

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW021419\
 Data File : VW008666.D
 Acq On : 14 Feb 2019 15:13
 Operator : SY/VA
 Sample : K1349-05
 Misc : 7.37G/5ML/MSVOA W/SOIL
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 OR-01-021319-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\82W021319S.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	4.123	356	364	372	rBV2	24228	69789	3.46%	0.344%
2	4.909	480	493	508	rBV2	36882	148086	7.34%	0.731%
3	7.141	848	859	872	rBV2	57480	175165	8.69%	0.864%
4	7.878	970	980	985	rBV	284652	667035	33.08%	3.292%
5	7.946	985	991	1001	rVB	532392	1174037	58.22%	5.794%
6	8.250	1030	1041	1042	rBV	25895	45607	2.26%	0.225%
7	8.305	1042	1050	1060	rVB2	278923	640962	31.78%	3.163%
8	8.842	1127	1138	1148	rBV	735008	1462943	72.55%	7.220%
9	9.329	1207	1218	1225	rBV3	21736	58504	2.90%	0.289%
10	9.750	1282	1287	1294	rVB6	11441	23341	1.16%	0.115%
11	10.079	1333	1341	1347	rBV6	8997	27820	1.38%	0.137%
12	10.323	1372	1381	1387	rVV	1147836	2016566	100.00%	9.952%
13	10.390	1387	1392	1398	rVV	110616	203885	10.11%	1.006%
14	10.494	1402	1409	1417	rVB6	15900	43598	2.16%	0.215%
15	10.665	1432	1437	1440	rBV2	19273	32626	1.62%	0.161%
16	10.701	1440	1443	1449	rVV	28477	46551	2.31%	0.230%
17	10.933	1478	1481	1487	rVB3	11913	21611	1.07%	0.107%
18	11.231	1524	1530	1536	rVB	82340	141297	7.01%	0.697%
19	11.335	1542	1547	1551	rVB4	15662	27193	1.35%	0.134%
20	11.439	1551	1564	1571	rVB8	33513	86685	4.30%	0.428%
21	11.512	1571	1576	1581	rBV3	12407	21397	1.06%	0.106%
22	11.628	1587	1595	1606	rBV	945606	1620656	80.37%	7.998%
23	11.731	1606	1612	1615	rBV2	29250	51203	2.54%	0.253%
24	11.835	1621	1629	1634	rBV	96252	206881	10.26%	1.021%
25	11.975	1647	1652	1658	rBV8	9373	21896	1.09%	0.108%
26	12.042	1658	1663	1670	rVV9	8405	23805	1.18%	0.117%
27	12.164	1670	1683	1689	rVV2	81487	229405	11.38%	1.132%
28	12.256	1690	1698	1701	rVB3	26661	58034	2.88%	0.286%
29	12.341	1707	1712	1713	rBV3	29546	39872	1.98%	0.197%
30	12.615	1751	1757	1765	rBV	783949	1339947	66.45%	6.613%
31	12.701	1767	1771	1776	rVB8	11570	21484	1.07%	0.106%
32	12.780	1779	1784	1792	rVB4	69141	148356	7.36%	0.732%
33	12.865	1792	1798	1802	rBV2	95672	172197	8.54%	0.850%
34	12.902	1802	1804	1806	rVV	33609	45600	2.26%	0.225%

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35	12.945	1806	1811	1816	rVB3	105715	211879	10.51%	1.046%
36	13.103	1830	1837	1843	rVV	72583	163153	8.09%	0.805%
37	13.164	1843	1847	1852	rVV5	38756	65697	3.26%	0.324%
38	13.256	1857	1862	1866	rVV	157643	245104	12.15%	1.210%
39	13.298	1866	1869	1872	rVB3	17148	23719	1.18%	0.117%
40	13.384	1880	1883	1887	rVB3	25804	34894	1.73%	0.172%
41	13.445	1887	1893	1897	rBV4	62956	112214	5.56%	0.554%
42	13.493	1897	1901	1906	rVV5	31649	57183	2.84%	0.282%
43	13.560	1906	1912	1921	rVV2	854858	1482348	73.51%	7.316%
44	13.719	1930	1938	1940	rBV8	44123	101586	5.04%	0.501%
45	13.755	1940	1944	1947	rVV2	163667	267087	13.24%	1.318%
46	13.792	1947	1950	1960	rVB3	107529	245905	12.19%	1.214%
47	13.877	1960	1964	1969	rBV3	134667	210841	10.46%	1.041%
48	13.951	1971	1976	1981	rVV2	69706	119692	5.94%	0.591%
49	14.018	1982	1987	1989	rVV	73763	131277	6.51%	0.648%
50	14.042	1989	1991	1995	rVV	88310	133397	6.62%	0.658%
51	14.097	1995	2000	2005	rVB2	119816	221456	10.98%	1.093%
52	14.213	2013	2019	2024	rVB2	140815	235738	11.69%	1.163%
53	14.274	2024	2029	2035	rVB6	21998	52033	2.58%	0.257%
54	14.341	2035	2040	2044	rBV2	48547	91141	4.52%	0.450%
55	14.389	2044	2048	2051	rVV5	60257	95501	4.74%	0.471%
56	14.426	2051	2054	2057	rVV2	62526	92569	4.59%	0.457%
57	14.463	2057	2060	2063	rVB	96511	124311	6.16%	0.614%
58	14.505	2063	2067	2070	rBV3	63049	77855	3.86%	0.384%
59	14.548	2070	2074	2080	rVB4	56344	101498	5.03%	0.501%
60	14.609	2080	2084	2086	rBV5	23343	33394	1.66%	0.165%
61	14.646	2086	2090	2094	rVB2	54726	73691	3.65%	0.364%
62	14.694	2094	2098	2102	rBV2	124671	245896	12.19%	1.214%
63	14.731	2102	2104	2109	rVB2	116635	155829	7.73%	0.769%
64	14.810	2109	2117	2122	rBV3	258833	607734	30.14%	2.999%
65	14.859	2123	2125	2133	rVB4	59832	98622	4.89%	0.487%
66	14.963	2137	2142	2148	rVB2	171660	277825	13.78%	1.371%
67	15.030	2149	2153	2155	rBV5	29294	46741	2.32%	0.231%
68	15.066	2155	2159	2164	rBV3	73969	123135	6.11%	0.608%
69	15.182	2172	2178	2186	rVB3	99335	230501	11.43%	1.138%
70	15.255	2186	2190	2197	rVB8	27197	64775	3.21%	0.320%
71	15.334	2198	2203	2205	rBV4	48834	85603	4.24%	0.422%

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72	15.371	2205	2209	2214	rVB	136364	216798	10.75%	1.070%
73	15.426	2214	2218	2222	rBV	71811	107391	5.33%	0.530%
74	15.481	2222	2227	2229	rBV5	61547	99311	4.92%	0.490%
75	15.560	2237	2240	2249	rVB4	81796	156837	7.78%	0.774%
76	15.664	2250	2257	2264	rBV3	85178	199602	9.90%	0.985%
77	15.737	2266	2269	2275	rVV6	23515	39629	1.97%	0.196%
78	15.834	2275	2285	2286	rVV4	54606	117880	5.85%	0.582%
79	15.871	2286	2291	2298	rVV	174338	324407	16.09%	1.601%
80	15.956	2302	2305	2309	rVV5	16959	26129	1.30%	0.129%
81	16.029	2312	2317	2324	rVB4	34924	86978	4.31%	0.429%
82	16.188	2333	2343	2349	rBV5	84925	225779	11.20%	1.114%
83	16.383	2367	2375	2380	rBV3	77372	167319	8.30%	0.826%
84	16.450	2380	2386	2392	rBV	134931	273263	13.55%	1.349%
85	16.651	2411	2419	2427	rBV	147467	364860	18.09%	1.801%

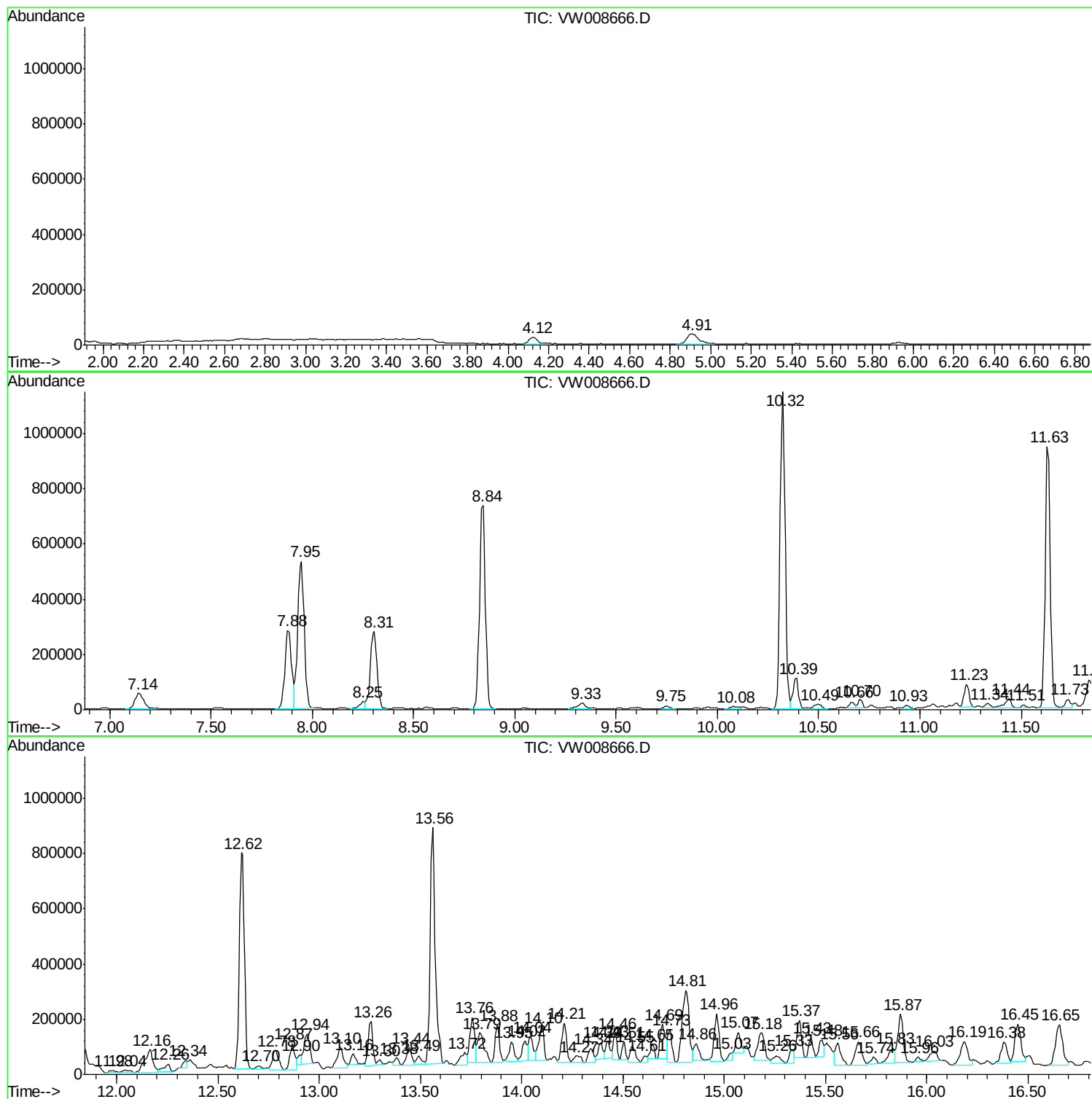
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W021319S.M
Quant Title : SW846 8260

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



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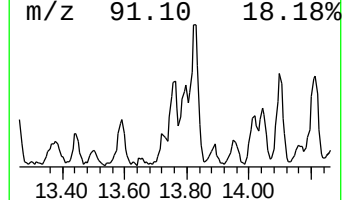
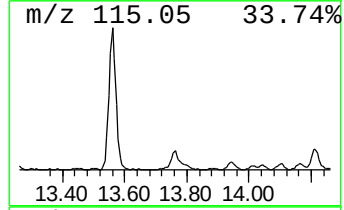
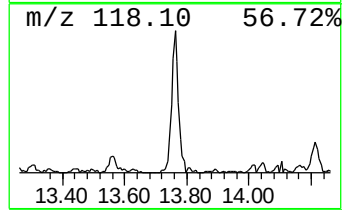
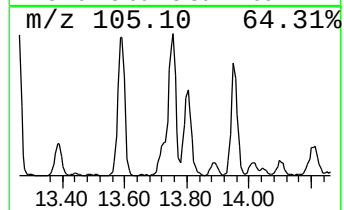
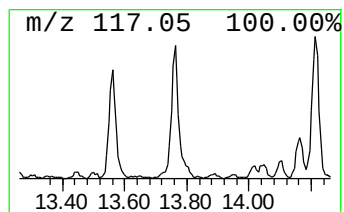
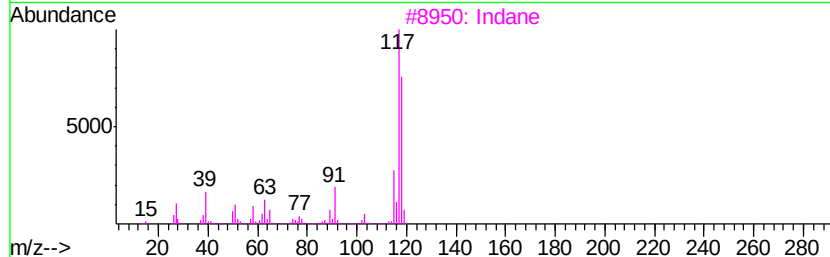
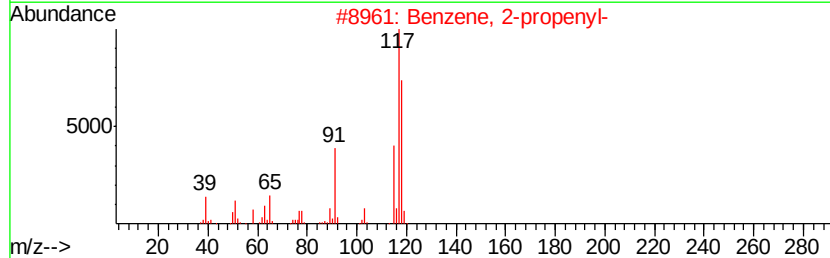
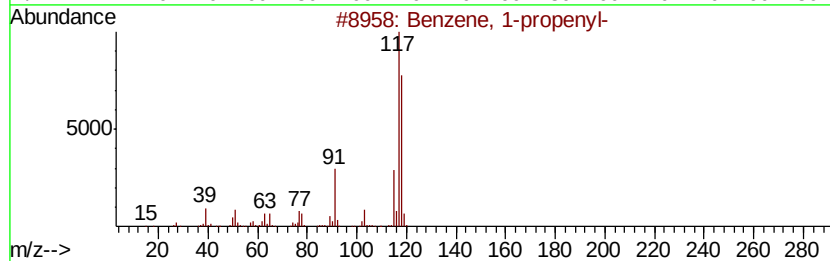
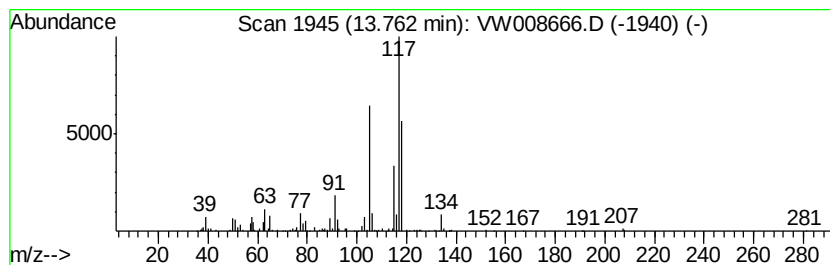
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W021319S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-propenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	9.01 ug/l	267087	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
2		Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
3		Indane	118	C9H10	000496-11-7	64
4		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	64
5		Benzene, cyclopropyl-	118	C9H10	000873-49-4	58



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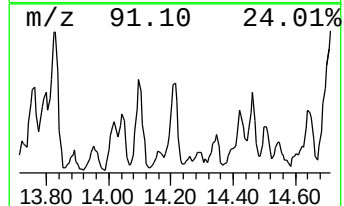
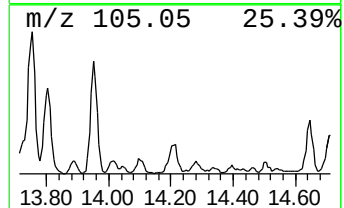
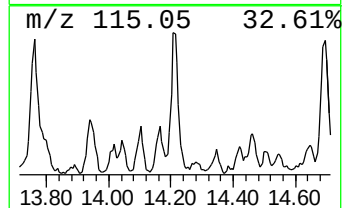
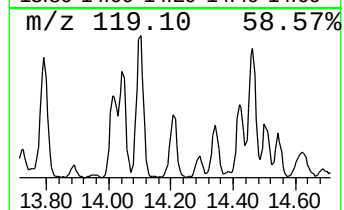
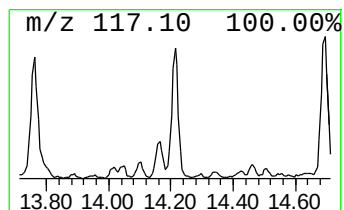
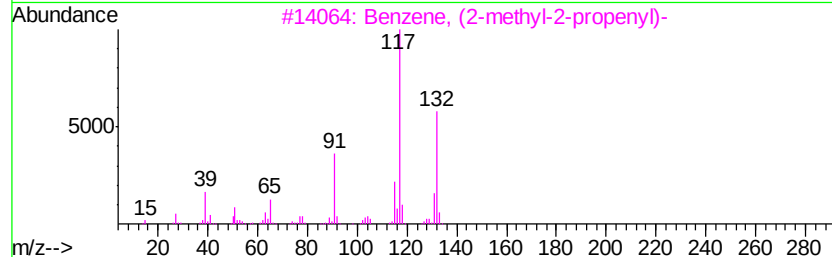
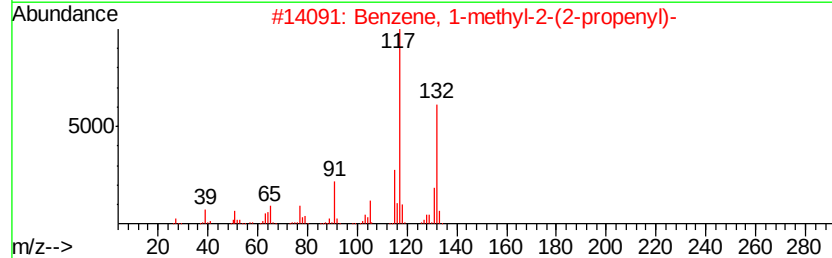
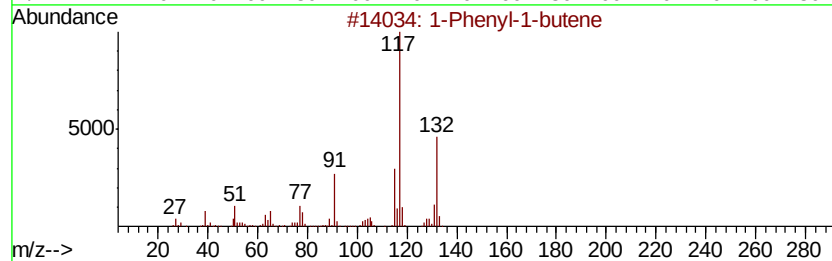
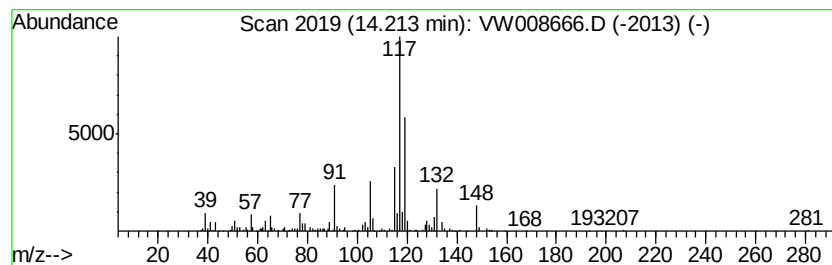
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 1-Phenyl-1-butene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	7.95 ug/l	235738	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	64
2		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	62
3		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	58
4		Indan, 1-methyl-	132	C10H12	000767-58-8	58
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	58



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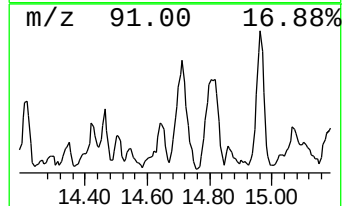
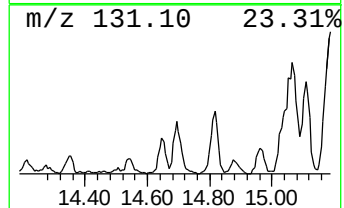
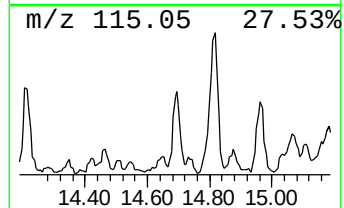
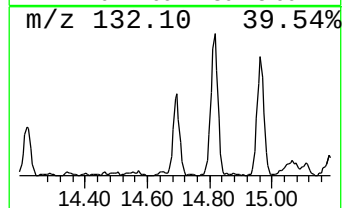
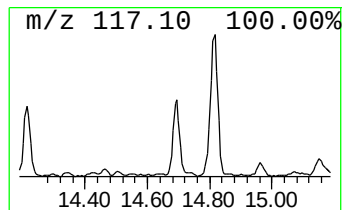
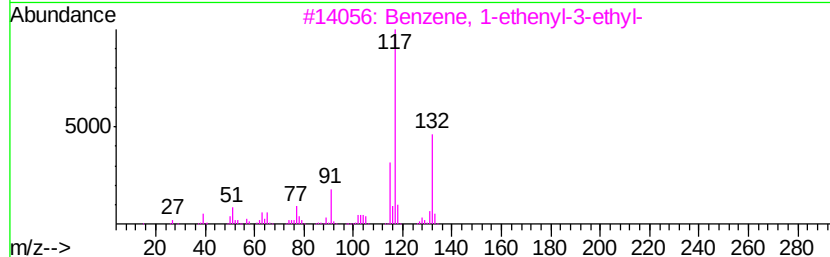
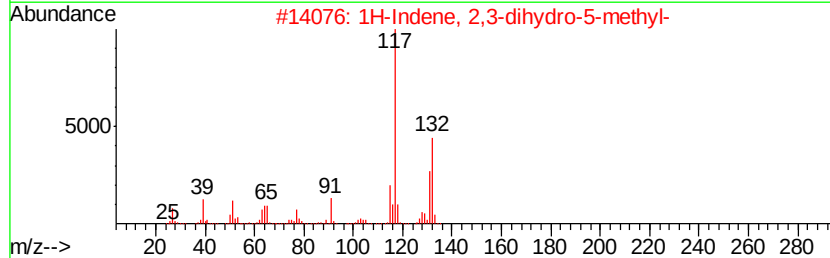
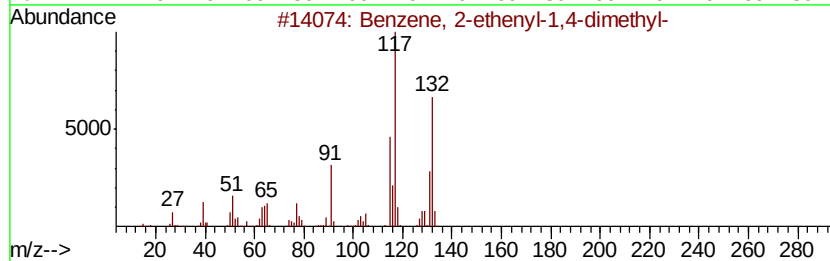
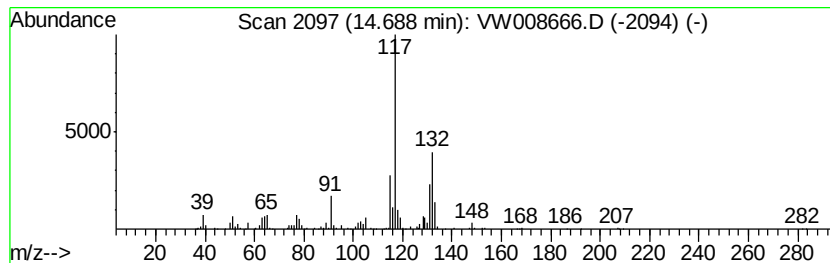
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	8.29 ug/l	245896	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 2-ethenyl-1,4-dimethyl-	132 C10H12	002039-89-6 91
2		1H-Indene, 2,3-dihydro-5-methyl-	132 C10H12	000874-35-1 90
3		Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4 87
4		Benzene, (2-methyl-2-propenyl)-	132 C10H12	003290-53-7 87
5		1-Phenyl-1-butene	132 C10H12	000824-90-8 87



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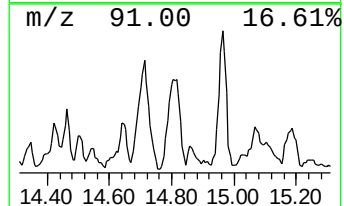
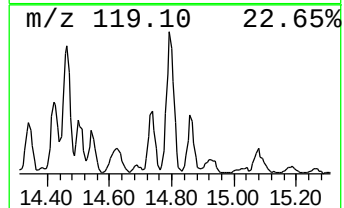
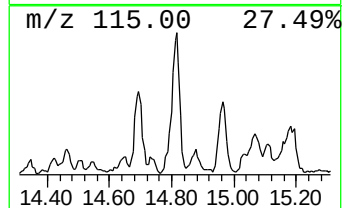
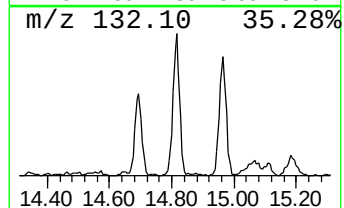
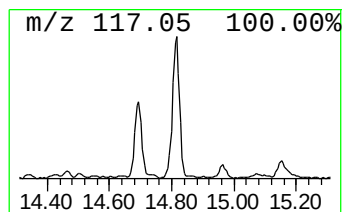
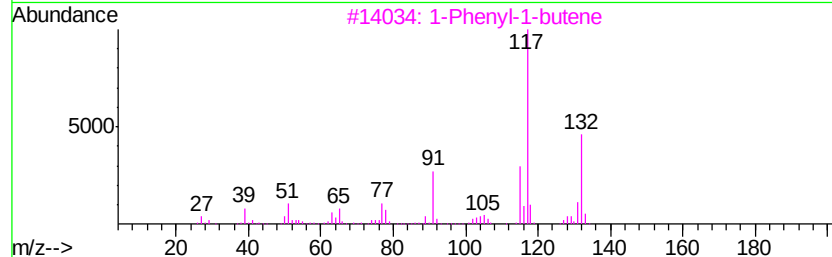
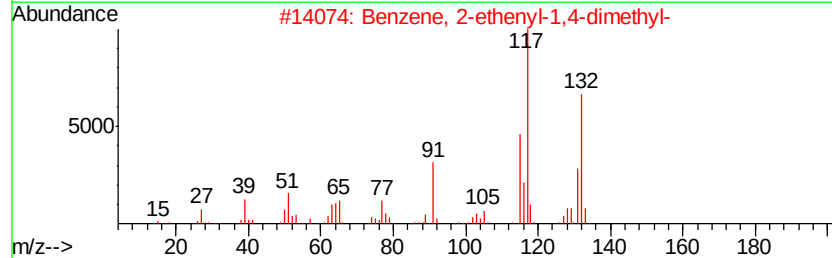
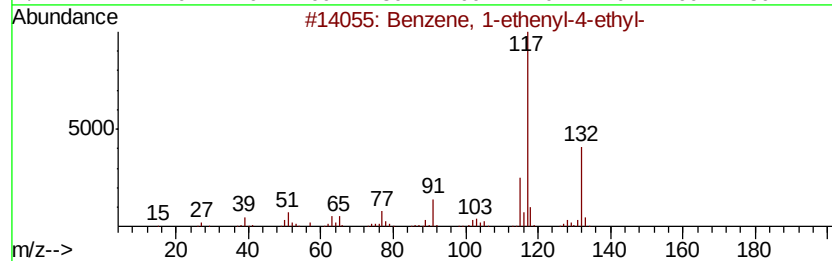
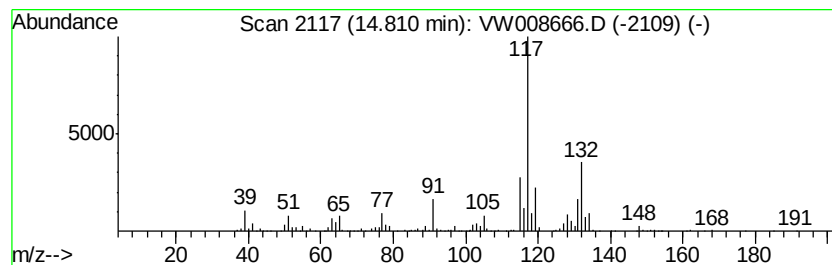
Instrument :
MSVOA_W
ClientSampleId :
OR-01-021319-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W021319S.M
Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.81	20.50 ug/l	607734	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-4-ethyl-	132 C10H12	003454-07-7 94
2		Benzene, 2-ethenyl-1,4-dimethyl-	132 C10H12	002039-89-6 90
3		1-Phenyl-1-butene	132 C10H12	000824-90-8 87
4		Benzene, 2-butenyl-	132 C10H12	001560-06-1 87
5		3-Phenylbut-1-ene	132 C10H12	000934-10-1 81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW021419\
Data File : VW008666.D
Acq On : 14 Feb 2019 15:13
Operator : SY/VA
Sample : K1349-05
Misc : 7.37G/5ML/MSVOA W/SOIL
ALS Vial : 9 Sample Multiplier: 1

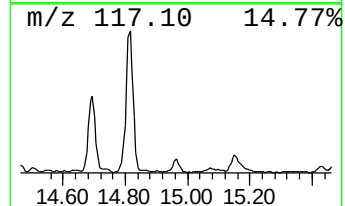
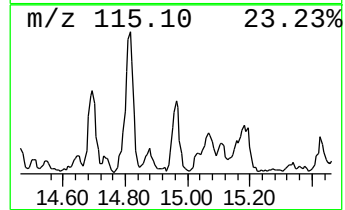
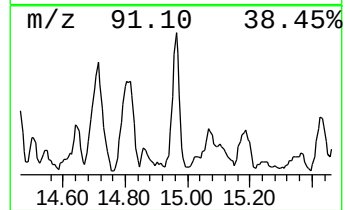
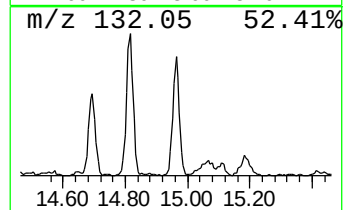
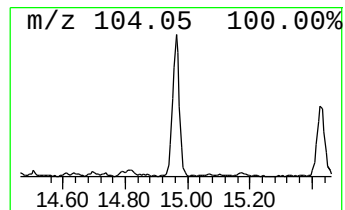
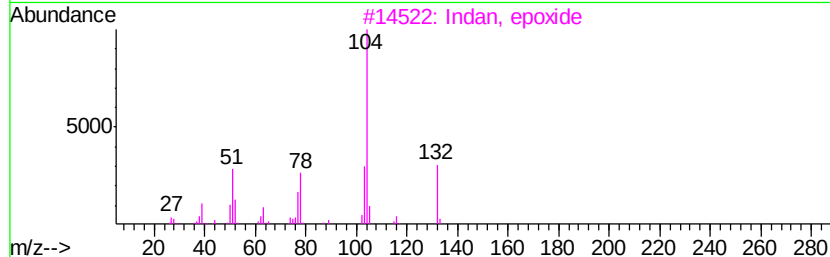
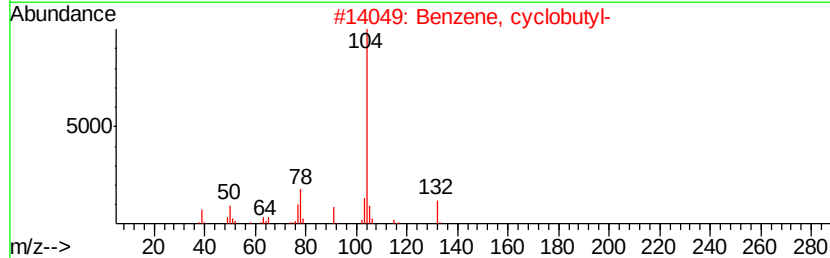
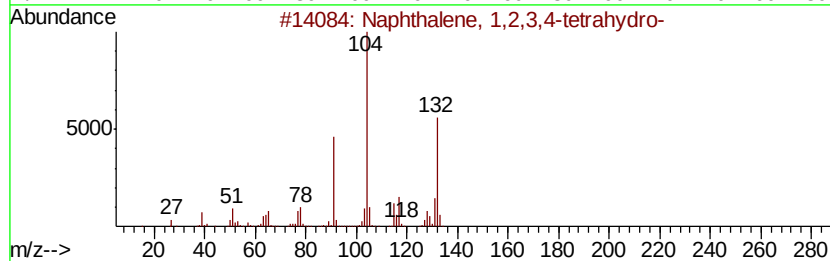
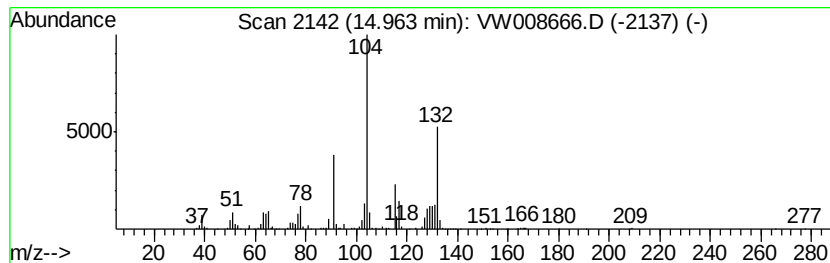
Instrument :
MSVOA_W
ClientSampleId :
OR-01-021319-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W021319S.M
Quant Title : SW846 8260

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

Peak Number 5 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	9.37 ug/l	277825	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-	132 C10H12	000119-64-2 95
2		Benzene, cyclobutyl-	132 C10H12	004392-30-7 50
3		Indan, epoxide	132 C9H8O	000768-22-9 49
4		2H-Inden-2-one, 1,3-dihydro-	132 C9H8O	000615-13-4 49
5		2H-Cyclopenta[d]pyridazine, 2-me...	132 C8H8N2	022291-85-6 38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW021419\
Data File : VW008666.D
Acq On : 14 Feb 2019 15:13
Operator : SY/VA
Sample : K1349-05
Misc : 7.37G/5ML/MSVOA W/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
OR-01-021319-A

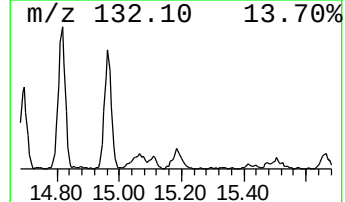
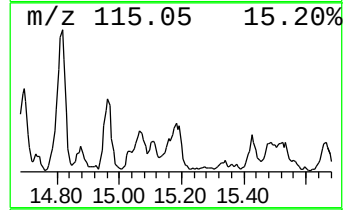
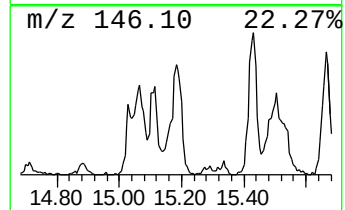
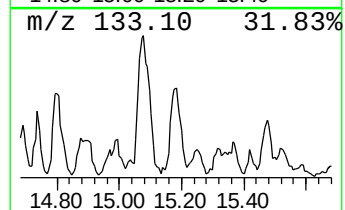
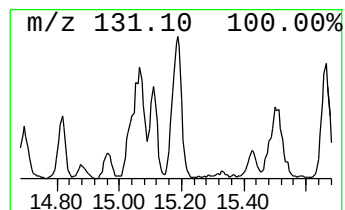
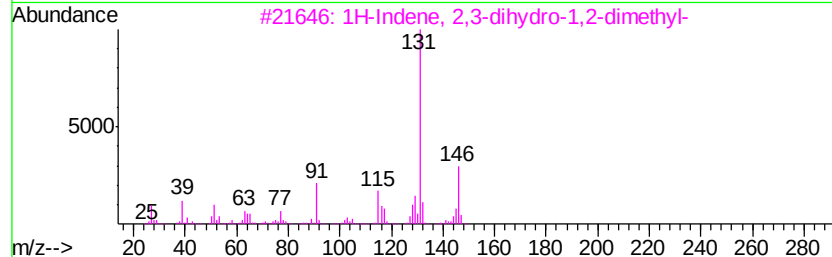
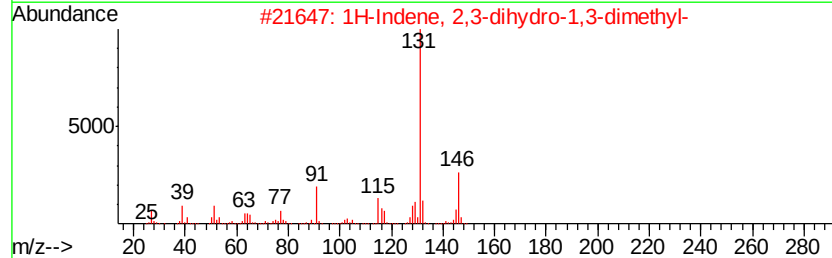
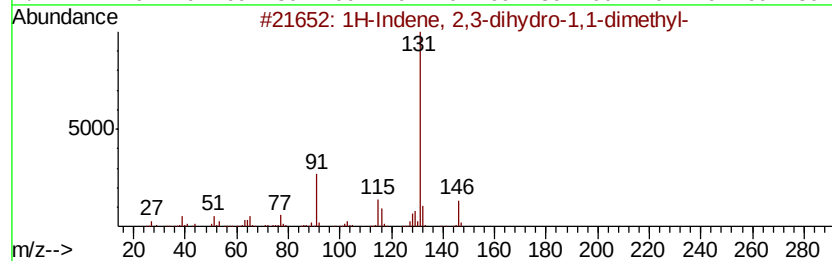
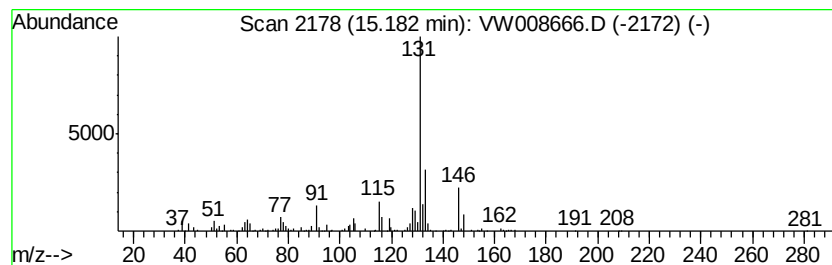
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W021319S.M
Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 2,3-dihydro-1,1-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.18	7.77 ug/l	230501	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	81
2		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	81
3		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	81
4		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	81
5		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW021419\
Data File : VW008666.D
Acq On : 14 Feb 2019 15:13
Operator : SY/VA
Sample : K1349-05
Misc : 7.37G/5ML/MSVOA W/SOIL
ALS Vial : 9 Sample Multiplier: 1

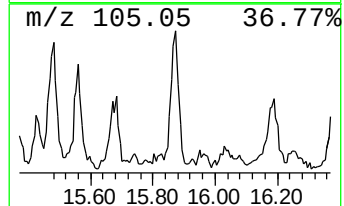
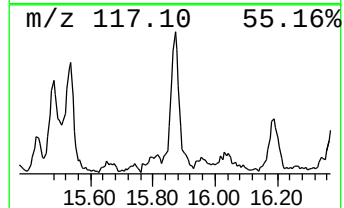
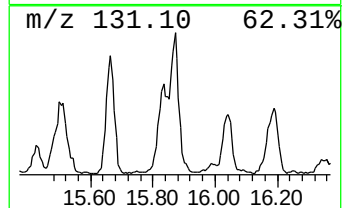
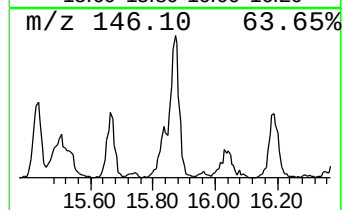
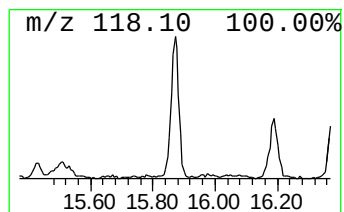
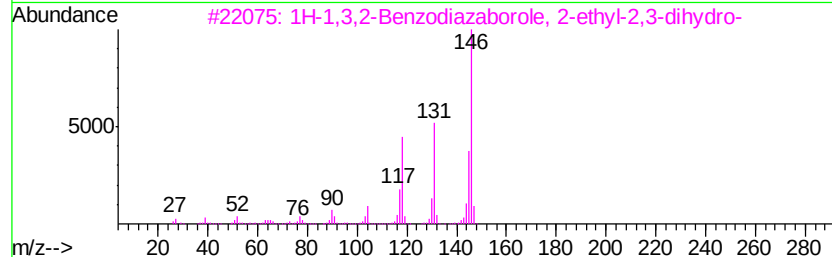
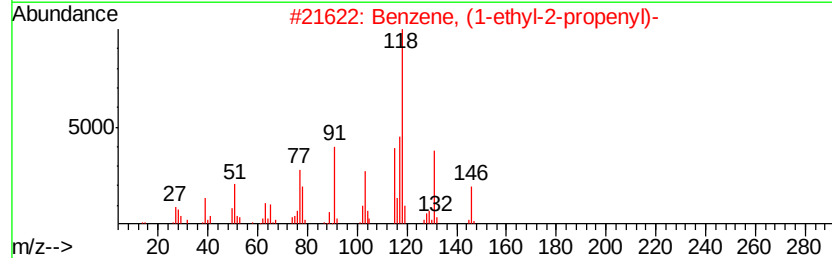
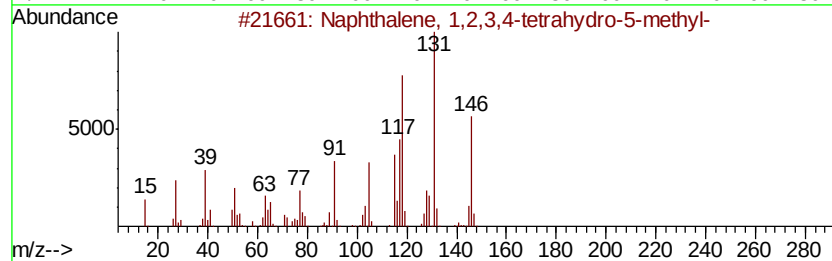
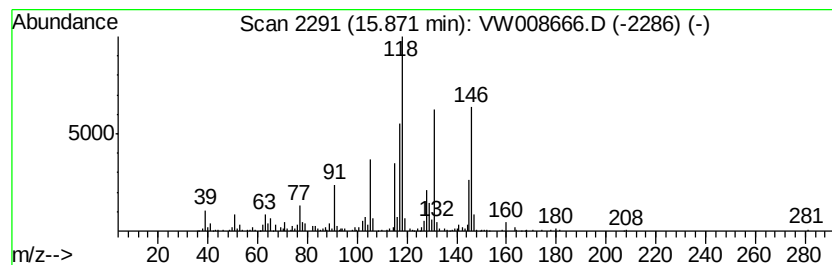
Instrument :
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ClientSampleId :
OR-01-021319-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W021319S.M
Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetra... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.87	10.94 ug/l	324407	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 81
2		Benzene, (1-ethyl-2-propenyl)-	146 C11H14	019947-22-9 74
3		1H-1,3,2-Benzodiazaborole, 2-eth...	146 C8H11BN2	074663-81-3 64
4		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 62
5		2,3,4,5,6,7-Hexahydro-1H-cyclope...	146 C11H14	1000189-31-0 49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW021419\
Data File : VW008666.D
Acq On : 14 Feb 2019 15:13
Operator : SY/VA
Sample : K1349-05
Misc : 7.37G/5ML/MSVOA W/SOIL
ALS Vial : 9 Sample Multiplier: 1

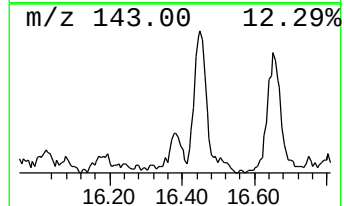
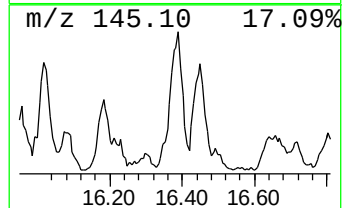
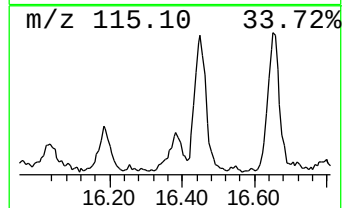
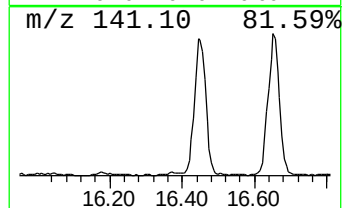
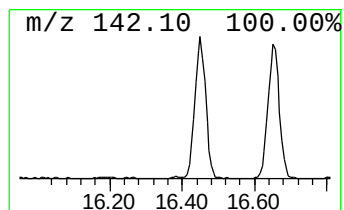
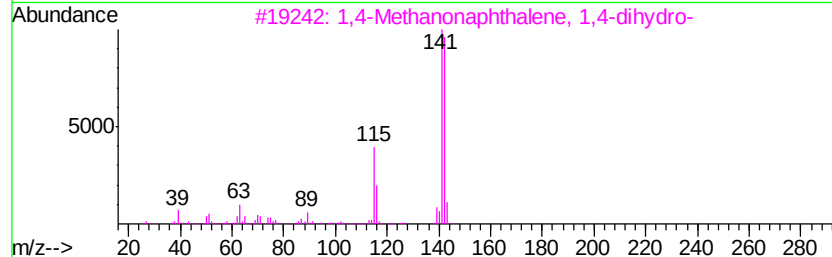
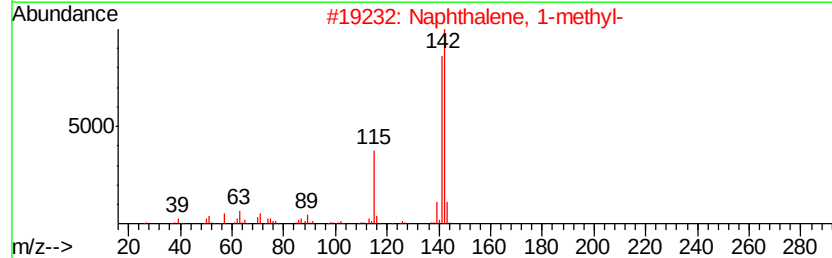
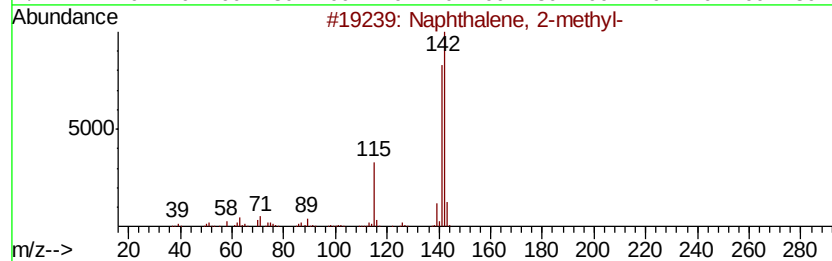
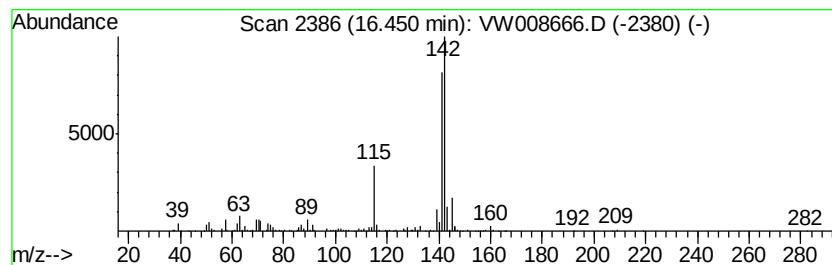
Instrument :
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ClientSampleId :
OR-01-021319-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W021319S.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.45	9.22 ug/l	273263	1,4-Dichlorobenzene-d4	13.56	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	93
4	Benzocycloheptatriene	142	C11H10	000264-09-5	93
5	Naphthalene, 1-[(2-propenyloxy)m...	198	C14H14O	074685-39-5	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW021419\
Data File : VW008666.D
Acq On : 14 Feb 2019 15:13
Operator : SY/VA
Sample : K1349-05
Misc : 7.37G/5ML/MSVOA_W/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
OR-01-021319-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W021319S.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, 1-propenyl-	13.76	9.0	ug/l	267087	4	13.56	1482350	50.0
1-Phenyl-1-butene	14.21	8.0	ug/l	235738	4	13.56	1482350	50.0
Benzene, 2-etheny...	14.69	8.3	ug/l	245896	4	13.56	1482350	50.0
Benzene, 1-etheny...	14.81	20.5	ug/l	607734	4	13.56	1482350	50.0
Naphthalene, 1,2,...	14.96	9.4	ug/l	277825	4	13.56	1482350	50.0
1H-Indene, 2,3-di...	15.18	7.8	ug/l	230501	4	13.56	1482350	50.0
Naphthalene, 1,2,...	15.87	10.9	ug/l	324407	4	13.56	1482350	50.0
Naphthalene, 1-me...	16.45	9.2	ug/l	273263	4	13.56	1482350	50.0