

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW110518\
 Data File : VW006672.D
 Acq On : 06 Nov 2018 08:30
 Operator : SY/AP
 Sample : J5812-05
 Misc : 5.83G/5ML/MSVOA W/SOIL
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 TP1-B

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\82W110218S.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	4.135	354	366	381	rBV2	12291	35864	3.25%	0.445%
2	4.909	481	493	506	rBV3	10631	36452	3.30%	0.452%
3	7.165	850	863	865	rBV	5519	15853	1.44%	0.197%
4	7.891	972	982	986	rBV	172226	400933	36.34%	4.975%
5	7.952	986	992	1003	rVB	326341	722035	65.45%	8.959%
6	8.311	1042	1051	1064	rVB	145824	321945	29.18%	3.995%
7	8.848	1130	1139	1149	rBV	456333	900129	81.59%	11.169%
8	10.329	1373	1382	1388	rBV	625486	1103224	100.00%	13.689%
9	10.390	1388	1392	1402	rVB	33092	56020	5.08%	0.695%
10	11.634	1588	1596	1607	rBV	619800	1028065	93.19%	12.756%
11	12.530	1736	1743	1747	rBV6	7870	13805	1.25%	0.171%
12	12.621	1752	1758	1768	rVB	445823	725442	65.76%	9.001%
13	13.566	1905	1913	1921	rVB	575494	934333	84.69%	11.593%
14	13.761	1940	1945	1956	rVB2	14670	27933	2.53%	0.347%
15	13.883	1960	1965	1969	rBV3	12484	22514	2.04%	0.279%
16	14.085	1993	1998	2006	rVB4	10215	24879	2.26%	0.309%
17	14.213	2013	2019	2024	rBV2	13134	23317	2.11%	0.289%
18	14.395	2043	2049	2052	rBV4	14087	26480	2.40%	0.329%
19	14.560	2072	2076	2081	rVB5	10087	16399	1.49%	0.203%
20	14.700	2094	2099	2102	rBV2	17494	28688	2.60%	0.356%
21	14.731	2102	2104	2110	rVB4	9897	15763	1.43%	0.196%
22	14.816	2110	2118	2123	rBV3	45882	94836	8.60%	1.177%
23	14.865	2123	2126	2131	rVB4	7998	13602	1.23%	0.169%
24	15.072	2157	2160	2166	rVB5	9946	17335	1.57%	0.215%
25	15.279	2188	2194	2200	rBV8	5800	14247	1.29%	0.177%
26	15.377	2205	2210	2216	rVB	56389	98873	8.96%	1.227%
27	15.481	2223	2227	2232	rBV5	13438	26101	2.37%	0.324%
28	15.664	2251	2257	2265	rBV4	14824	39203	3.55%	0.486%
29	15.743	2266	2270	2275	rBV6	6768	11780	1.07%	0.146%
30	15.834	2279	2285	2289	rVV6	7149	16370	1.48%	0.203%
31	15.962	2303	2306	2311	rVB7	7456	12000	1.09%	0.149%
32	16.182	2335	2342	2347	rBV8	12066	29583	2.68%	0.367%
33	16.389	2369	2376	2380	rBV	28171	62493	5.66%	0.775%
34	16.456	2380	2387	2402	rVB	130381	304871	27.63%	3.783%

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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

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35	16.657	2412	2420	2431	rVB	381542	838119	75.97%	10.399%
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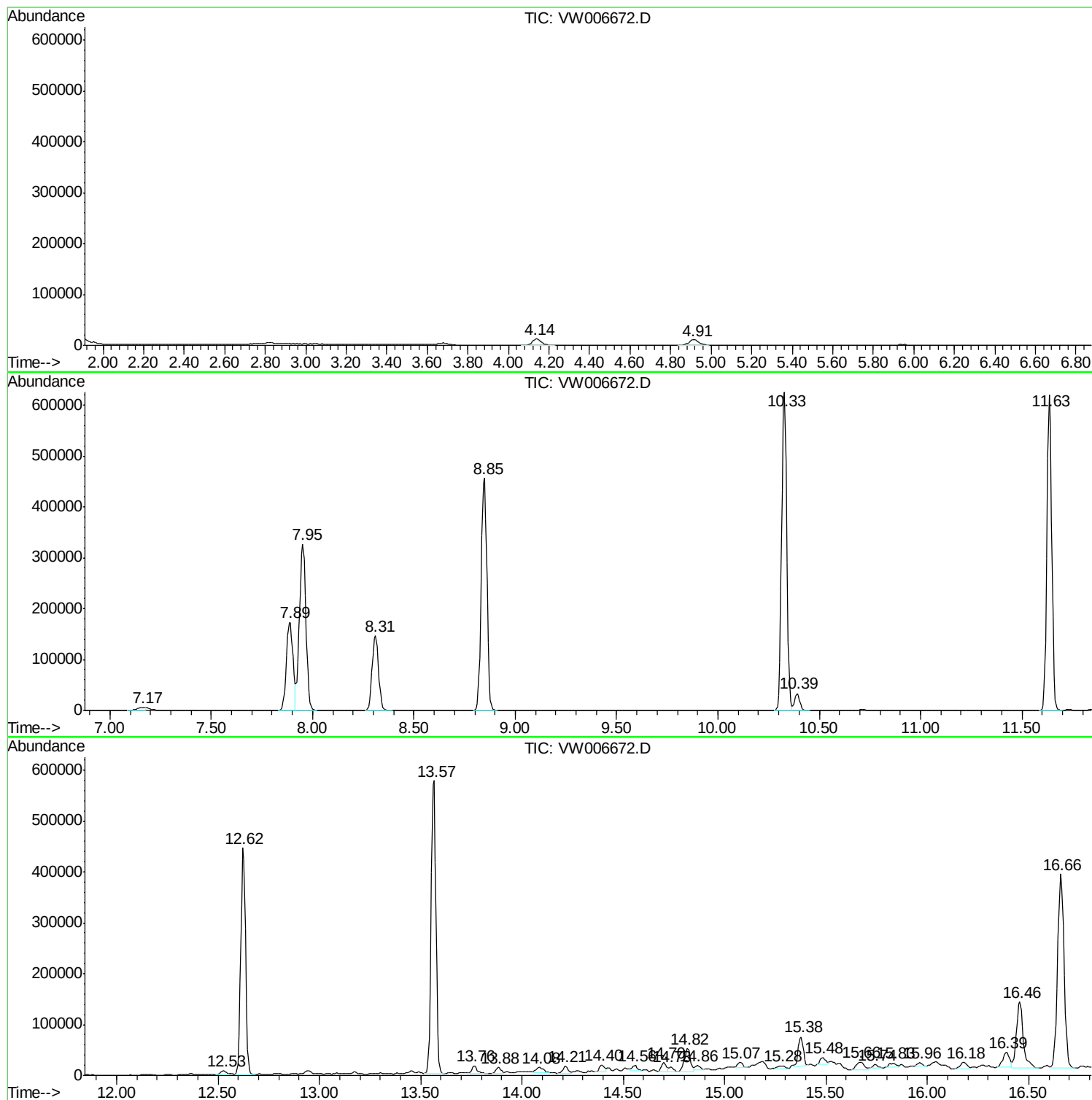
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W110218S.M
Quant Title : SW846 8260

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



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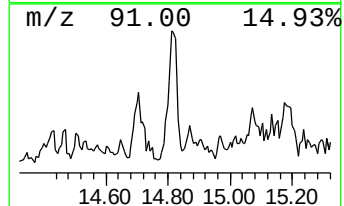
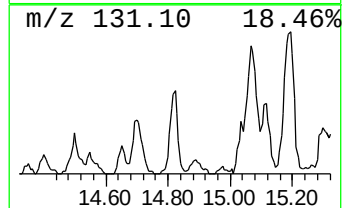
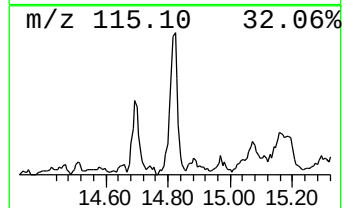
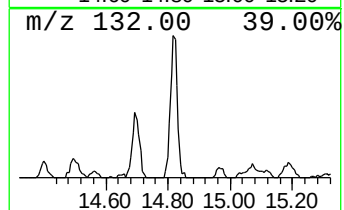
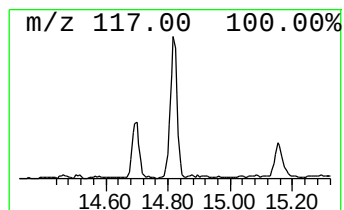
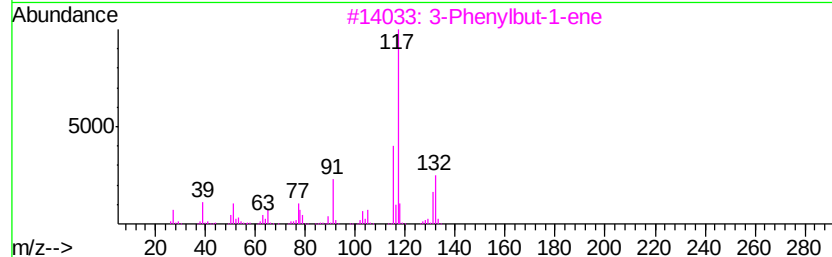
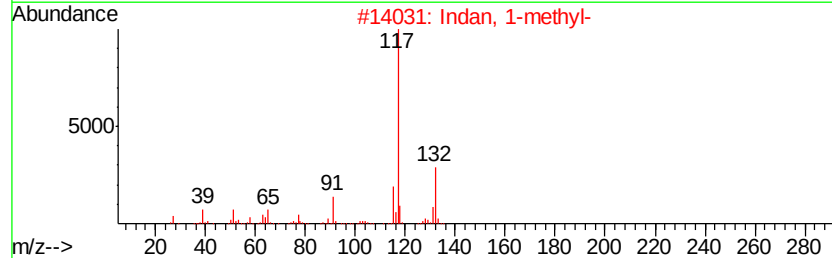
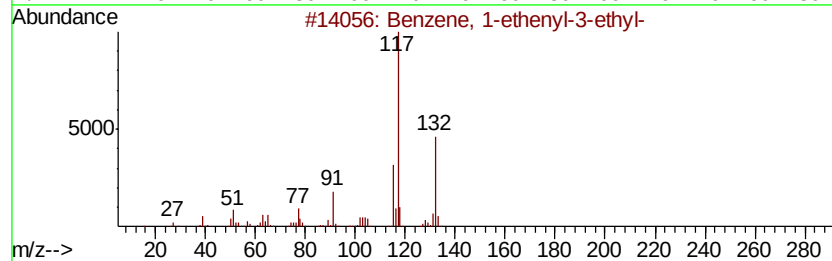
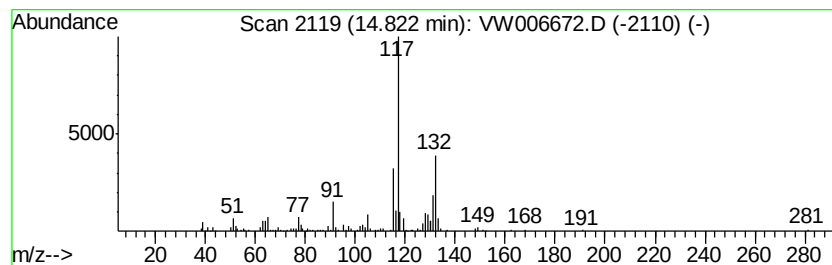
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Peak Number 1 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	5.08 ug/l	94836	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
2		Indan, 1-methyl-	132	C10H12	000767-58-8	87
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
4		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	87



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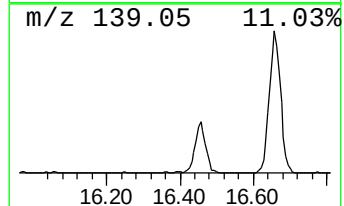
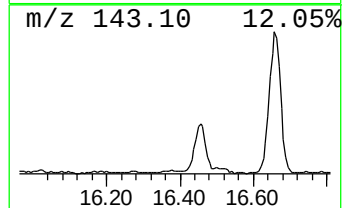
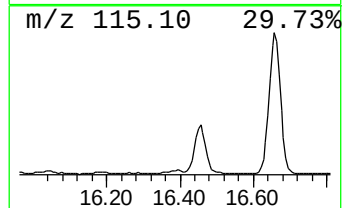
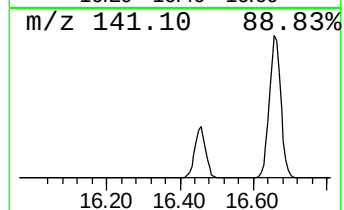
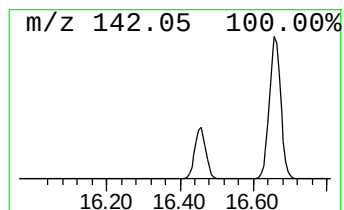
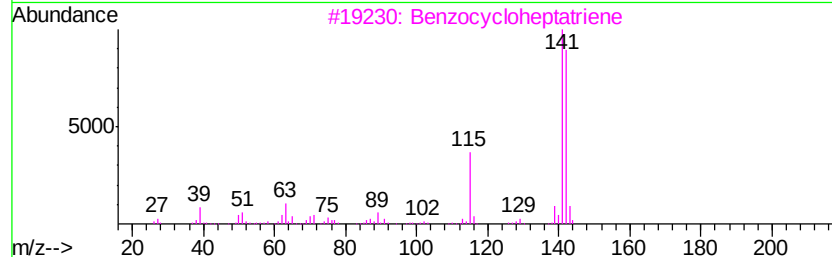
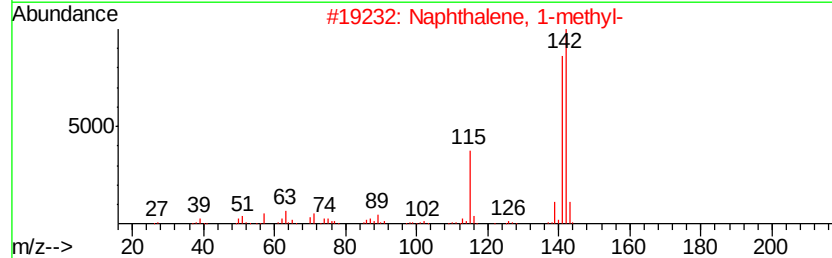
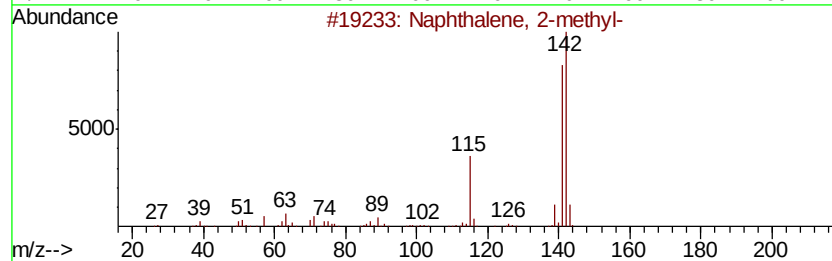
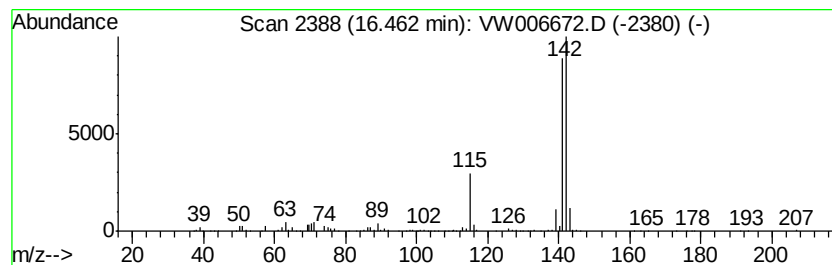
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.46	16.31 ug/l	304871	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
3		Benzocycloheptatriene	142	C11H10	000264-09-5	91
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-etheny...	14.82	5.1	ug/l	94836	4	13.57	934333	50.0
Naphthalene, 2-me...	16.46	16.3	ug/l	304871	4	13.57	934333	50.0