

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW040419\
 Data File : VW009692.D
 Acq On : 04 Apr 2019 13:18
 Operator : SY/VA
 Sample : K2197-01
 Misc : 5.40G/5ML/MSVOA W/SOIL
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 CL-02-040219-A

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\82W032919S.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	2.635	113	120	124	rBV7	9743	25594	1.78%	0.255%
2	2.806	143	148	152	rBV	17883	33516	2.34%	0.333%
3	3.660	286	288	298	rVB5	10360	18834	1.31%	0.187%
4	4.117	354	363	372	rBV	35158	92659	6.46%	0.922%
5	4.909	483	493	512	rVB2	24655	106387	7.41%	1.058%
6	7.141	848	859	871	rBV	230222	641208	44.68%	6.379%
7	7.878	968	980	985	rBV	174553	414547	28.89%	4.124%
8	7.945	985	991	1001	rVB	430638	951706	66.32%	9.468%
9	8.299	1031	1049	1060	rBV	184585	439155	30.60%	4.369%
10	8.842	1129	1138	1149	rBV	638265	1264899	88.14%	12.584%
11	9.335	1207	1219	1225	rBV5	11245	30089	2.10%	0.299%
12	9.957	1315	1321	1332	rVB6	6478	17020	1.19%	0.169%
13	10.091	1332	1343	1347	rBV5	8083	21309	1.48%	0.212%
14	10.323	1372	1381	1387	rBV	685959	1215140	84.67%	12.089%
15	10.384	1387	1391	1398	rVB	66069	121770	8.49%	1.211%
16	10.500	1404	1410	1419	rVB4	16625	38622	2.69%	0.384%
17	10.664	1431	1437	1440	rVV3	27021	49118	3.42%	0.489%
18	10.701	1440	1443	1448	rVV	20666	37003	2.58%	0.368%
19	10.762	1448	1453	1458	rVB4	21801	39652	2.76%	0.394%
20	10.939	1477	1482	1488	rVB4	8408	15428	1.08%	0.153%
21	11.061	1492	1502	1507	rVV5	20006	45245	3.15%	0.450%
22	11.109	1507	1510	1513	rVV5	11340	18513	1.29%	0.184%
23	11.177	1513	1521	1525	rVV4	28668	69863	4.87%	0.695%
24	11.231	1525	1530	1535	rVB3	27407	49585	3.46%	0.493%
25	11.335	1543	1547	1552	rBV5	8230	15121	1.05%	0.150%
26	11.439	1552	1564	1571	rVB6	28171	76466	5.33%	0.761%
27	11.512	1571	1576	1582	rBV3	9901	16172	1.13%	0.161%
28	11.634	1585	1596	1605	rBV	837151	1435073	100.00%	14.277%
29	11.725	1605	1611	1615	rBV5	13831	26390	1.84%	0.263%
30	11.841	1621	1630	1634	rBV2	28719	66942	4.66%	0.666%
31	12.164	1673	1683	1692	rBV6	20437	58255	4.06%	0.580%
32	12.621	1751	1758	1767	rVB2	464732	802204	55.90%	7.981%
33	12.938	1807	1810	1816	rVB5	13457	22849	1.59%	0.227%
34	13.103	1829	1837	1844	rBV5	7493	16784	1.17%	0.167%

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Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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35	13.249	1858	1861	1866	rBV2	8667	14385	1.00%	0.143%
36	13.560	1904	1912	1924	rBV	698705	1137362	79.25%	11.315%
37	13.755	1940	1944	1949	rBV5	11372	20288	1.41%	0.202%
38	13.877	1960	1964	1968	rBV6	9567	15871	1.11%	0.158%
39	14.078	1992	1997	2011	rVB2	73690	153683	10.71%	1.529%
40	14.200	2012	2017	2024	rBV5	17014	38679	2.70%	0.385%
41	14.426	2050	2054	2057	rVB6	12700	18064	1.26%	0.180%
42	14.505	2063	2067	2069	rBV4	10370	16782	1.17%	0.167%
43	14.700	2094	2099	2101	rBV3	9920	17584	1.23%	0.175%
44	14.731	2101	2104	2109	rVB4	19530	27223	1.90%	0.271%
45	14.810	2110	2117	2123	rBV5	23151	55003	3.83%	0.547%
46	14.956	2137	2141	2146	rBV8	10118	20157	1.40%	0.201%
47	15.371	2205	2209	2214	rVB	19528	28346	1.98%	0.282%
48	15.505	2225	2231	2239	rVB2	15758	36103	2.52%	0.359%
49	16.450	2381	2386	2393	rVB2	15709	31155	2.17%	0.310%
50	16.651	2410	2419	2428	rBV2	56662	128189	8.93%	1.275%

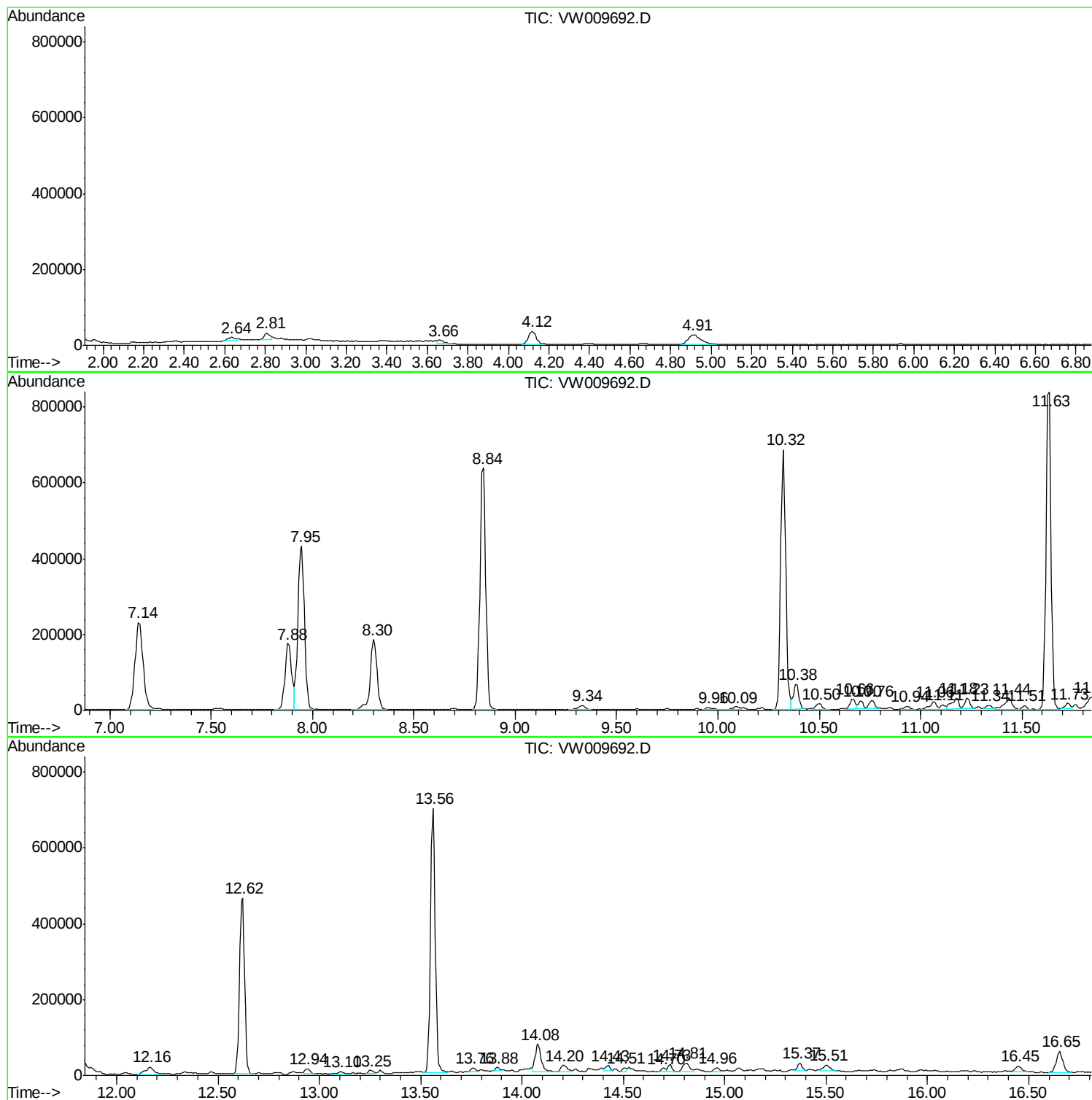
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W032919S.M
Quant Title : SW846 8260

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



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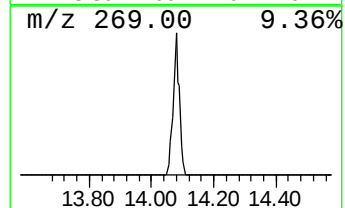
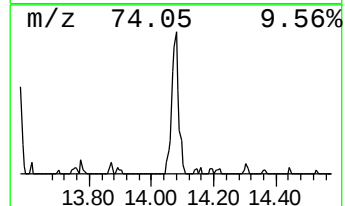
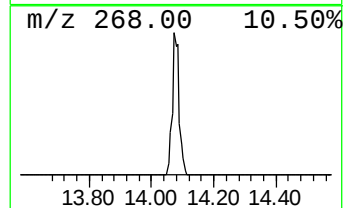
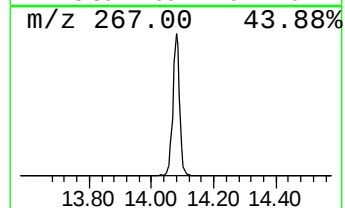
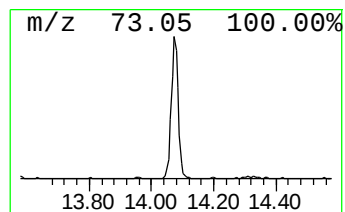
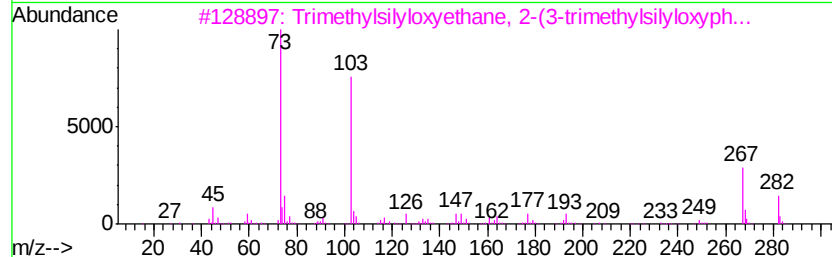
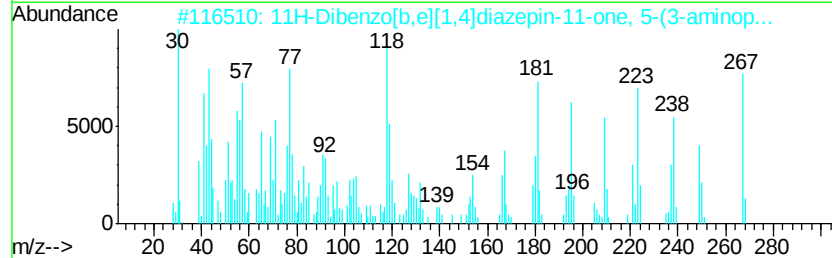
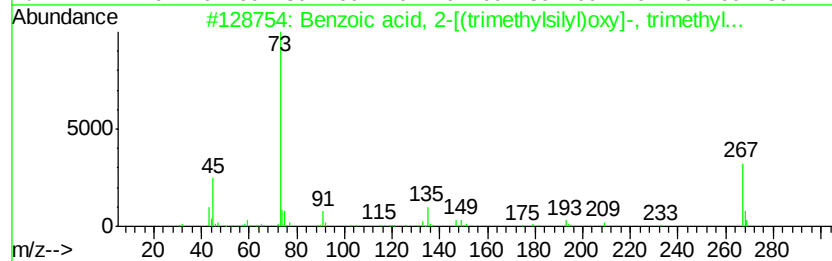
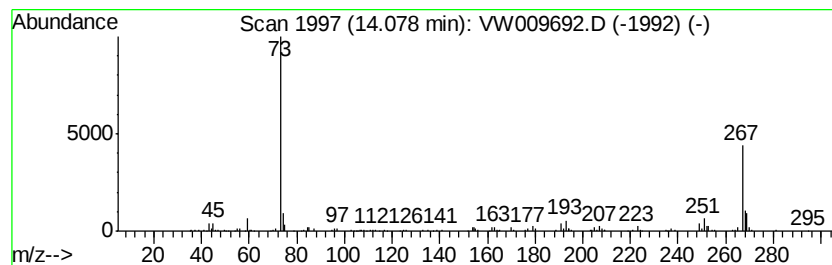
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W032919S.M
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Peak Number 1 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.08	6.76 ug/l	153683	1,4-Dichlorobenzene-d4	13.56		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzoic acid, 2-[(trimethylsilyl)oxy]-, trimethyl...	282	C13H22O3Si2	003789-85-3	45
2		11H-Dibenzo[b,e][1,4]diazepin-11-one, 5-(3-aminopropyl)-	267	C16H17N3O	013450-73-2	38
3		Trimethylsilyloxymethane, 2-(3-trimethylsilyloxyphenyl)-	282	C14H26O2Si2	1000365-49-0	25
4		Benzeneethanamine, N-[(pentafluorophenyl)methyl]-	475	C21H26F5N02Si2	055429-85-1	25
5		Silane, (1-cyclopenten-1-yl)methyl-	154	C9H18Si	075311-60-3	10



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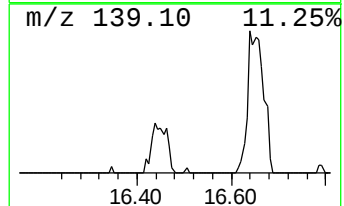
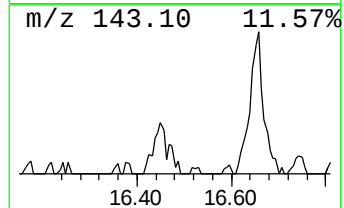
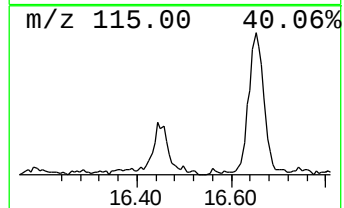
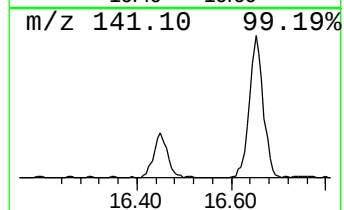
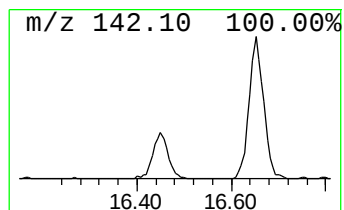
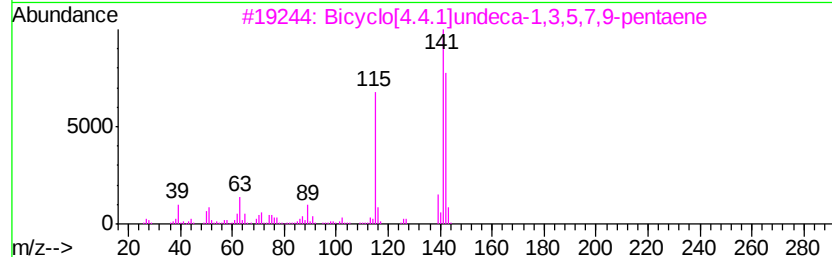
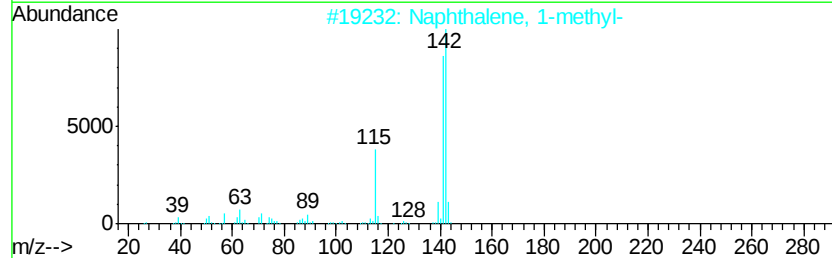
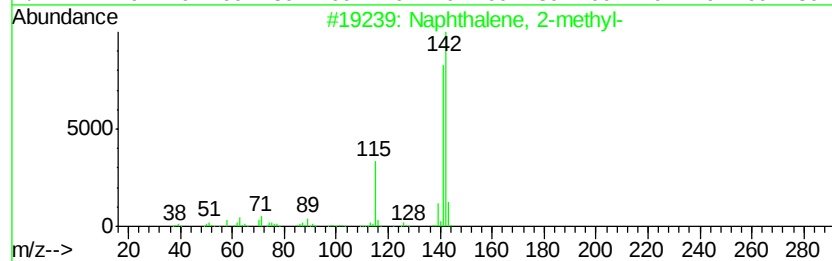
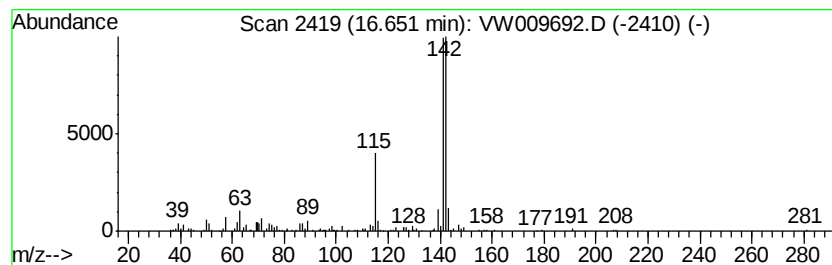
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Peak Number 2 Naphthalene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.65	5.64 ug/l	128189	1,4-Dichlorobenzene-d4	13.56		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
3		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	94
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	93



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	14.08	6.8	ug/l	153683	4	13.56	1137360	50.0
Naphthalene, 1-me...	16.65	5.6	ug/l	128189	4	13.56	1137360	50.0