

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX042619\
 Data File : VX009184.D
 Acq On : 26 Apr 2019 13:03
 Operator : JC/SP
 Sample : K2581-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-03

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_X\METHOD\82X041519W.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	1.343	26	31	33	rBV2	11107	19689	1.72%	0.297%
2	2.611	227	239	244	rBV	9939	20072	1.75%	0.303%
3	5.500	700	713	726	rBV	118946	346761	30.22%	5.227%
4	5.659	728	739	757	rVV	218353	619100	53.96%	9.331%
5	6.061	795	805	814	rBV	149004	394746	34.41%	5.950%
6	6.147	815	819	826	rVB2	8444	20302	1.77%	0.306%
7	6.860	925	936	954	rBV	352764	831098	72.44%	12.527%
8	8.713	1233	1240	1248	rBV	743157	1147286	100.00%	17.292%
9	8.780	1248	1251	1259	rVB4	6547	12234	1.07%	0.184%
10	10.109	1463	1469	1482	rBV	678381	928744	80.95%	13.998%
11	10.359	1502	1510	1515	rVB	9602	17473	1.52%	0.263%
12	10.701	1561	1566	1575	rVB3	9005	13789	1.20%	0.208%
13	11.134	1632	1637	1648	rBV	530520	700713	61.08%	10.561%
14	11.804	1743	1747	1753	rVB	13560	18661	1.63%	0.281%
15	12.024	1779	1783	1787	rBV2	13885	18175	1.58%	0.274%
16	12.079	1787	1792	1798	rVB	633317	799151	69.66%	12.045%
17	12.469	1852	1856	1862	rBV4	8001	15210	1.33%	0.229%
18	12.975	1934	1939	1943	rBV2	29930	45073	3.93%	0.679%
19	13.347	1996	2000	2006	rVB4	12360	23373	2.04%	0.352%
20	13.475	2011	2021	2028	rBV4	12372	30386	2.65%	0.458%
21	13.834	2072	2080	2087	rBV2	18933	53200	4.64%	0.802%
22	13.907	2088	2092	2095	rBV4	10817	16481	1.44%	0.248%
23	14.285	2151	2154	2159	rVB7	10909	18240	1.59%	0.275%
24	14.371	2165	2168	2172	rBV6	7499	13296	1.16%	0.200%
25	14.542	2192	2196	2200	rBV3	15823	24418	2.13%	0.368%
26	14.670	2214	2217	2224	rBV3	27615	55175	4.81%	0.832%
27	14.731	2224	2227	2231	rVV	16659	19255	1.68%	0.290%
28	14.810	2238	2240	2242	rBV2	12372	12619	1.10%	0.190%
29	14.846	2243	2246	2249	rVB4	19393	24037	2.10%	0.362%
30	14.901	2253	2255	2260	rVB5	21548	26149	2.28%	0.394%
31	15.249	2308	2312	2314	rBV3	16565	26060	2.27%	0.393%
32	15.273	2314	2316	2320	rVB2	30787	36350	3.17%	0.548%
33	15.486	2347	2351	2357	rVB2	60044	85627	7.46%	1.291%
34	15.554	2358	2362	2365	rBV3	25555	47314	4.12%	0.713%

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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 >
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35	15.688	2381	2384	2387	rVB	24641	29739	2.59%	0.448%
36	15.816	2402	2405	2411	rVB3	77477	124638	10.86%	1.879%

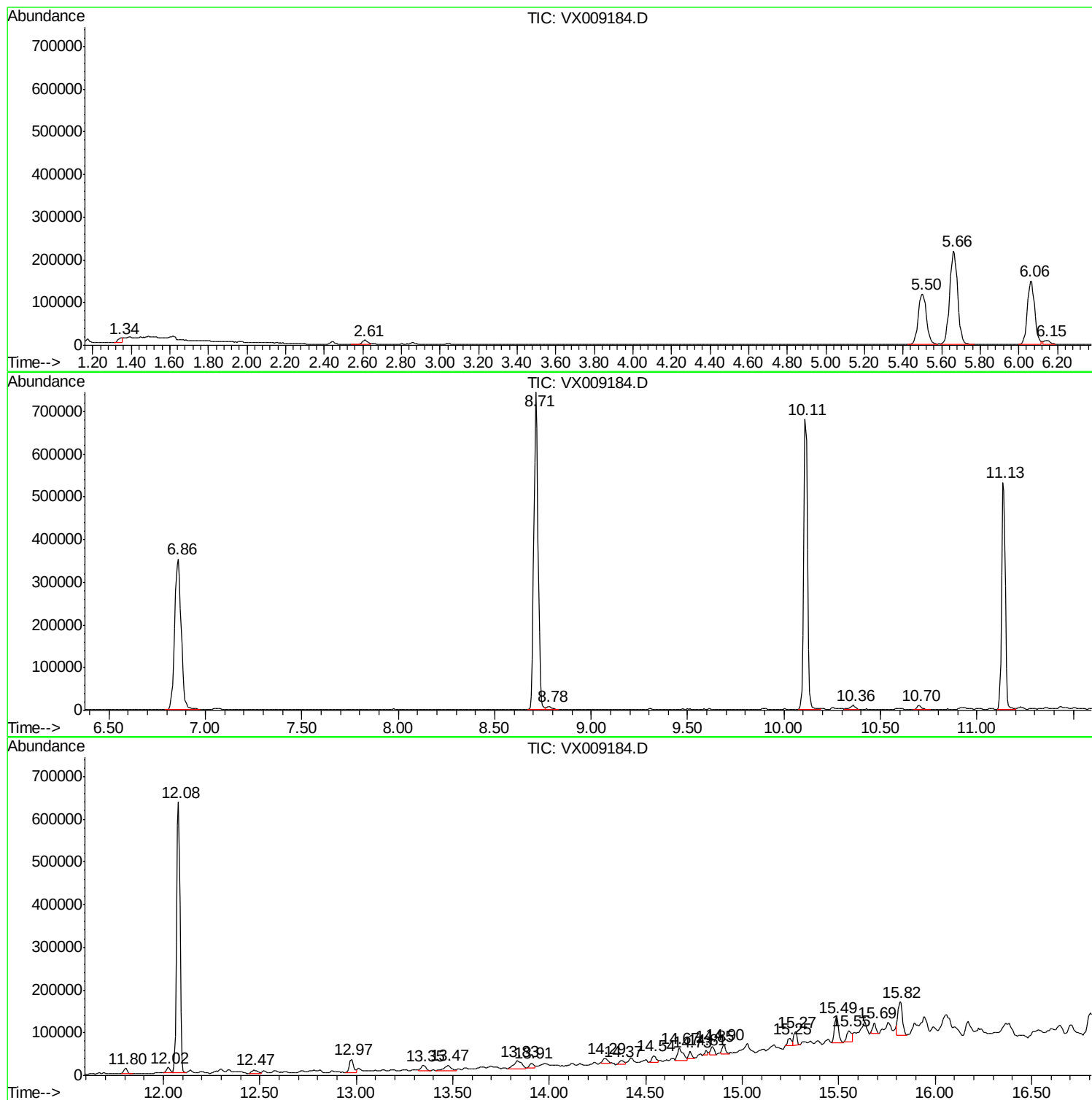
Sum of corrected areas: 6634634

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X041519W.M
Quant Title : SW846 8260

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



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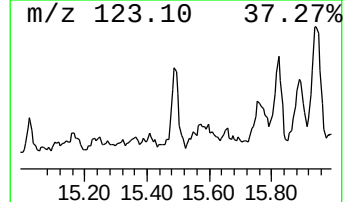
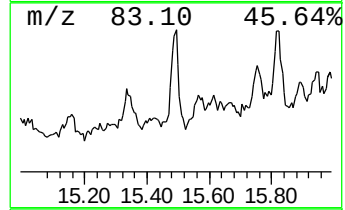
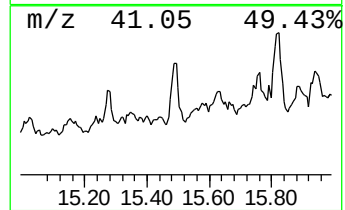
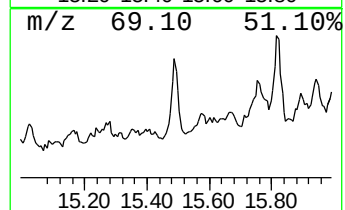
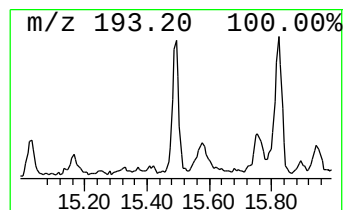
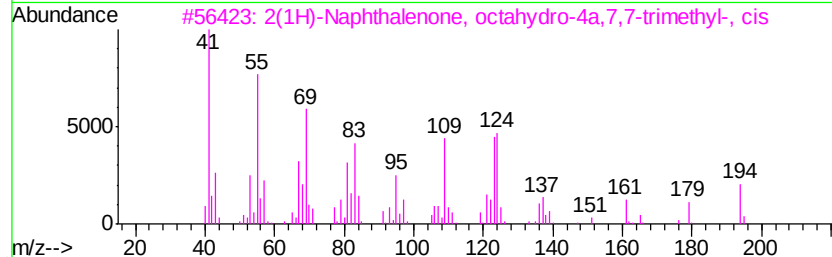
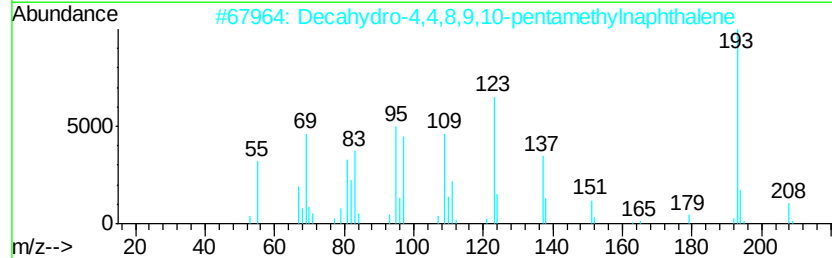
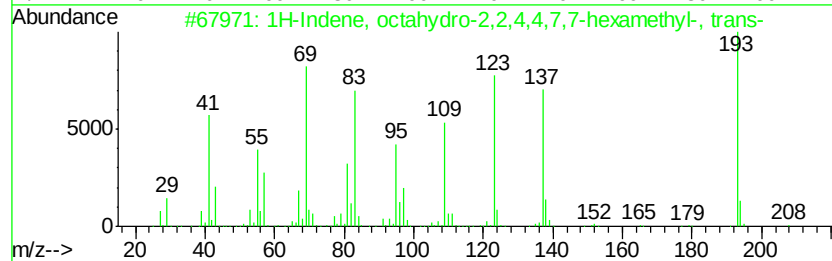
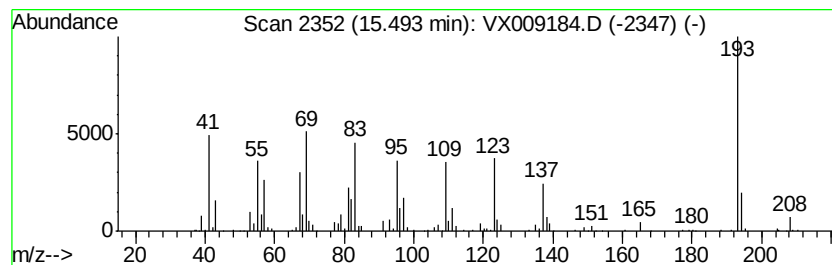
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ClientSampled :
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Peak Number 1 1H-Indene, octahydro-2,2,4,... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.49	5.36 ug/l	85627	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, octahydro-2,2,4,4,7,7...	208 C15H28	054832-83-6	59
2	Decahydro-4,4,8,9,10-pentamethyl...	208 C15H28	080655-44-3	52
3	2(1H)-Naphthalenone, octahydro-4...	194 C13H22O	007056-56-6	37
4	1-(p-Fluorophenyl)-4-piperidone	193 C11H12FNO	1000238-56-7	35
5	Borinic acid, diethyl-, 3,3,5-tr...	208 C13H25BO	057387-76-5	27



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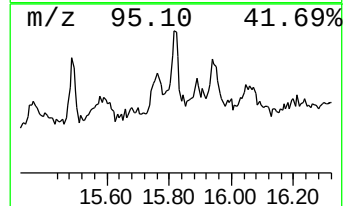
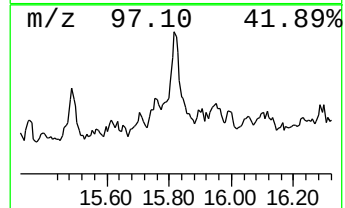
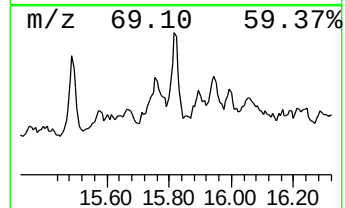
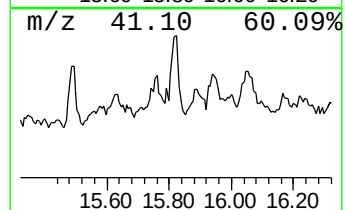
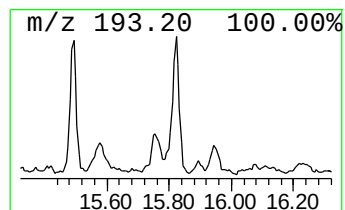
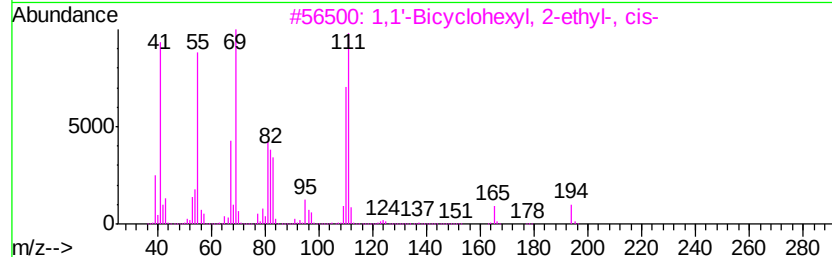
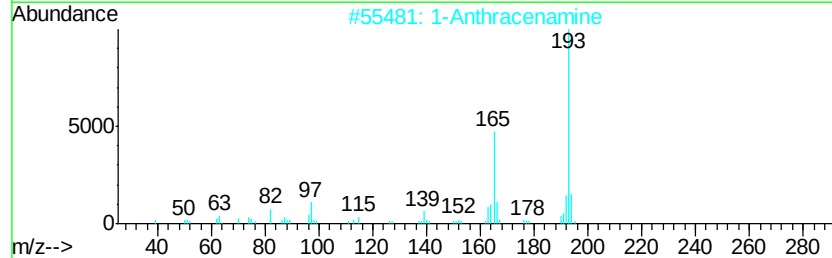
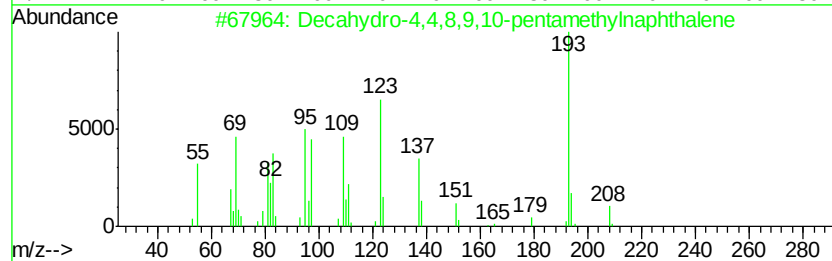
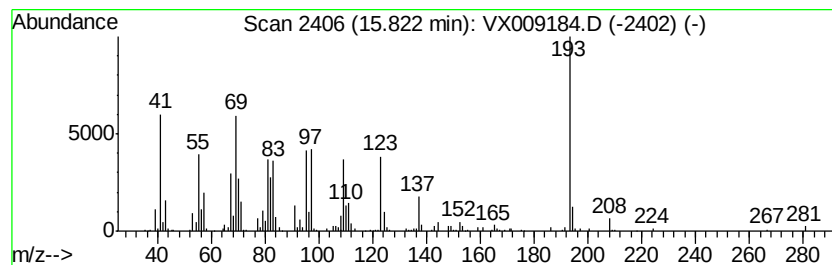
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Decahydro-4,4,8,9,10-pentam... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.82	7.80 ug/l	124638	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Decahydro-4,4,8,9,10-pentamethyl...	208 C15H28	080655-44-3 92
2		1-Anthracenamine	193 C14H11N	000610-49-1 22
3		1,1'-Bicyclohexyl, 2-ethyl-, cis-	194 C14H26	050991-12-3 18
4		1,7-Octadien-3-one, 2-methyl-6-m...	150 C10H14O	041702-60-7 15
5		Bicyclo[2.2.1]heptane, 2,2,3-tri...	138 C10H18	020536-40-7 15



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1H-Indene, octahy...	15.49	5.4	ug/l	85627	4	12.08	799151	50.0
Decahydro-4,4,8,9...	15.82	7.8	ug/l	124638	4	12.08	799151	50.0