

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW112618\
 Data File : VW007077.D
 Acq On : 27 Nov 2018 06:01
 Operator : VAY/AP
 Sample : J6065-01
 Misc : 5.83G/5ML/MSVOA_W/SOIL
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 USTS-1-2-3

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\82W112118S.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	4.915	480	494	514	rBV4	35775	143668	11.00%	0.991%
2	5.434	567	579	591	rBV2	9636	32257	2.47%	0.223%
3	6.982	825	833	842	rVV3	7772	21824	1.67%	0.151%
4	7.884	969	981	986	rBV	74790	183155	14.02%	1.264%
5	7.951	986	992	1000	rVV	173332	394374	30.20%	2.721%
6	8.031	1000	1005	1014	rVB4	9222	23144	1.77%	0.160%
7	8.159	1017	1026	1032	rBV2	16732	36493	2.79%	0.252%
8	8.305	1042	1050	1062	rVV	81566	180773	13.84%	1.247%
9	8.433	1062	1071	1078	rVV3	15245	39155	3.00%	0.270%
10	8.512	1078	1084	1089	rVV4	15480	35043	2.68%	0.242%
11	8.573	1089	1094	1104	rVV	13765	31672	2.43%	0.219%
12	8.841	1130	1138	1150	rBV	228909	473583	36.26%	3.268%
13	9.073	1168	1176	1180	rBV4	7743	17580	1.35%	0.121%
14	9.122	1180	1184	1194	rVB4	5866	13326	1.02%	0.092%
15	9.335	1205	1219	1225	rBV3	39984	114144	8.74%	0.788%
16	9.390	1225	1228	1237	rVB4	7362	13532	1.04%	0.093%
17	9.610	1256	1264	1271	rVB2	13553	30793	2.36%	0.212%
18	9.750	1281	1287	1297	rVB2	14670	27061	2.07%	0.187%
19	9.854	1298	1304	1308	rBV3	13078	22139	1.70%	0.153%
20	9.902	1308	1312	1316	rVV2	11070	20755	1.59%	0.143%
21	9.963	1316	1322	1324	rVV6	7651	17114	1.31%	0.118%
22	9.994	1324	1327	1335	rVB5	10998	23683	1.81%	0.163%
23	10.091	1335	1343	1356	rBV5	13556	44955	3.44%	0.310%
24	10.250	1364	1369	1374	rVB3	17977	31907	2.44%	0.220%
25	10.323	1374	1381	1388	rBV	285390	537026	41.12%	3.706%
26	10.390	1388	1392	1398	rVV	34792	64418	4.93%	0.445%
27	10.512	1404	1412	1418	rVB7	22184	65968	5.05%	0.455%
28	10.609	1423	1428	1432	rVB6	7953	13885	1.06%	0.096%
29	10.664	1432	1437	1441	rBV	26616	45691	3.50%	0.315%
30	10.750	1441	1451	1462	rVB5	43721	140854	10.79%	0.972%
31	10.847	1462	1467	1472	rVB4	10571	21012	1.61%	0.145%
32	10.939	1472	1482	1487	rBV6	10121	24518	1.88%	0.169%
33	11.036	1487	1498	1505	rBV4	17381	60376	4.62%	0.417%
34	11.109	1505	1510	1512	rBV3	9696	15092	1.16%	0.104%

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35	11.146	1512	1516	1518	rVV4	13004	21536	1.65%	0.149%
36	11.183	1518	1522	1525	rVV	18291	30345	2.32%	0.209%
37	11.231	1525	1530	1535	rVB	32042	56092	4.30%	0.387%
38	11.292	1535	1540	1543	rBV4	11704	17602	1.35%	0.121%
39	11.341	1543	1548	1552	rVB3	15019	23338	1.79%	0.161%
40	11.439	1560	1564	1572	rVB3	24293	49097	3.76%	0.339%
41	11.512	1572	1576	1580	rBV2	10463	21005	1.61%	0.145%
42	11.634	1590	1596	1603	rBV	284039	482721	36.96%	3.331%
43	11.731	1608	1612	1615	rBV4	9321	16467	1.26%	0.114%
44	11.768	1615	1618	1621	rVB4	10915	14394	1.10%	0.099%
45	11.817	1621	1626	1631	rBV7	10605	22768	1.74%	0.157%
46	11.878	1631	1636	1640	rBV2	25424	44716	3.42%	0.309%
47	12.036	1655	1662	1665	rBV6	7324	16976	1.30%	0.117%
48	12.140	1673	1679	1685	rBV4	45289	104677	8.02%	0.722%
49	12.256	1695	1698	1702	rVB	20200	27819	2.13%	0.192%
50	12.341	1702	1712	1722	rBV6	42248	174829	13.39%	1.206%
51	12.621	1752	1758	1766	rVB	270457	446928	34.22%	3.084%
52	12.707	1767	1772	1776	rBV5	20102	36295	2.78%	0.250%
53	12.786	1781	1785	1792	rVB6	38481	80570	6.17%	0.556%
54	12.865	1792	1798	1801	rBV5	17056	34352	2.63%	0.237%
55	12.932	1802	1809	1816	rVV8	23610	70783	5.42%	0.488%
56	13.079	1830	1833	1836	rBV4	13777	20738	1.59%	0.143%
57	13.133	1838	1842	1845	rVV2	22172	39293	3.01%	0.271%
58	13.170	1845	1848	1856	rVB5	32162	55980	4.29%	0.386%
59	13.298	1866	1869	1873	rBV4	12528	20638	1.58%	0.142%
60	13.487	1897	1900	1904	rVB4	15621	19047	1.46%	0.131%
61	13.560	1907	1912	1920	rVB	335021	538793	41.26%	3.718%
62	13.633	1920	1924	1926	rBV4	12537	19619	1.50%	0.135%
63	13.725	1934	1939	1942	rBV	53589	83631	6.40%	0.577%
64	13.761	1942	1945	1949	rVB	41231	55936	4.28%	0.386%
65	13.804	1949	1952	1959	rVB5	12609	29202	2.24%	0.202%
66	13.889	1960	1966	1971	rBV2	82277	144189	11.04%	0.995%
67	14.017	1984	1987	1993	rVB4	13360	22364	1.71%	0.154%
68	14.103	1993	2001	2006	rBV	351475	542181	41.52%	3.741%
69	14.164	2007	2011	2014	rBV	33769	46487	3.56%	0.321%
70	14.213	2014	2019	2025	rVB2	301912	485238	37.16%	3.348%
71	14.273	2025	2029	2035	rVB3	31807	60146	4.61%	0.415%

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72	14.353	2035	2042	2045	rBV	40845	79565	6.09%	0.549%
73	14.426	2045	2054	2062	rVB	285441	524449	40.16%	3.619%
74	14.505	2062	2067	2070	rBV	79148	116855	8.95%	0.806%
75	14.548	2070	2074	2082	rVB2	57597	110780	8.48%	0.764%
76	14.645	2085	2090	2094	rVB2	101636	169817	13.00%	1.172%
77	14.700	2094	2099	2102	rBV3	93673	175391	13.43%	1.210%
78	14.737	2102	2105	2110	rVB	104751	165105	12.64%	1.139%
79	14.816	2110	2118	2122	rBV2	389549	790429	60.52%	5.454%
80	14.865	2122	2126	2132	rVB2	119599	202728	15.52%	1.399%
81	15.072	2139	2160	2164	rBV2	524215	1305977	100.00%	9.012%
82	15.109	2164	2166	2172	rVV2	256633	464511	35.57%	3.205%
83	15.188	2172	2179	2189	rVB3	418152	999829	76.56%	6.899%
84	15.334	2197	2203	2212	rVB6	83347	199175	15.25%	1.374%
85	15.426	2212	2218	2221	rBV3	39843	75750	5.80%	0.523%
86	15.481	2221	2227	2230	rVV2	106611	225883	17.30%	1.559%
87	15.529	2230	2235	2250	rVB3	157054	462571	35.42%	3.192%
88	15.663	2250	2257	2264	rBV	217470	445305	34.10%	3.073%
89	15.743	2266	2270	2273	rVB	16368	22096	1.69%	0.152%
90	15.834	2274	2285	2295	rBV4	115817	415250	31.80%	2.865%
91	15.956	2300	2305	2310	rVB	67133	115525	8.85%	0.797%
92	16.029	2310	2317	2323	rBV3	78860	187102	14.33%	1.291%
93	16.182	2333	2342	2347	rBV3	45343	118148	9.05%	0.815%
94	16.230	2347	2350	2358	rVB6	15519	32312	2.47%	0.223%
95	16.352	2365	2370	2372	rBV	16660	28309	2.17%	0.195%
96	16.444	2381	2385	2389	rVV3	9255	17330	1.33%	0.120%
97	16.499	2391	2394	2406	rVB6	14385	30662	2.35%	0.212%
98	16.651	2411	2419	2428	rBV6	23918	73475	5.63%	0.507%

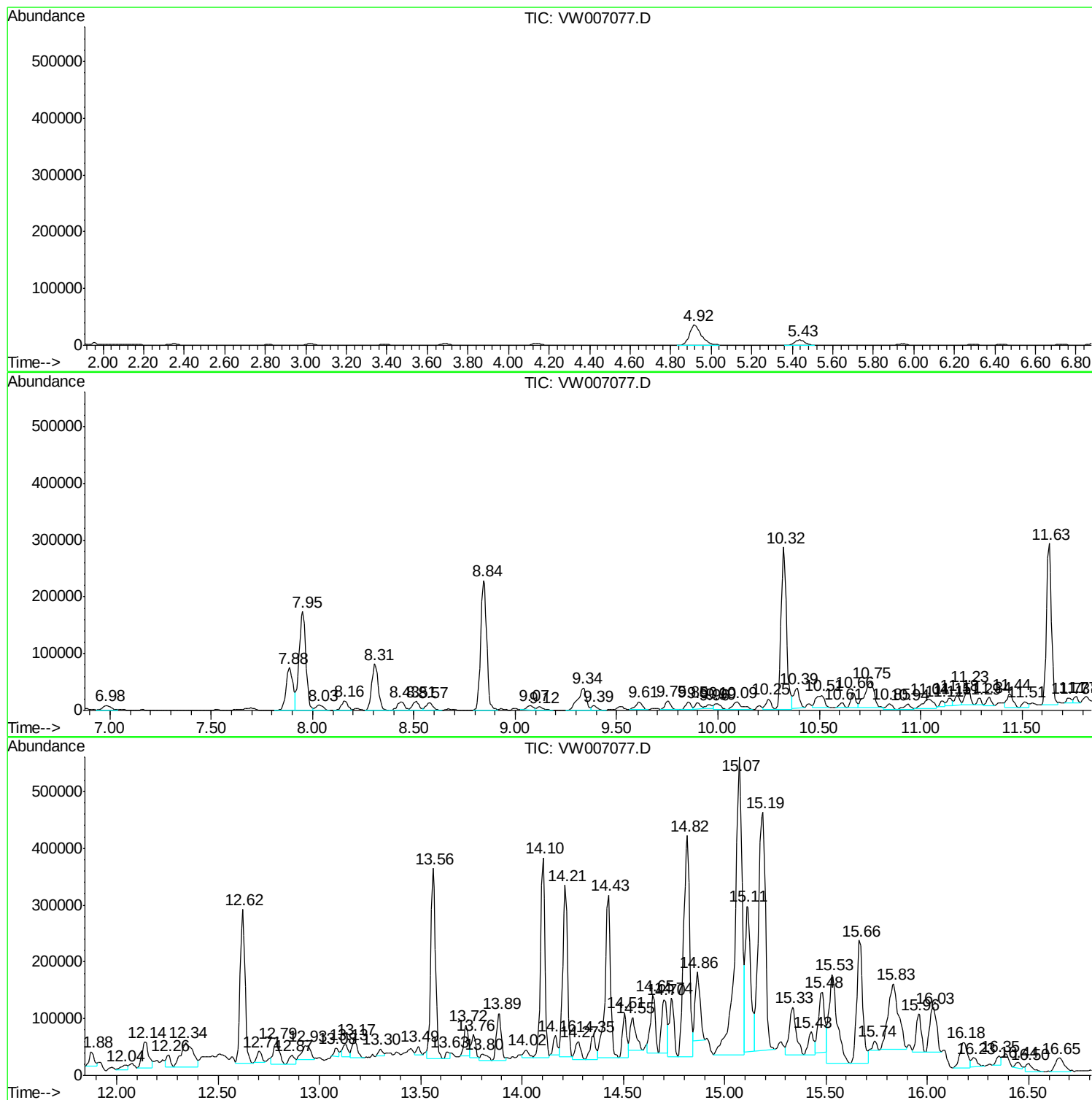
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W112118S.M
Quant Title : SW846 8260

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



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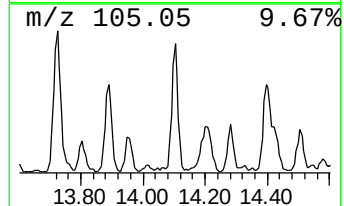
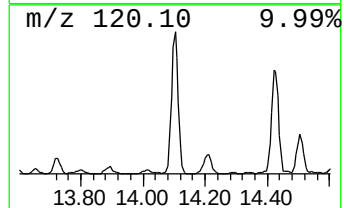
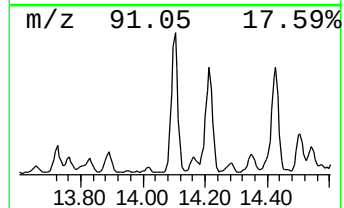
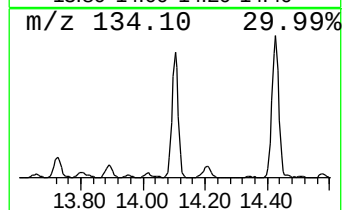
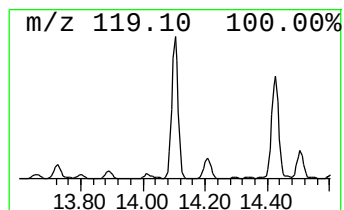
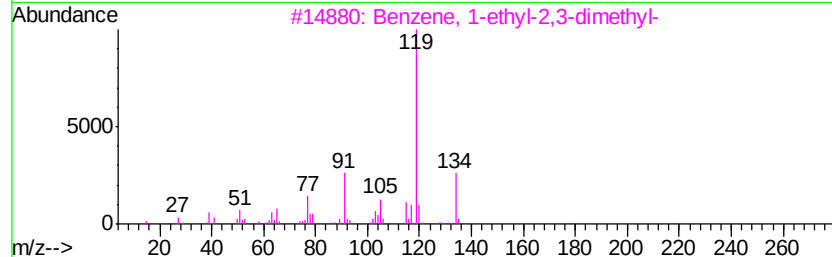
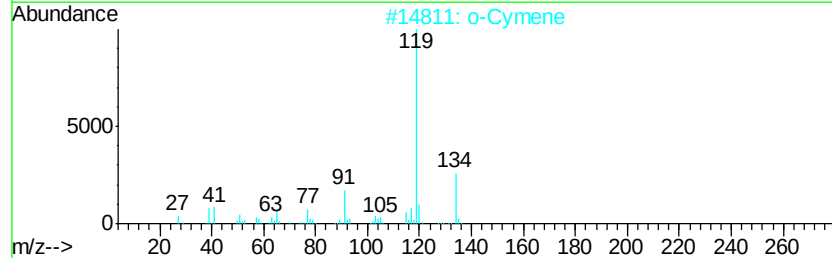
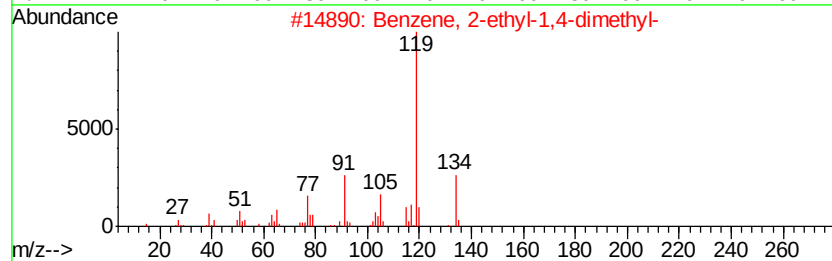
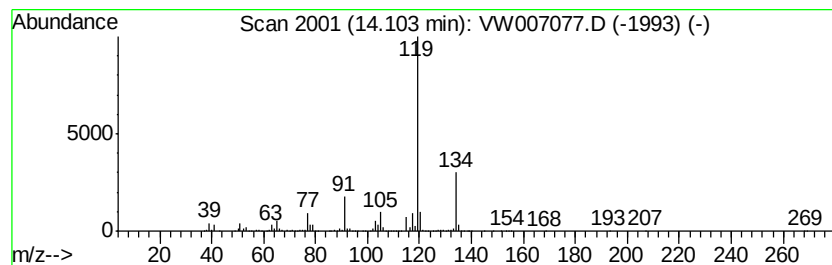
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W112118S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.10	50.31 ug/l	542181	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2		o-Cymene	134	C10H14	000527-84-4	95
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



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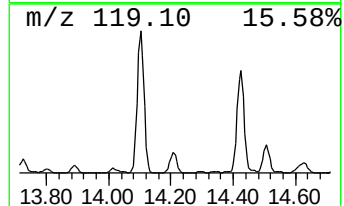
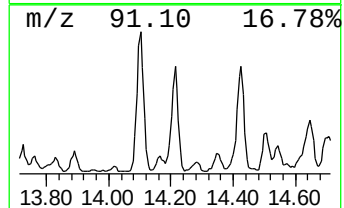
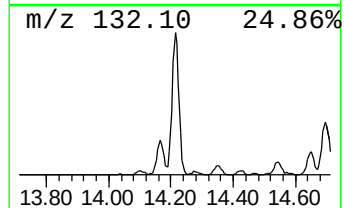
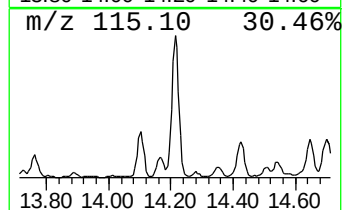
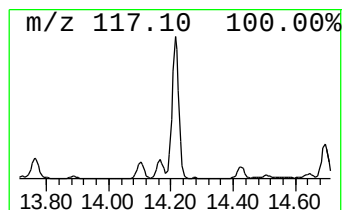
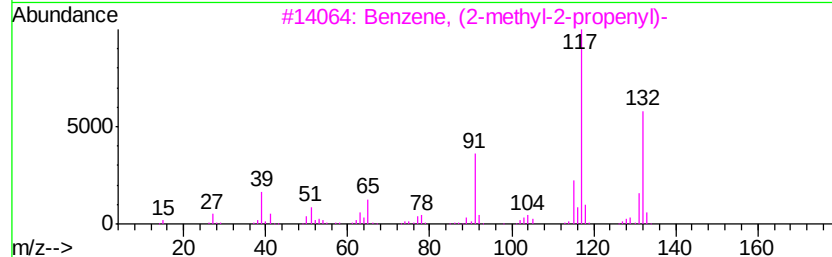
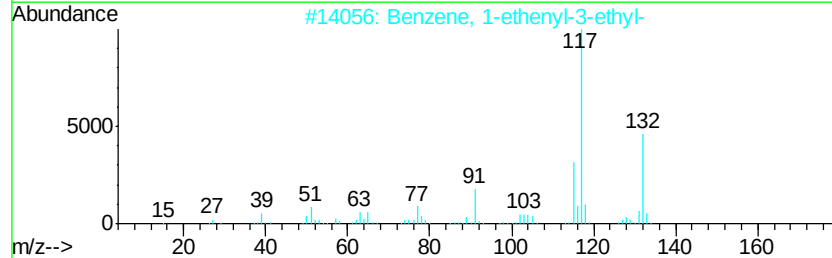
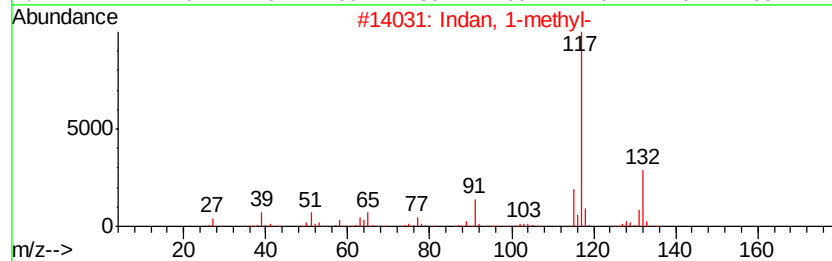
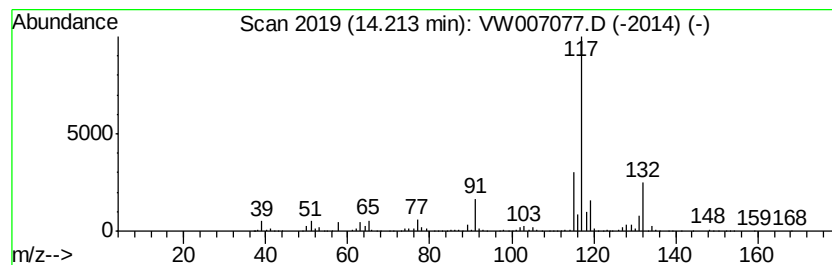
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Peak Number 2 Indan, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	45.03 ug/l	485238	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indan, 1-methyl-	132	C10H12	000767-58-8	87
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86
3		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	83
4		Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	80
5		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	80



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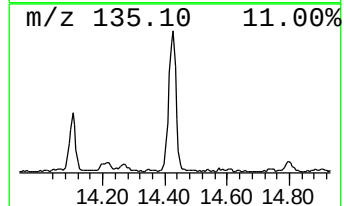
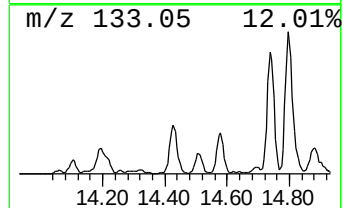
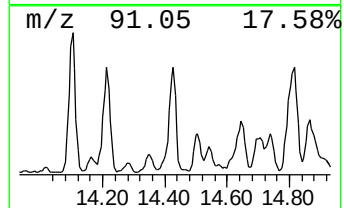
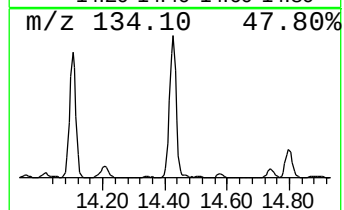
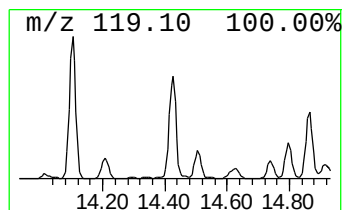
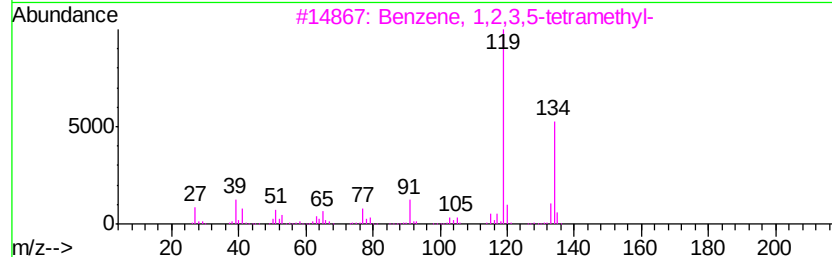
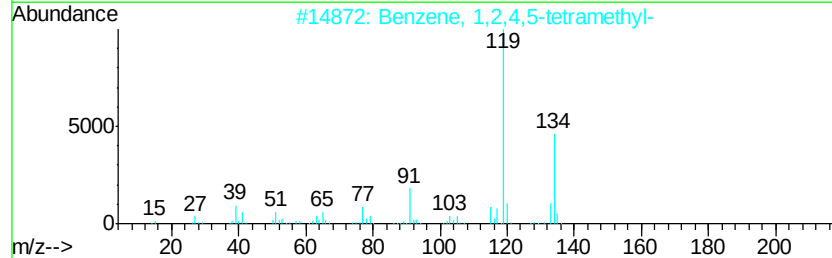
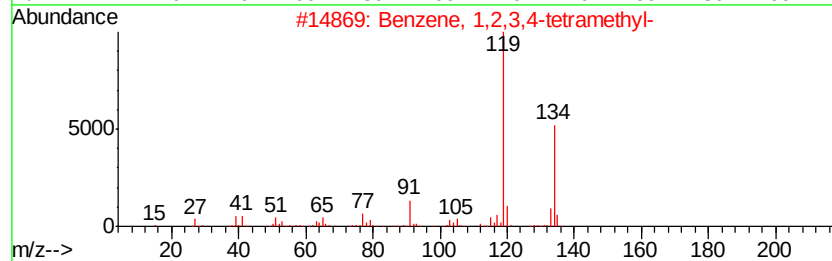
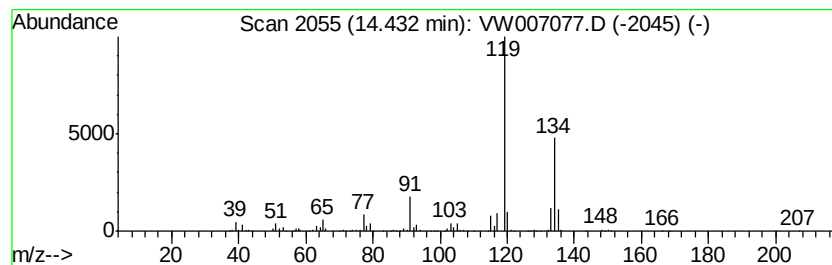
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Peak Number 3 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.43	48.67 ug/l	524449	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1,2,3,4-tetramethyl-	134 C10H14	000488-23-3	95
2	Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2	95
3	Benzene, 1,2,3,5-tetramethyl-	134 C10H14	000527-53-7	95
4	Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2	91
5	Benzene, 1-methyl-3-(1-methyleth...	134 C10H14	000535-77-3	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW112618\
Data File : VW007077.D
Acq On : 27 Nov 2018 06:01
Operator : VAY/AP
Sample : J6065-01
Misc : 5.83G/5ML/MSVOA_W/SOIL
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
USTS-1-2-3

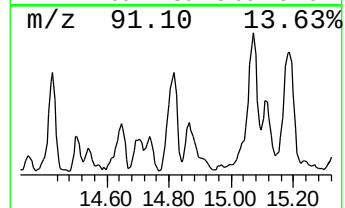
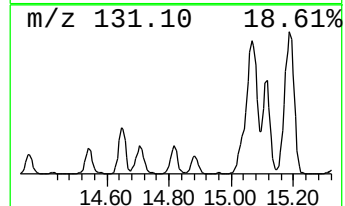
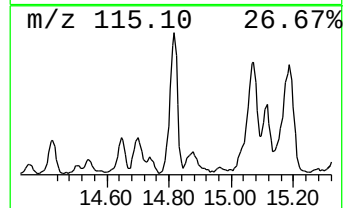
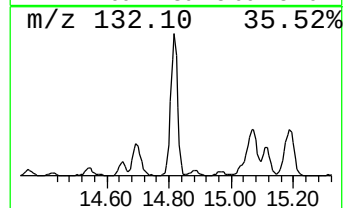
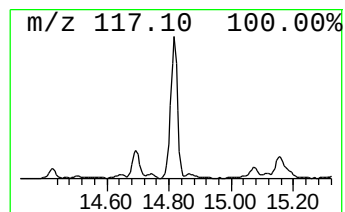
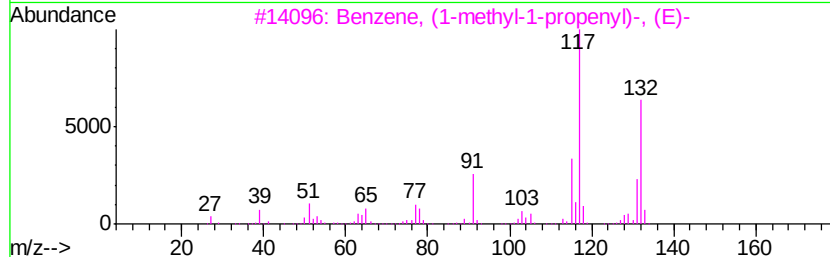
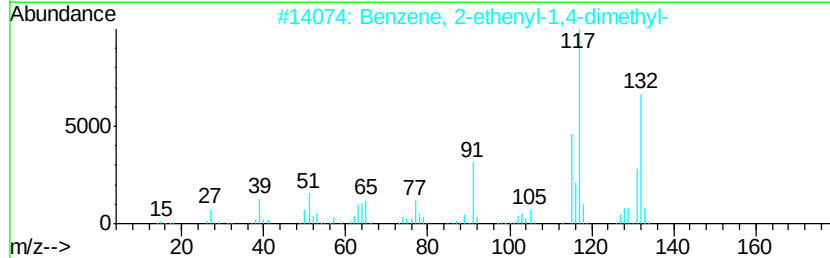
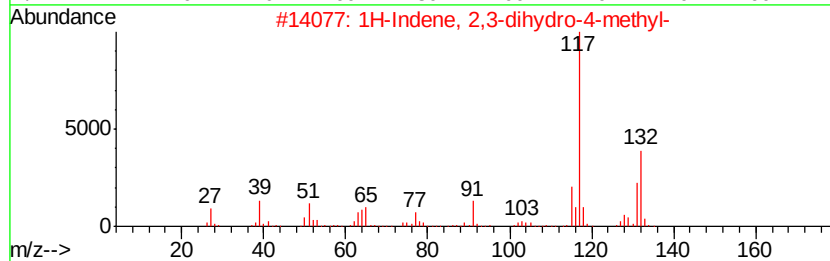
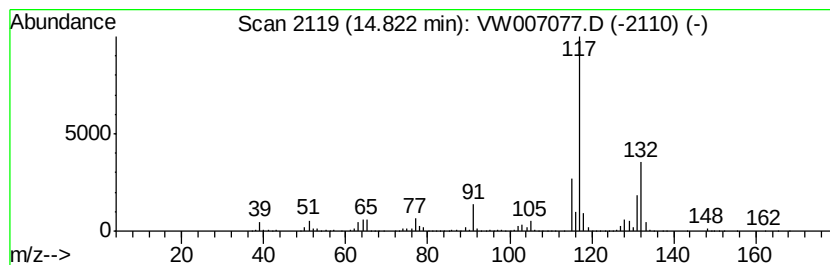
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W112118S.M
Quant Title : SW846 8260

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

Peak Number 4 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	73.35 ug/l	790429	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91
4		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	91
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW112618\
Data File : VW007077.D
Acq On : 27 Nov 2018 06:01
Operator : VAY/AP
Sample : J6065-01
Misc : 5.83G/5ML/MSVOA_W/SOIL
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
USTS-1-2-3

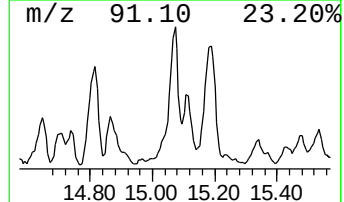
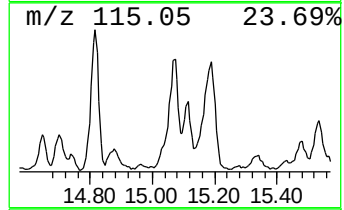
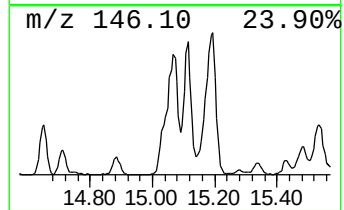
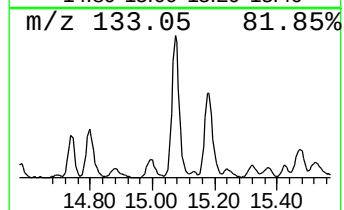
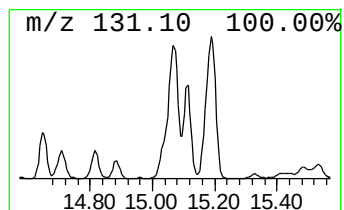
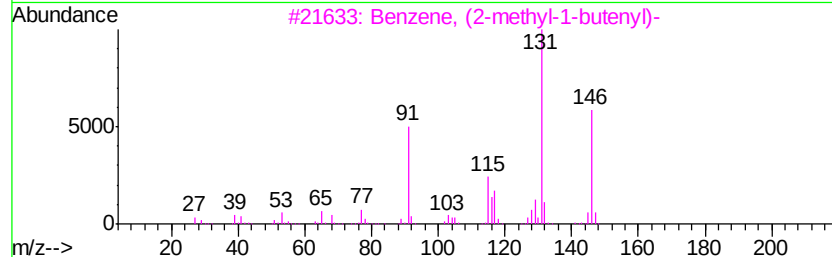
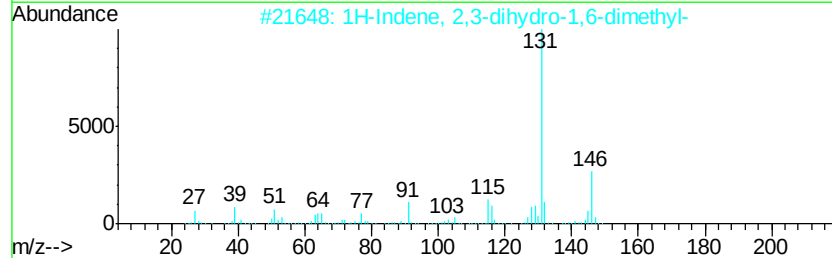
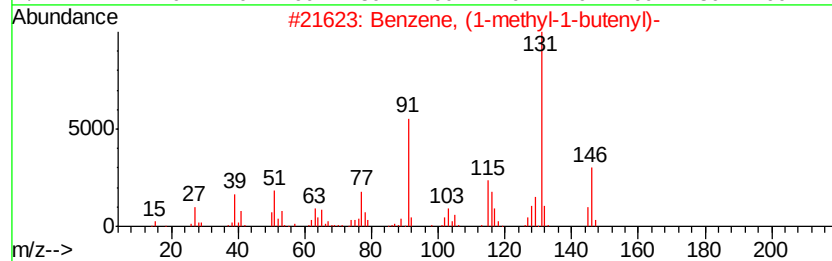
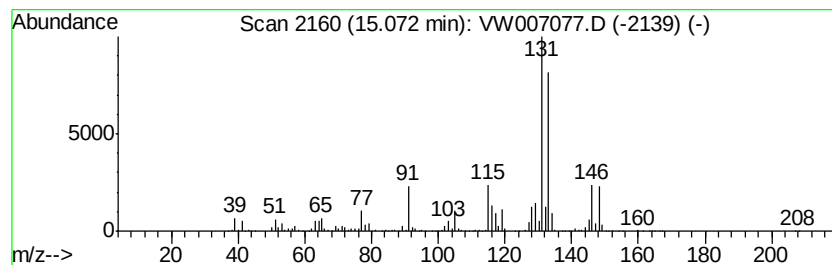
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W112118S.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, (1-methyl-1-butenyl)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	121.20 ug/l	1305980	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	90
2		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	86
3		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	64
4		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	64
5		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW112618\
Data File : VW007077.D
Acq On : 27 Nov 2018 06:01
Operator : VAY/AP
Sample : J6065-01
Misc : 5.83G/5ML/MSVOA_W/SOIL
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
USTS-1-2-3

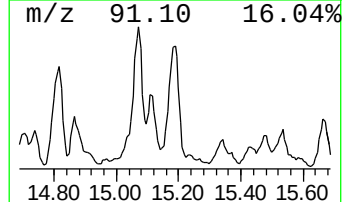
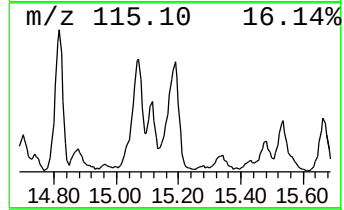
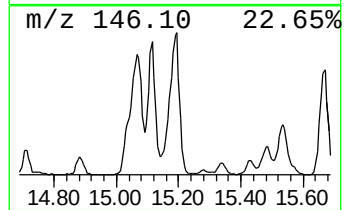
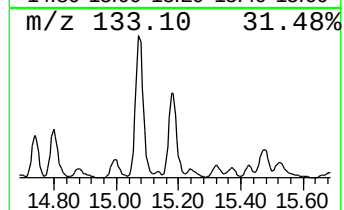
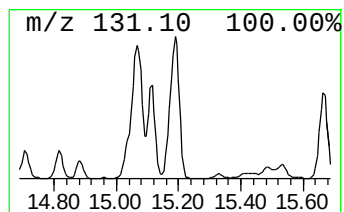
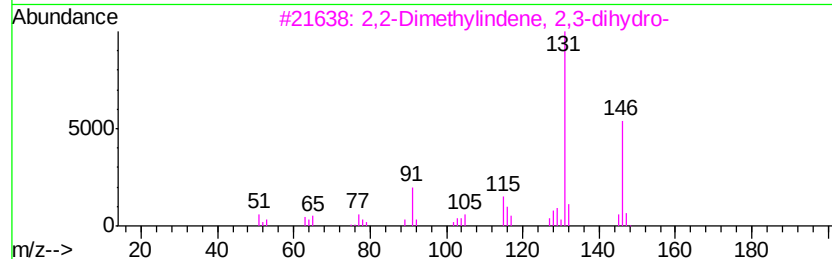
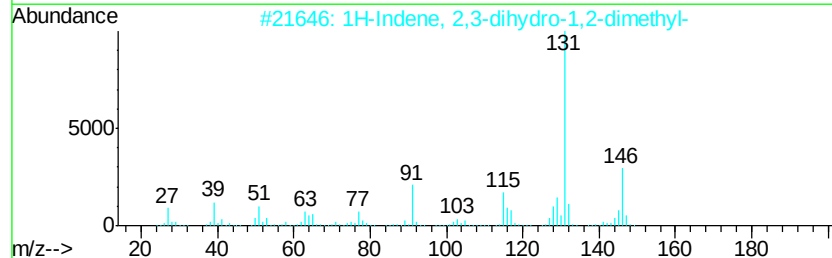
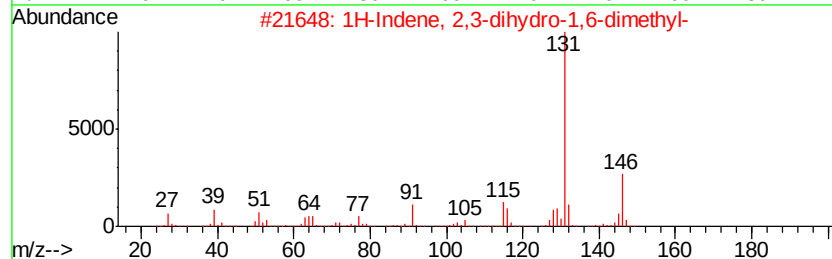
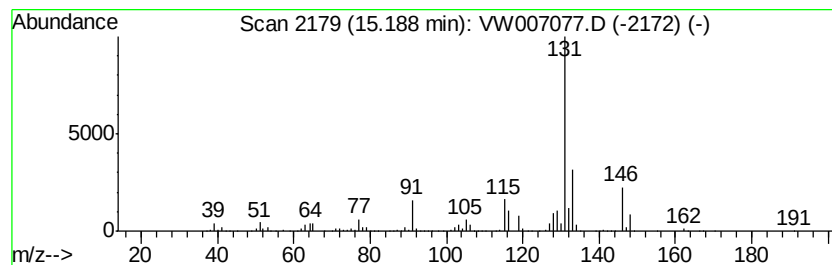
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W112118S.M
Quant Title : SW846 8260

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

Peak Number 7 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	92.78 ug/l	999829	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	93
2		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	87
3		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	87
4		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	87
5		Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW112618\
Data File : VW007077.D
Acq On : 27 Nov 2018 06:01
Operator : VAY/AP
Sample : J6065-01
Misc : 5.83G/5ML/MSVOA_W/SOIL
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
USTS-1-2-3

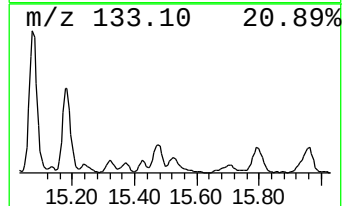
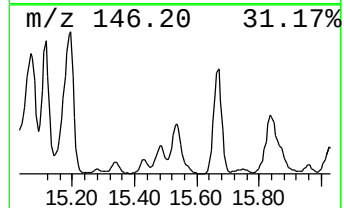
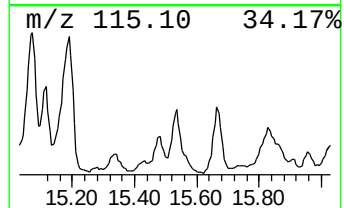
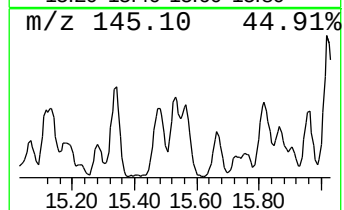
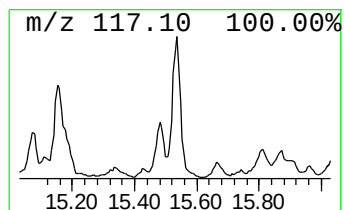
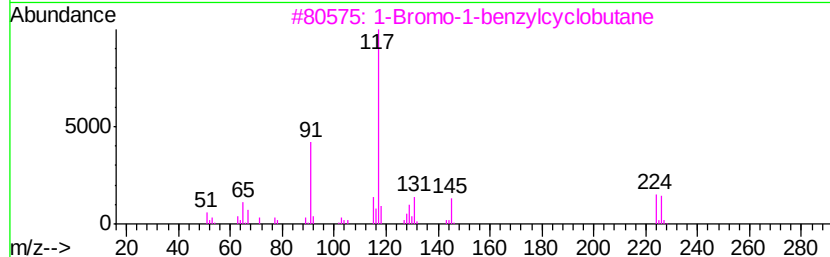
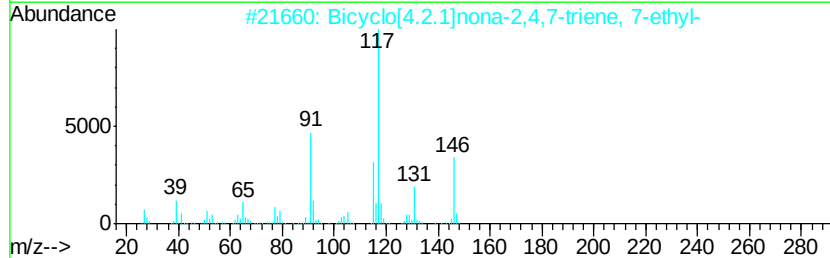
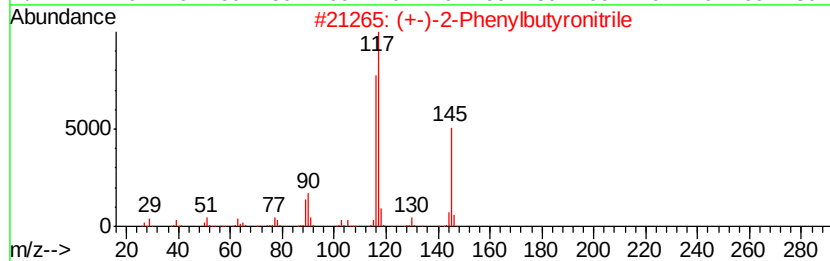
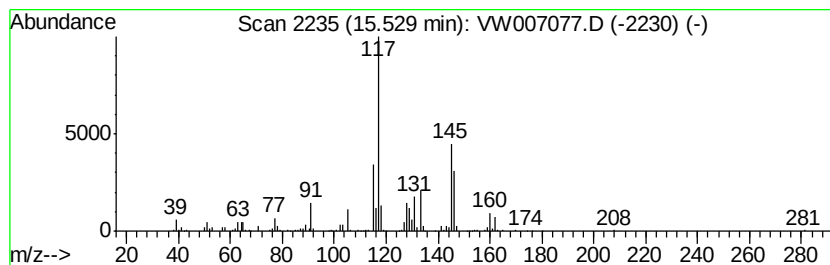
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W112118S.M
Quant Title : SW846 8260

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

Peak Number 8 unknown15.53 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.53	42.93 ug/l	462571	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(+)-2-Phenylbutyronitrile	145	C10H11N	069350-73-8	49
2		Bicyclo[4.2.1]nona-2,4,7-triene,...	146	C11H14	1000164-42-6	47
3		1-Bromo-1-benzylcyclobutane	224	C11H13Br	1000214-92-1	47
4		Benzene, 1-pentenyl-	146	C11H14	000826-18-6	46
5		1H-Indene, 1-ethyl-2,3-dihydro-	146	C11H14	004830-99-3	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW112618\
Data File : VW007077.D
Acq On : 27 Nov 2018 06:01
Operator : VAY/AP
Sample : J6065-01
Misc : 5.83G/5ML/MSVOA_W/SOIL
ALS Vial : 47 Sample Multiplier: 1

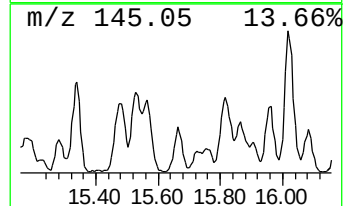
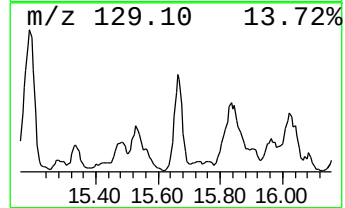
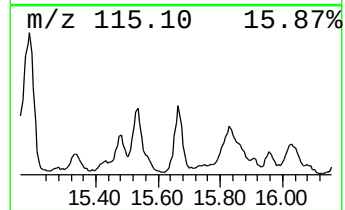
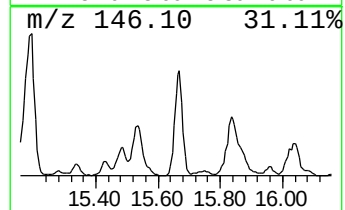
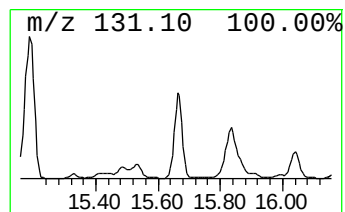
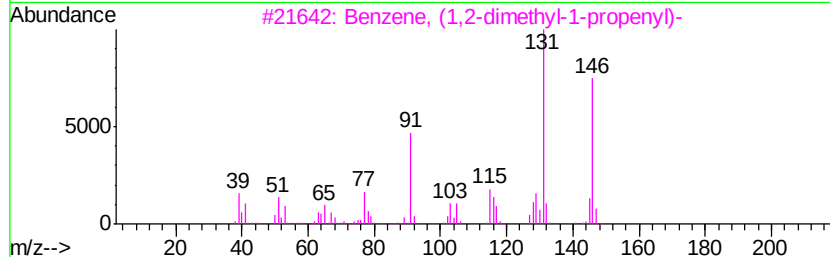
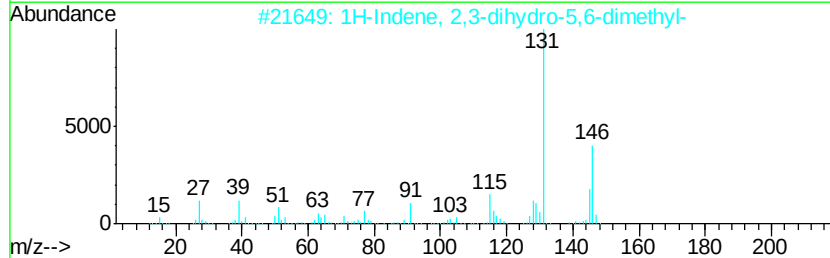
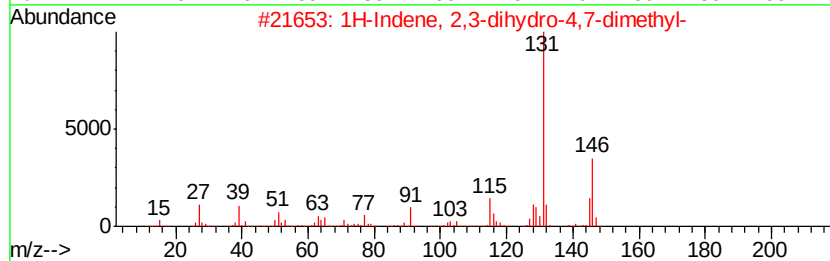
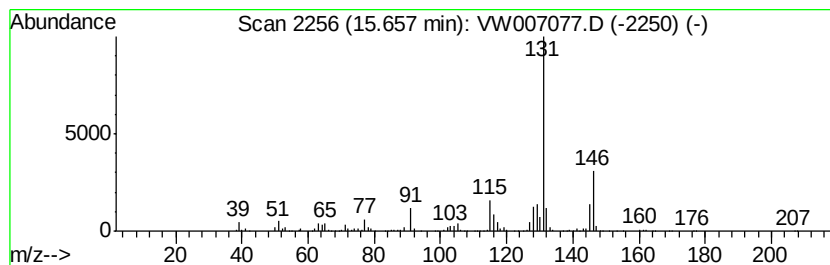
Instrument :
MSVOA_W
ClientSampleId :
USTS-1-2-3

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W112118S.M
Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.66	41.32 ug/l	445305	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9	97
2	1H-Indene, 2,3-dihydro-5,6-dimet...	146 C11H14	001075-22-5	94
3	Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3	93
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5	91
5	1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW112618\
Data File : VW007077.D
Acq On : 27 Nov 2018 06:01
Operator : VAY/AP
Sample : J6065-01
Misc : 5.83G/5ML/MSVOA_W/SOIL
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
USTS-1-2-3

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W112118S.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 2-ethyl-...	14.10	50.3	ug/l	542181	4	13.56	538793	50.0
Indan, 1-methyl-	14.21	45.0	ug/l	485238	4	13.56	538793	50.0
Benzene, 1,2,3,4-...	14.43	48.7	ug/l	524449	4	13.56	538793	50.0
1H-Indene, 2,3-di...	14.82	73.3	ug/l	790429	4	13.56	538793	50.0
Benzene, (1-methy...	15.07	121.2	ug/l	1305980	4	13.56	538793	50.0
1H-Indene, 2,3-di...	15.19	92.8	ug/l	999829	4	13.56	538793	50.0
unknown15.53	15.53	42.9	ug/l	462571	4	13.56	538793	50.0
1H-Indene, 2,3-di...	15.66	41.3	ug/l	445305	4	13.56	538793	50.0