8918	1.20%	0.199%
16	11.640	1640	1649	1657	rBV	351801	589771	79.18%	13.181%
17	12.628	1801	1811	1820	rBV2	222193	371930	49.93%	8.312%
18	12.945	1859	1863	1869	rVB3	9219	15736	2.11%	0.352%
19	13.421	1935	1941	1947	rBV4	3995	8666	1.16%	0.194%
20	13.567	1959	1965	1974	rVV	232786	383643	51.50%	8.574%
21	13.774	1994	1999	2002	rBV4	5863	9076	1.22%	0.203%
22	14.091	2045	2051	2059	rVB2	25683	42624	5.72%	0.953%
23	14.347	2085	2093	2100	rBV7	9511	20321	2.73%	0.454%
24	15.457	2270	2275	2280	rBV4	7510	13730	1.84%	0.307%
25	15.524	2281	2286	2292	rVB	9606	17094	2.29%	0.382%
26	15.896	2342	2347	2351	rBV8	7817	17357	2.33%	0.388%
27	15.987	2357	2362	2365	rBV6	5020	11220	1.51%	0.251%
28	16.396	2425	2429	2434	rBV6	13395	26312	3.53%	0.588%

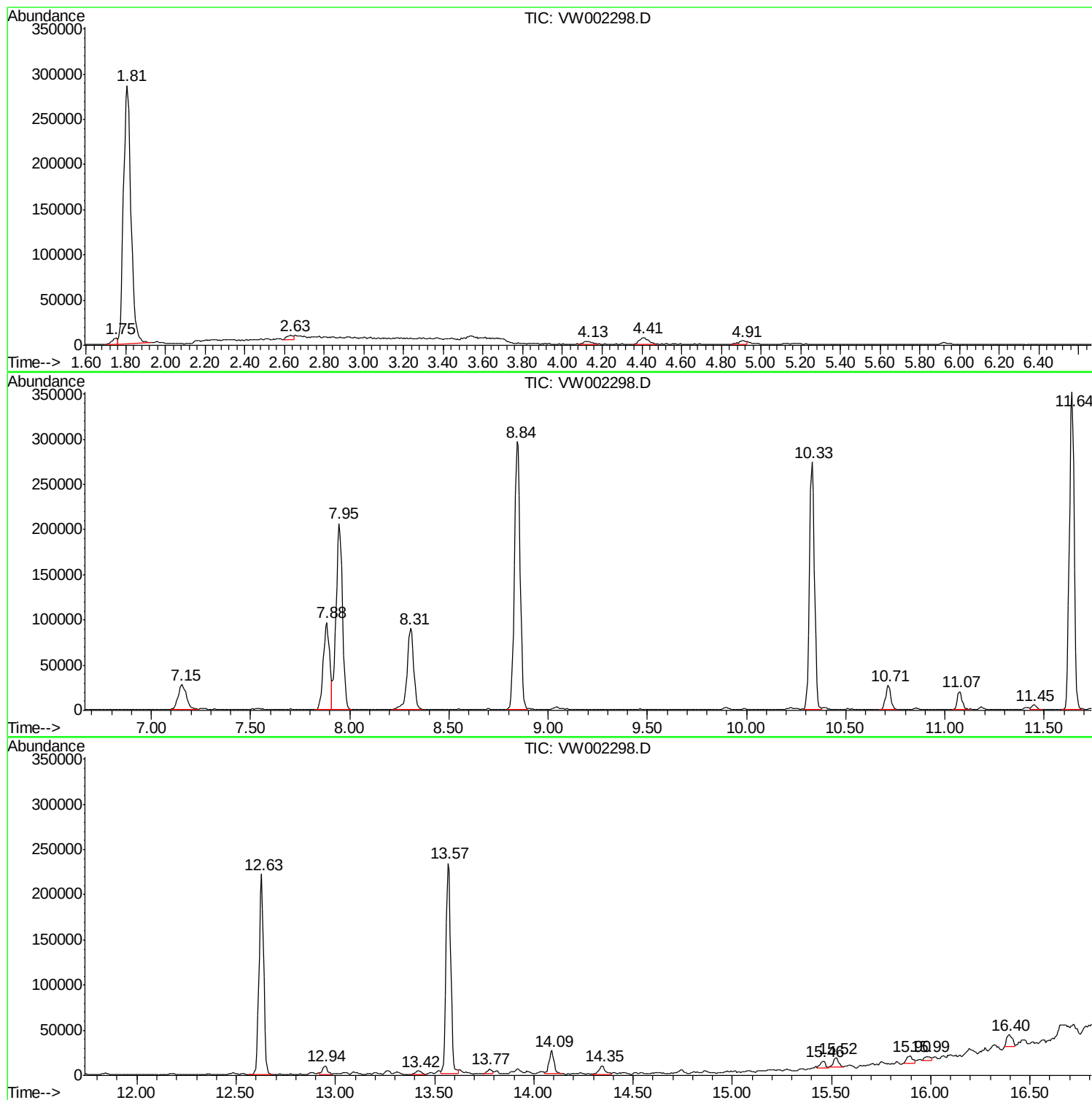
Sum of corrected areas: 4474514

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Quant Method : W:\HPCHEM1\MSVOA_W\METHOD\82W050818S.M
Quant Title : SW846 8260

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



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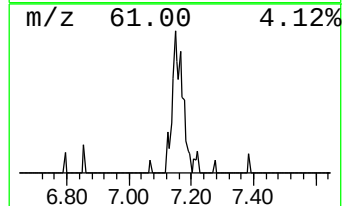
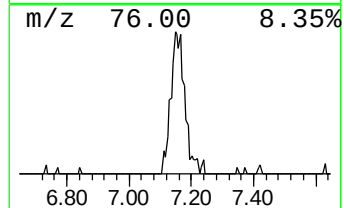
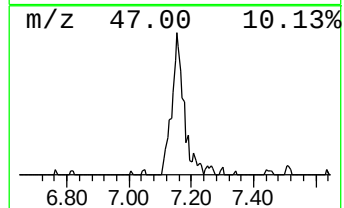
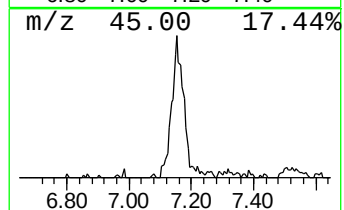
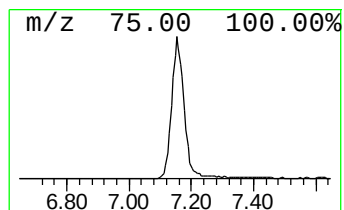
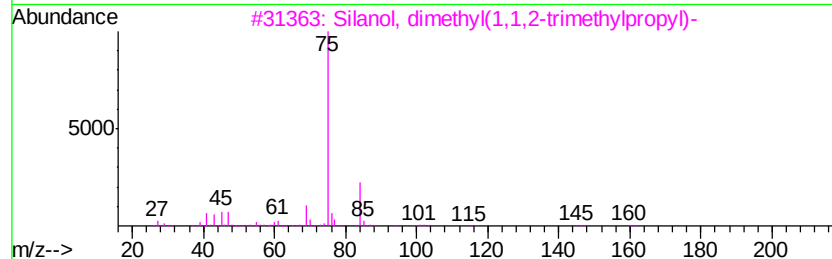
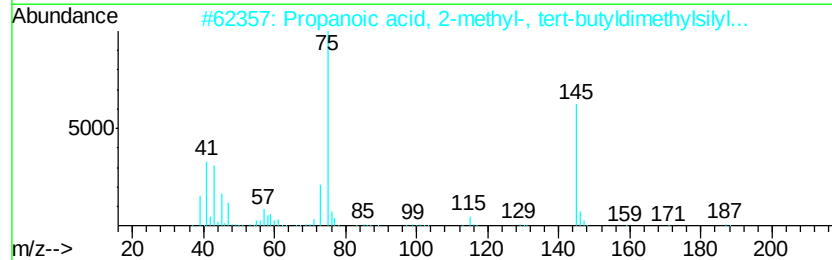
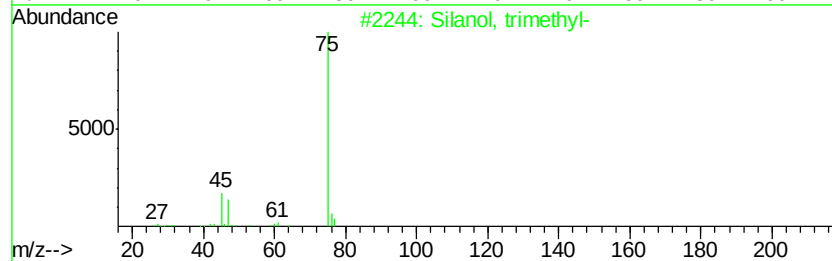
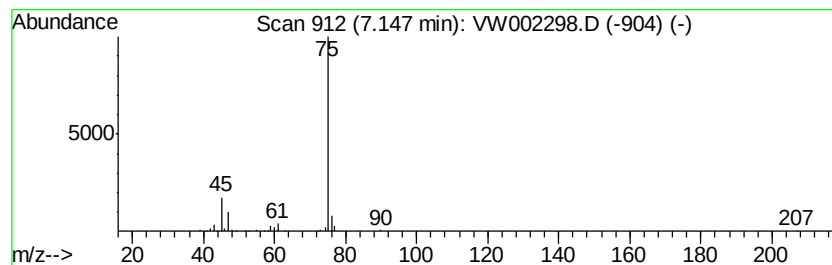
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Peak Number 2 Silanol, trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.15	8.76 ug/l	80790	Pentafluorobenzene	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Silanol, trimethyl-	90	C3H10OSi	001066-40-6	86
2		Propanoic acid, 2-methyl-, tert-...	202	C10H22O2Si	111864-21-2	78
3		Silanol, dimethyl(1,1,2-trimethyl-...	160	C8H20OSi	055644-10-5	78
4		tert-Butyldimethylsilanol	132	C6H16OSi	018173-64-3	64
5		Silanol, trimethyl-, acetate	132	C5H12O2Si	002754-27-0	64



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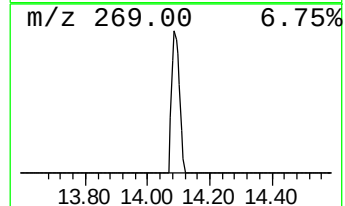
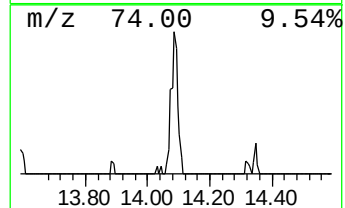
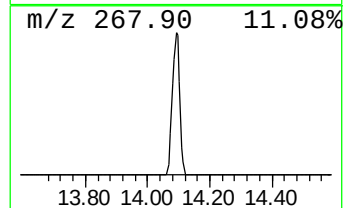
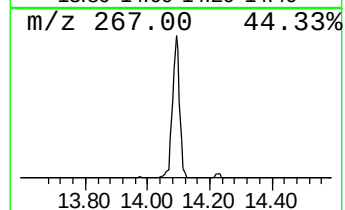
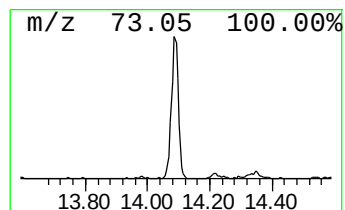
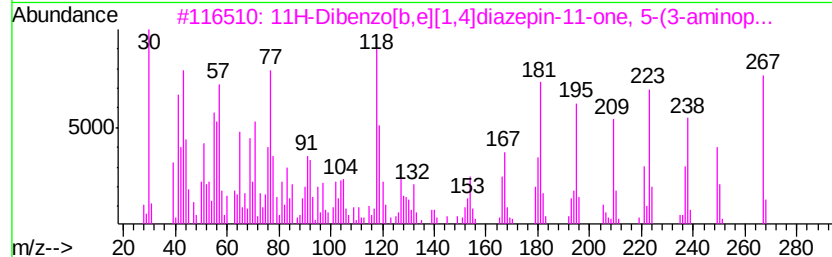
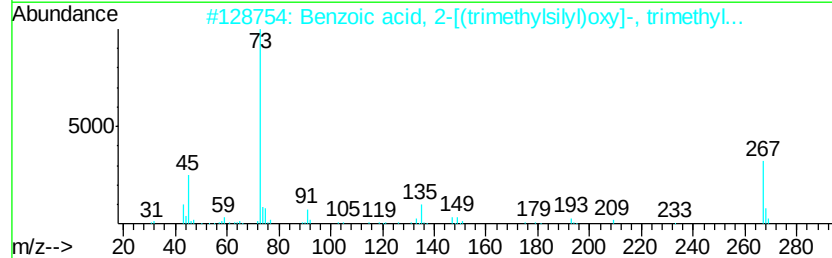
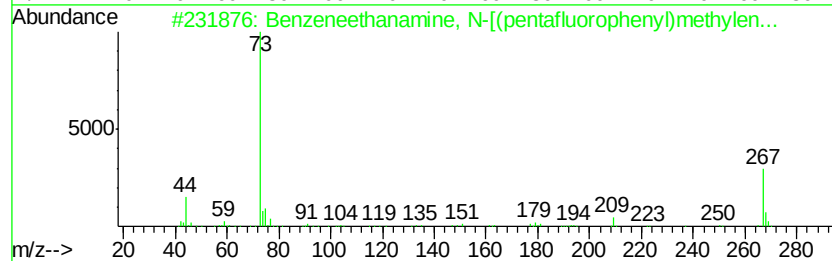
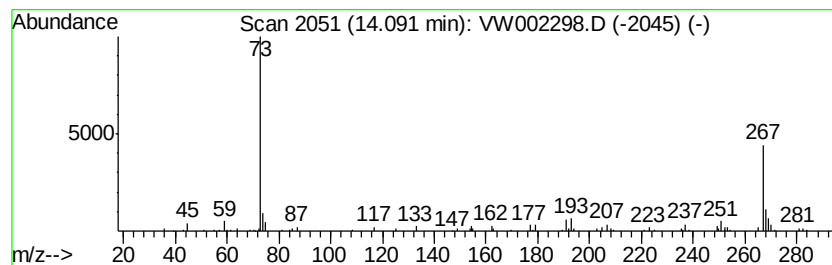
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Peak Number 3 Benzeneethanamine, N-[(pent... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.09	5.56 ug/l	42624	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzeneethanamine, N-[(pentafluoro...]	475	C21H26F5N02Si2	055429-85-1	50
2		Benzoic acid, 2-[(trimethylsilyl)oxy]-, trimethyl...	282	C13H22O3Si2	003789-85-3	50
3		11H-Dibenzo[b,e][1,4]diazepin-11-one, 5-(3-aminopropyl)-	267	C16H17N3O	013450-73-2	46
4		1-(3-Chlorophenyl)piperazine, 4-(3-chlorophenyl)-	268	C13H21ClN2Si	1000373-99-3	10
5		1-Propene-1-thiol	74	C3H6S	000925-89-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Silanol, trimethyl-	7.15	8.8	ug/l	80790	1	7.95	460917	50.0
Benzeneethanamine...	14.09	5.6	ug/l	42624	4	13.57	383643	50.0