

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW031319\
 Data File : VW009162.D
 Acq On : 13 Mar 2019 18:16
 Operator : SY/VA
 Sample : K1693-05
 Misc : 5.81G/5ML/MSVOA W/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 OR-02-031219-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\82W030719S.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	2.629	115	119	127	rBV3	9978	19788	1.34%	0.076%
2	4.116	355	363	374	rBV3	7108	21003	1.42%	0.080%
3	4.915	485	494	509	rBV4	16173	53673	3.62%	0.206%
4	7.146	849	860	873	rBV	26358	76313	5.15%	0.292%
5	7.884	972	981	986	rBV	184925	442881	29.88%	1.696%
6	7.945	986	991	1005	rVB	343107	751847	50.73%	2.879%
7	8.305	1039	1050	1061	rBV	200932	446204	30.11%	1.708%
8	8.841	1129	1138	1148	rBV	494286	994275	67.09%	3.807%
9	10.323	1372	1381	1387	rBV	761777	1331643	89.85%	5.099%
10	10.390	1387	1392	1402	rVB	163762	294526	19.87%	1.128%
11	10.707	1436	1444	1449	rBV	32636	64199	4.33%	0.246%
12	11.633	1588	1596	1604	rBV	651115	1100116	74.23%	4.212%
13	11.731	1605	1612	1621	rVB2	37182	63476	4.28%	0.243%
14	11.835	1621	1629	1639	rBV	149992	293781	19.82%	1.125%
15	12.164	1674	1683	1690	rBV	127407	226814	15.30%	0.868%
16	12.463	1727	1732	1737	rBV	31846	54283	3.66%	0.208%
17	12.621	1751	1758	1766	rBV	514342	864440	58.33%	3.310%
18	12.804	1782	1788	1793	rBV	105813	162784	10.98%	0.623%
19	12.865	1793	1798	1802	rBV	397667	666332	44.96%	2.551%
20	12.901	1802	1804	1807	rVV	185054	248409	16.76%	0.951%
21	12.944	1807	1811	1817	rVB2	247778	414480	27.97%	1.587%
22	13.042	1823	1827	1831	rBV	23722	35479	2.39%	0.136%
23	13.103	1831	1837	1844	rBV	274642	442905	29.88%	1.696%
24	13.170	1844	1848	1854	rVB5	25742	47288	3.19%	0.181%
25	13.255	1854	1862	1868	rBV	886280	1394927	94.12%	5.341%
26	13.383	1874	1883	1888	rBV2	62138	110592	7.46%	0.423%
27	13.444	1888	1893	1898	rBV	116691	189957	12.82%	0.727%
28	13.493	1898	1901	1906	rBV3	40689	63501	4.28%	0.243%
29	13.560	1906	1912	1915	rBV	571199	1002210	67.62%	3.837%
30	13.724	1933	1939	1940	rBV	99635	147410	9.95%	0.564%
31	13.755	1940	1944	1948	rVV2	569097	1010715	68.20%	3.870%
32	13.792	1948	1950	1960	rVB3	309366	638830	43.10%	2.446%
33	13.877	1960	1964	1970	rBV2	228900	352182	23.76%	1.348%
34	13.950	1971	1976	1981	rBV	169048	250033	16.87%	0.957%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	14.011	1981	1986	1989	rBV	197953	329878	22.26%	1.263%
36	14.041	1989	1991	1996	rVV	215238	268685	18.13%	1.029%
37	14.102	1996	2001	2005	rVB	296934	464361	31.33%	1.778%
38	14.163	2005	2011	2014	rBV	85749	137926	9.31%	0.528%
39	14.212	2014	2019	2024	rVB2	378600	614256	41.45%	2.352%
40	14.291	2024	2032	2036	rBV5	43945	116022	7.83%	0.444%
41	14.340	2036	2040	2044	rBV2	102944	158009	10.66%	0.605%
42	14.389	2044	2048	2050	rBV3	56590	76451	5.16%	0.293%
43	14.426	2050	2054	2057	rVV	127567	189555	12.79%	0.726%
44	14.462	2057	2060	2064	rVB	234114	291785	19.69%	1.117%
45	14.505	2064	2067	2070	rVB	117705	134010	9.04%	0.513%
46	14.548	2070	2074	2080	rVB5	79739	131305	8.86%	0.503%
47	14.645	2080	2090	2093	rBV3	126019	249399	16.83%	0.955%
48	14.694	2093	2098	2101	rBV	357860	604529	40.79%	2.315%
49	14.730	2101	2104	2109	rVB2	269441	419498	28.31%	1.606%
50	14.810	2109	2117	2123	rBV3	685052	1482041	100.00%	5.675%
51	14.962	2136	2142	2148	rVB	496805	785047	52.97%	3.006%
52	15.035	2149	2154	2155	rBV2	66128	95937	6.47%	0.367%
53	15.072	2155	2160	2163	rVV2	210030	417522	28.17%	1.599%
54	15.108	2164	2166	2171	rVV	119332	179716	12.13%	0.688%
55	15.188	2171	2179	2186	rVB3	208841	511573	34.52%	1.959%
56	15.334	2198	2203	2205	rBV4	72400	128881	8.70%	0.493%
57	15.371	2205	2209	2213	rVB	99226	157108	10.60%	0.602%
58	15.425	2213	2218	2223	rBV2	182280	305341	20.60%	1.169%
59	15.480	2223	2227	2229	rBV3	105224	174289	11.76%	0.667%
60	15.560	2237	2240	2249	rVB3	193566	339823	22.93%	1.301%
61	15.663	2249	2257	2263	rBV2	168550	369869	24.96%	1.416%
62	15.736	2265	2269	2274	rVB4	32307	55709	3.76%	0.213%
63	15.828	2278	2284	2285	rVV3	102978	165735	11.18%	0.635%
64	15.870	2286	2291	2301	rVB	393795	756345	51.03%	2.896%
65	15.956	2301	2305	2310	rBV6	37454	71394	4.82%	0.273%
66	16.035	2310	2318	2324	rBV3	81240	203466	13.73%	0.779%
67	16.187	2334	2343	2349	rBV2	188374	451610	30.47%	1.729%
68	16.297	2358	2361	2365	rVB5	16667	22628	1.53%	0.087%
69	16.383	2367	2375	2380	rVV3	134223	299491	20.21%	1.147%
70	16.450	2380	2386	2391	rVV	135401	296139	19.98%	1.134%
71	16.504	2391	2395	2400	rVB	92444	161983	10.93%	0.620%

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Integration Parameters: RTEINT.P

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Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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72	16.651	2412	2419	2426	rBV4	80076	222700	15.03%	0.853%
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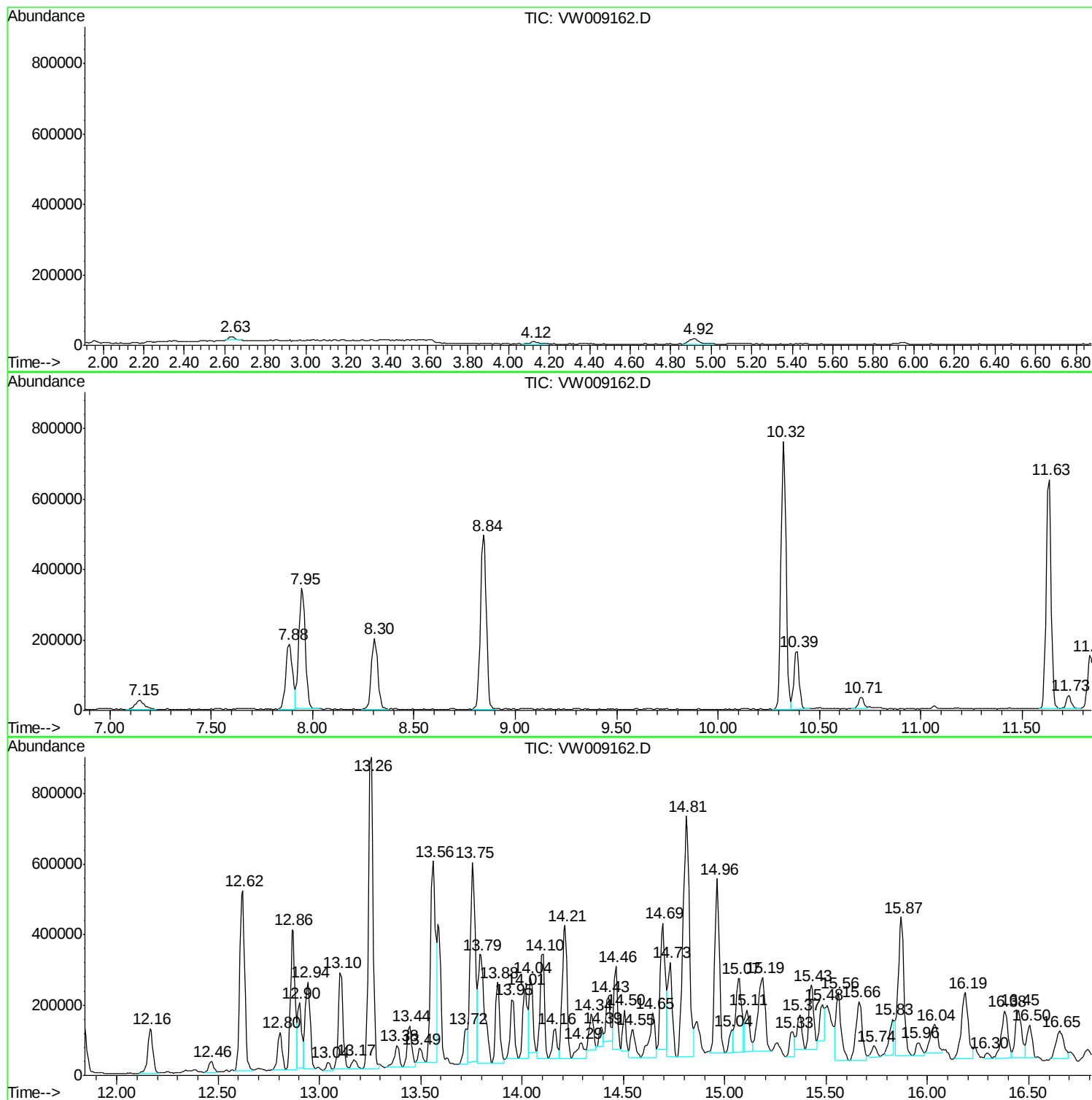
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Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW031319\
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Acq On : 13 Mar 2019 18:16
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ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
OR-02-031219-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W030719S.M
Quant Title : SW846 8260

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



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Misc : 5.81G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
OR-02-031219-A

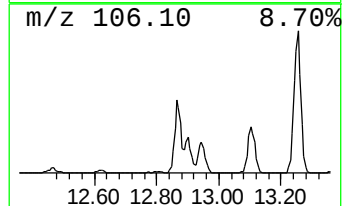
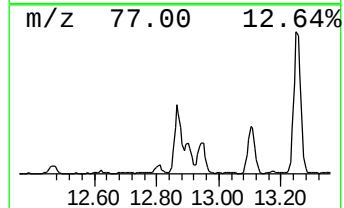
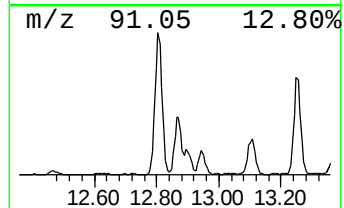
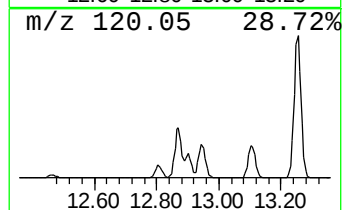
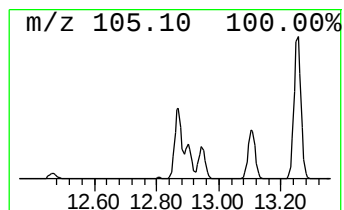
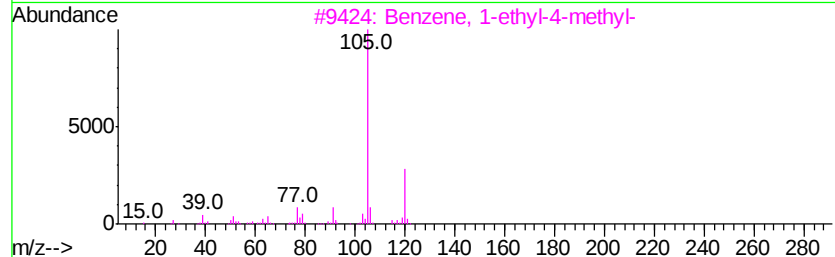
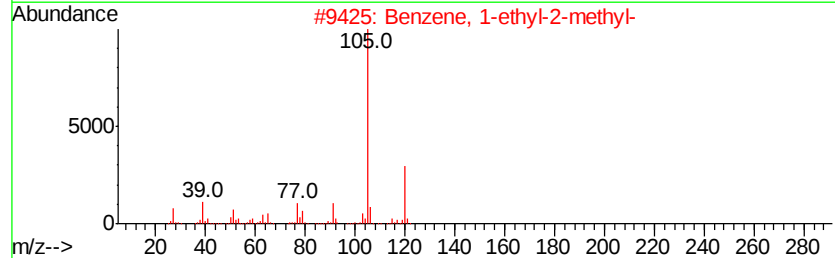
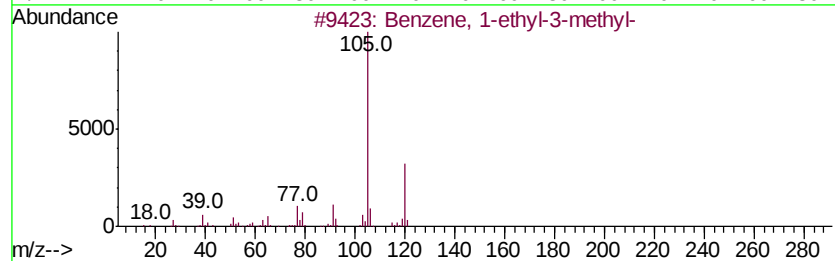
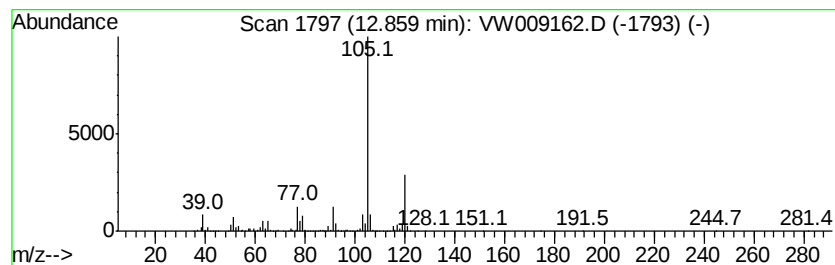
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W030719S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.86	33.24 ug/l	666332	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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Misc : 5.81G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
OR-02-031219-A

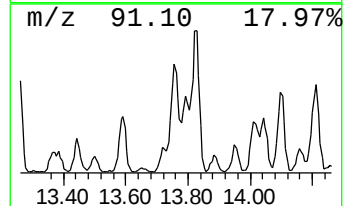
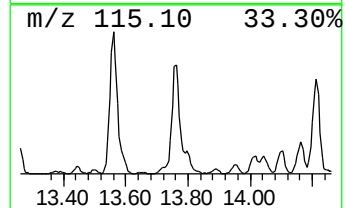
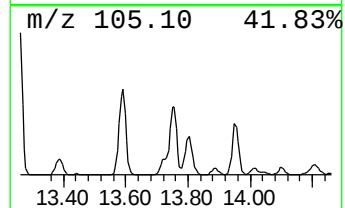
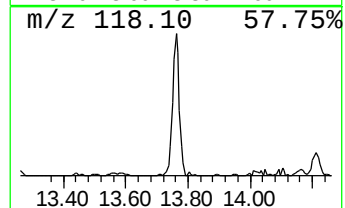
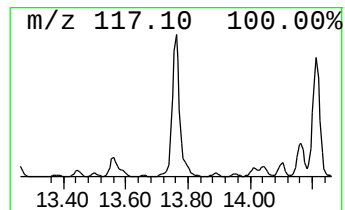
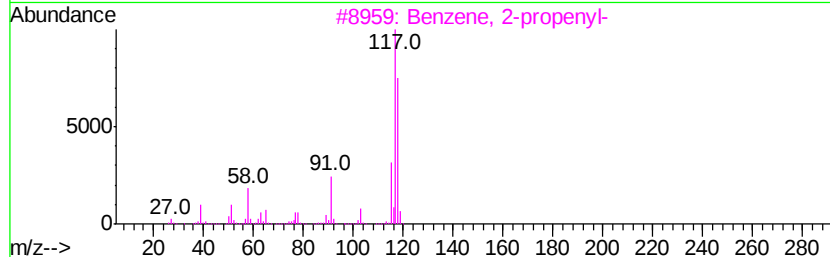
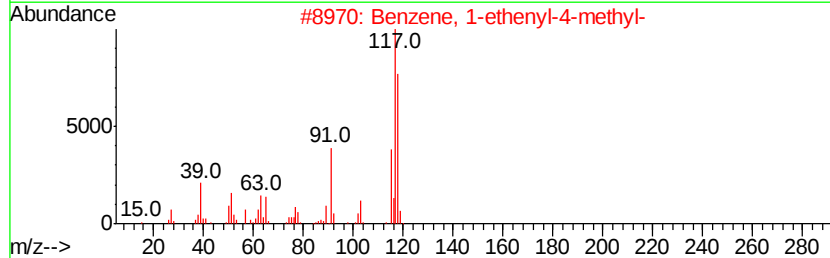
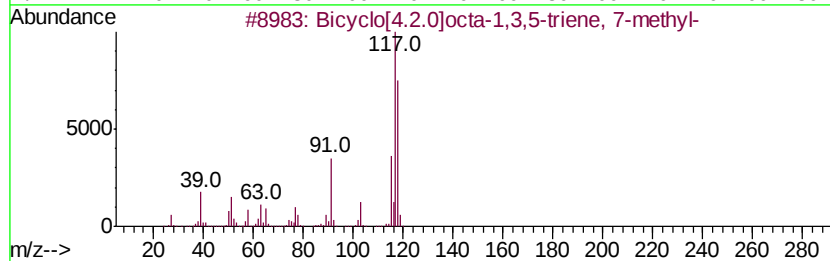
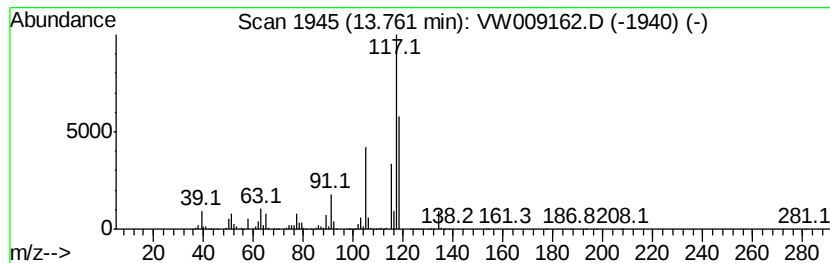
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W030719S.M
Quant Title : SW846 8260

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

Peak Number 2 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	50.42 ug/l	1010720	1,4-Dichlorobenzene-d4	13.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	055337-80-9	92
2		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	86
3		Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
4		Indane	118	C9H10	000496-11-7	70
5		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	70



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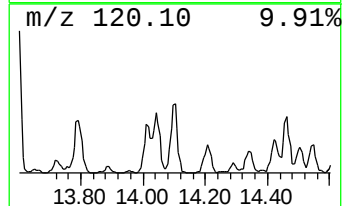
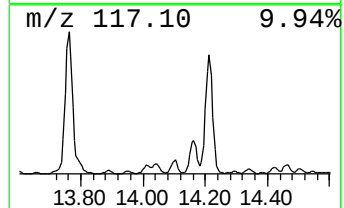
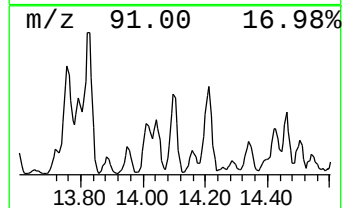
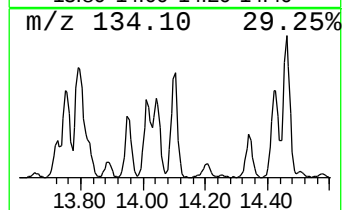
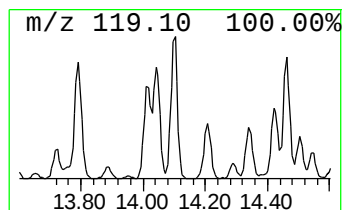
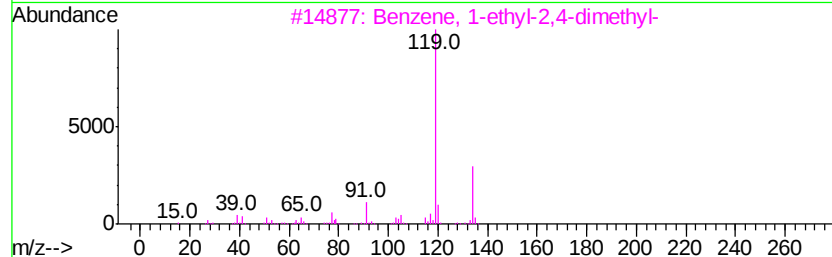
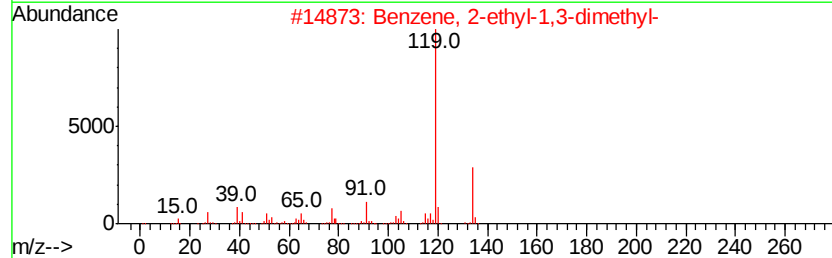
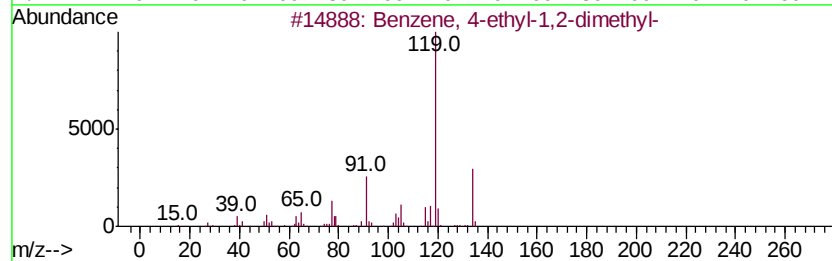
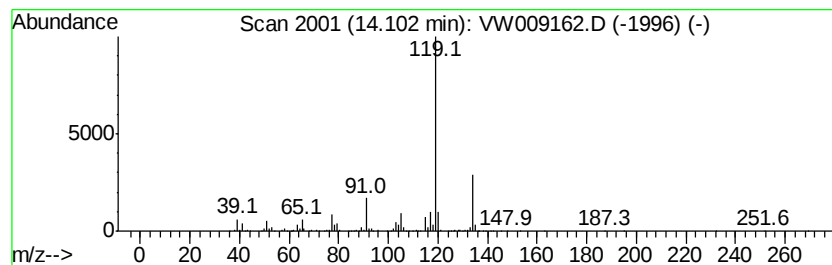
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.10	23.17 ug/l	464361	1,4-Dichlorobenzene-d4	13.56

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
5			o-Cymene	134	C10H14	000527-84-4	94



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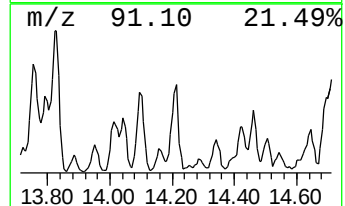
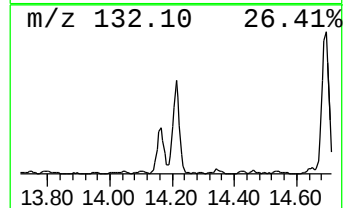
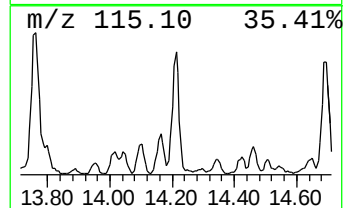
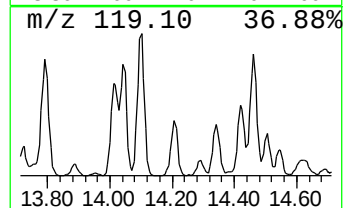
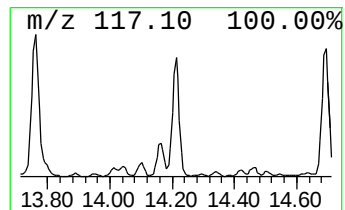
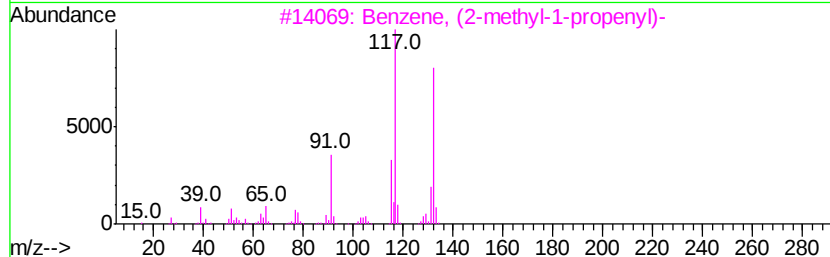
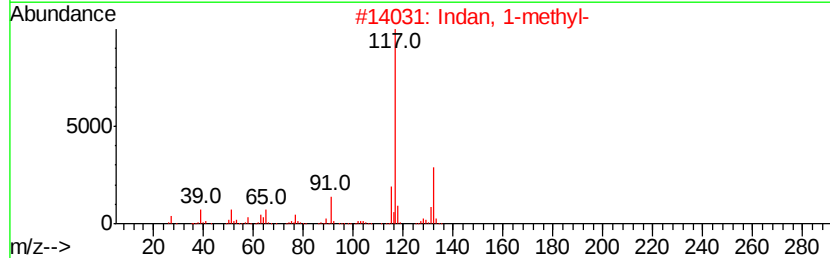
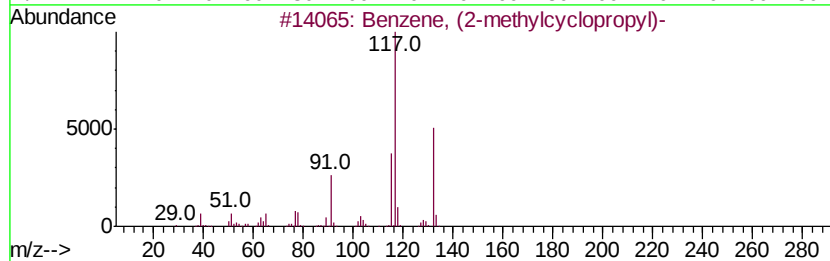
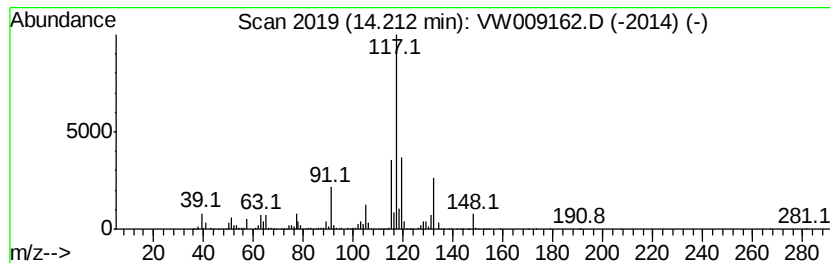
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, (2-methylcycloprop... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	30.65 ug/l	614256	1,4-Dichlorobenzene-d4	13.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	76
2		Indan, 1-methyl-	132	C10H12	000767-58-8	76
3		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	74
4		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	72
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	68



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW031319\
Data File : VW009162.D
Acq On : 13 Mar 2019 18:16
Operator : SY/VA
Sample : K1693-05
Misc : 5.81G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
OR-02-031219-A

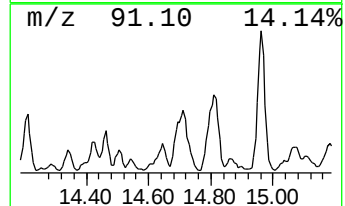
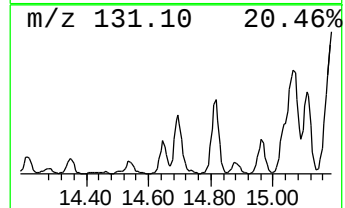
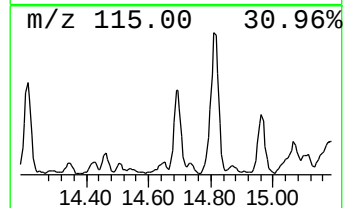
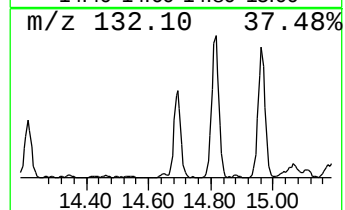
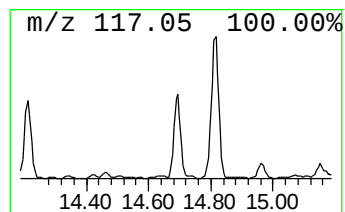
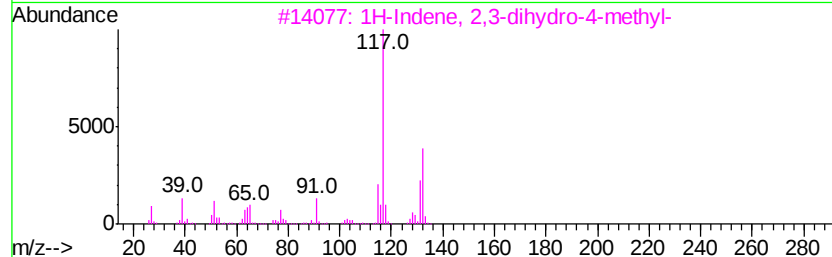
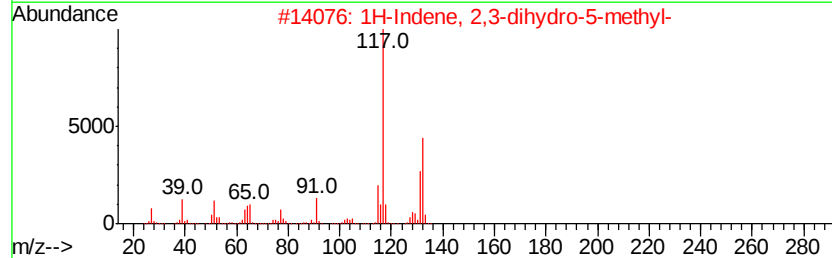
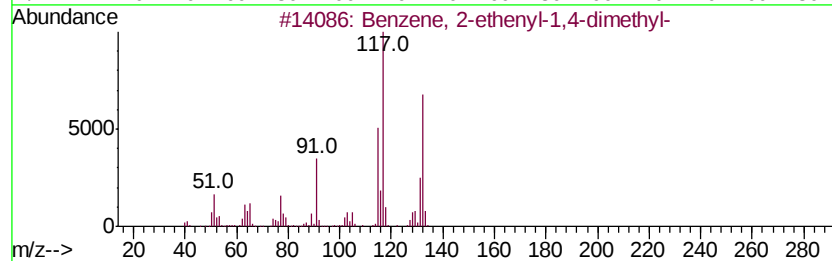
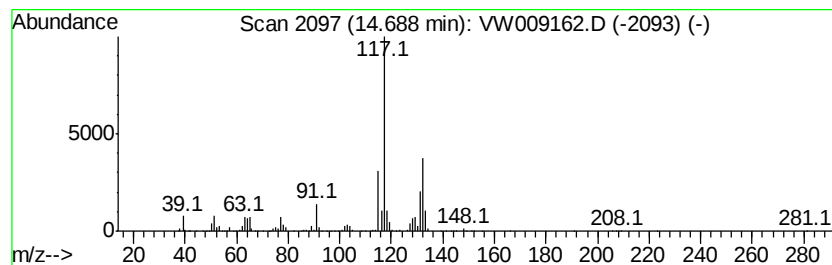
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	30.16 ug/l	604529	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	92
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW031319\
Data File : VW009162.D
Acq On : 13 Mar 2019 18:16
Operator : SY/VA
Sample : K1693-05
Misc : 5.81G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
OR-02-031219-A

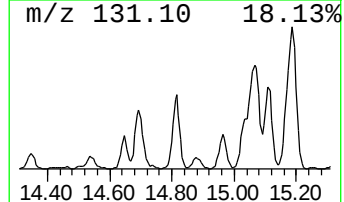
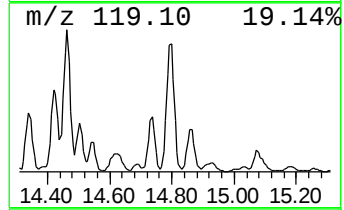
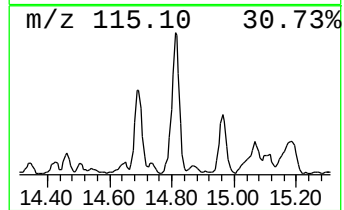
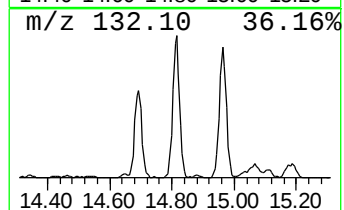
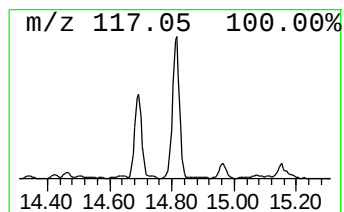
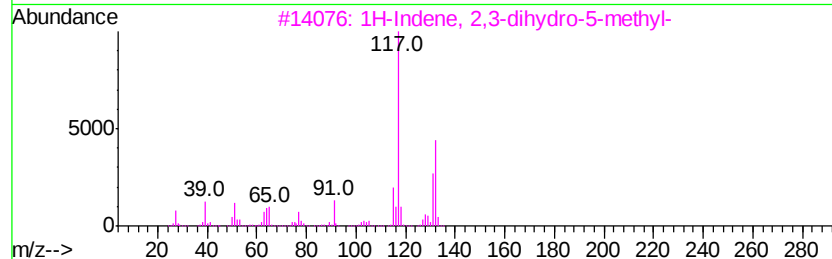
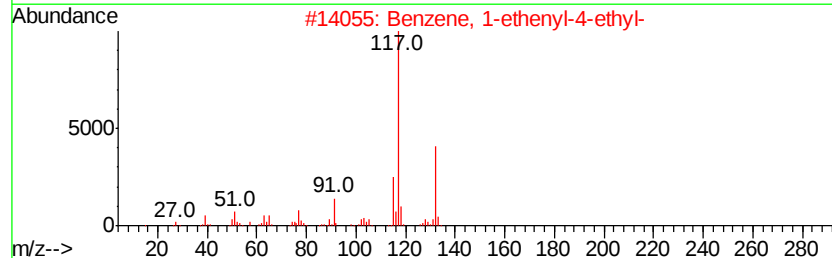
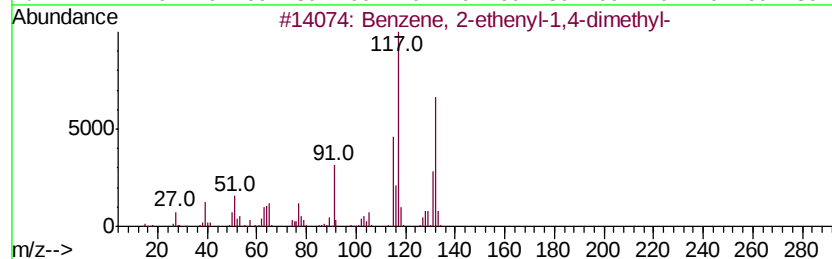
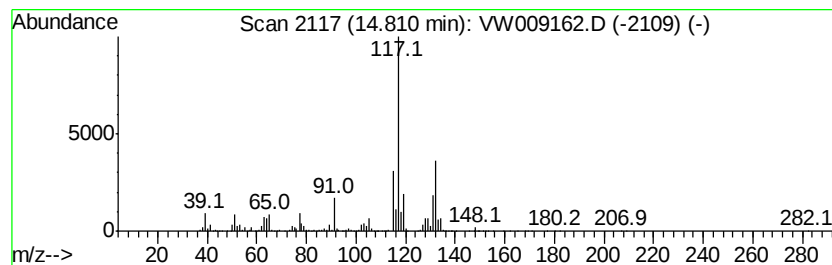
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.81	73.94 ug/l	1482040	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	94
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW031319\
Data File : VW009162.D
Acq On : 13 Mar 2019 18:16
Operator : SY/VA
Sample : K1693-05
Misc : 5.81G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
OR-02-031219-A

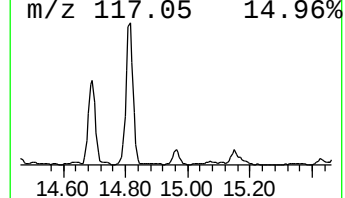
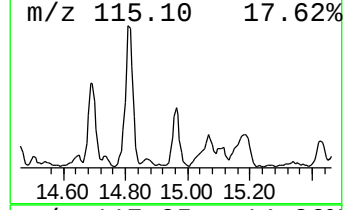
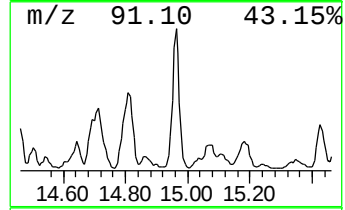
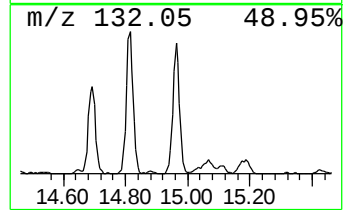
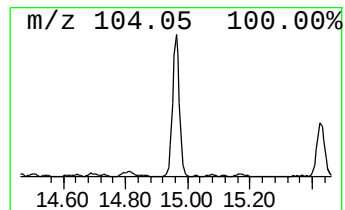
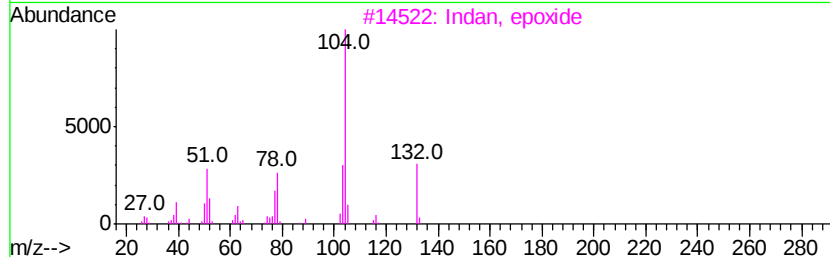
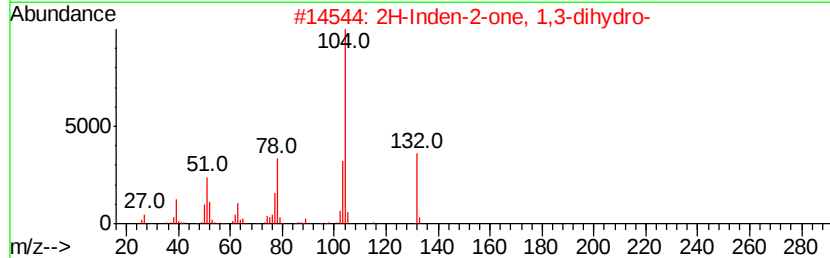
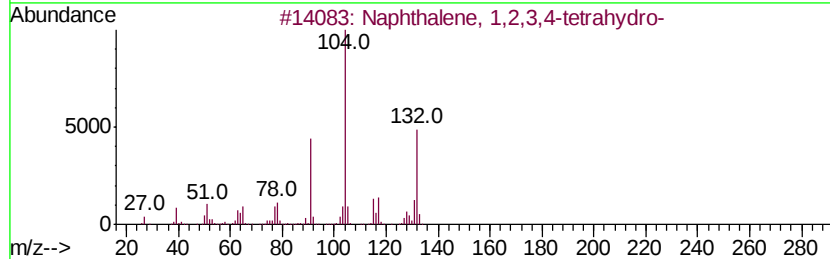
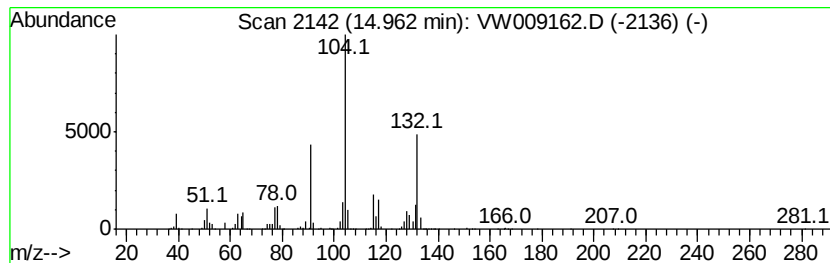
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	39.17 ug/l	785047	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	97
2		2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	58
3		Indan, epoxide	132	C9H8O	000768-22-9	52
4		2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	43
5		Benzenemethanol, 2-methyl-	122	C8H10O	000089-95-2	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW031319\
Data File : VW009162.D
Acq On : 13 Mar 2019 18:16
Operator : SY/VA
Sample : K1693-05
Misc : 5.81G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

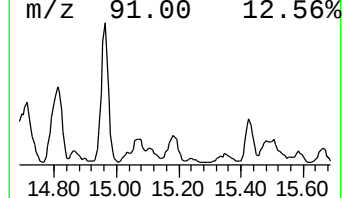
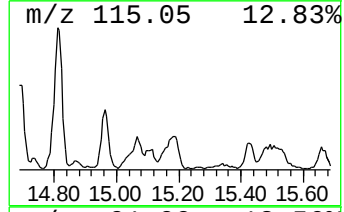
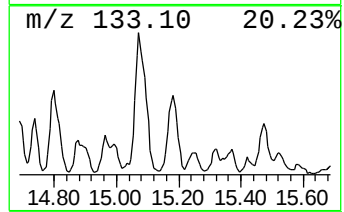
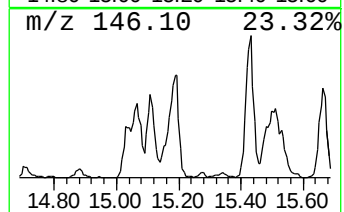
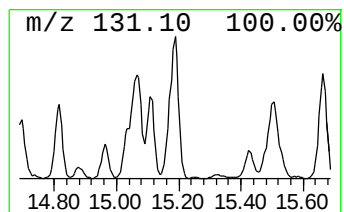
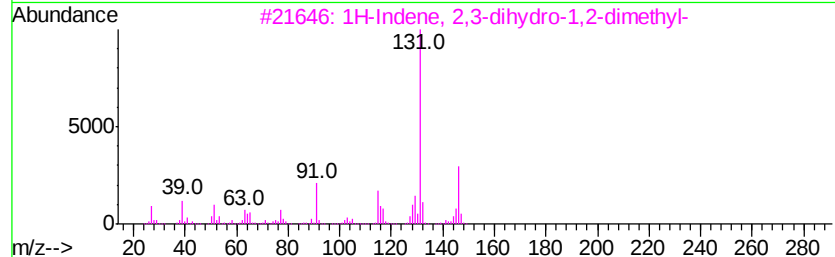
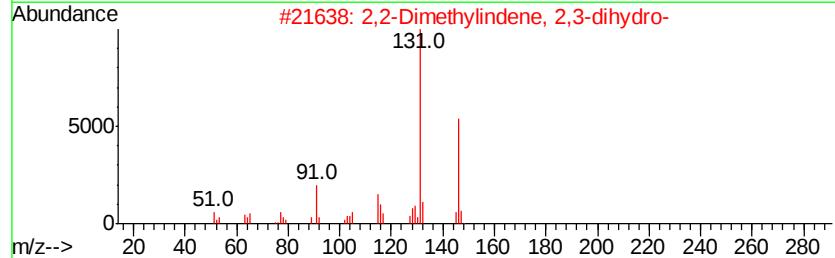
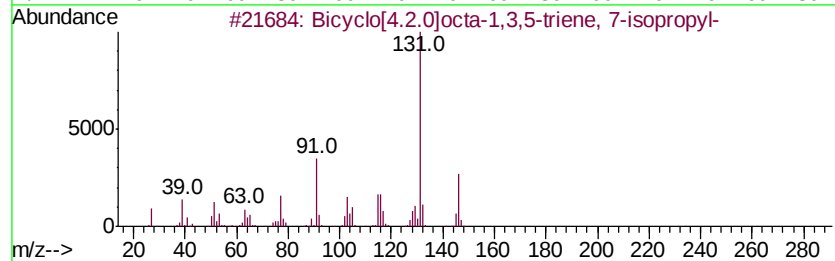
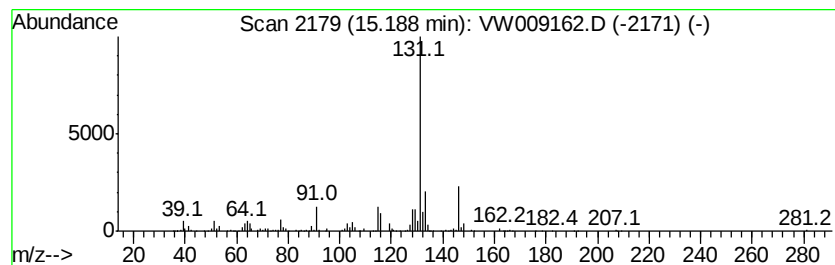
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ClientSampleId :
OR-02-031219-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W030719S.M
Quant Title : SW846 8260

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

Peak Number 8 2,2-Dimethylindene, 2,3-dih... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	25.52 ug/l	511573	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Bicyclo[4.2.0]octa-1,3,5-triene,...	146 C11H14	027087-54-3 90
2		2,2-Dimethylindene, 2,3-dihydro-	146 C11H14	020836-11-7 81
3		1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8 81
4		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 81
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW031319\
Data File : VW009162.D
Acq On : 13 Mar 2019 18:16
Operator : SY/VA
Sample : K1693-05
Misc : 5.81G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

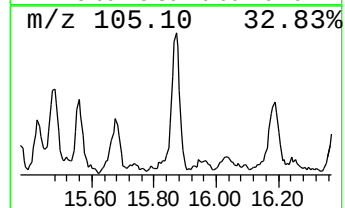
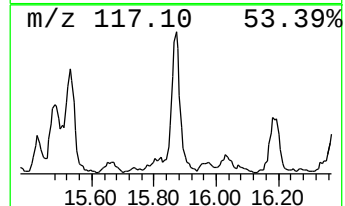
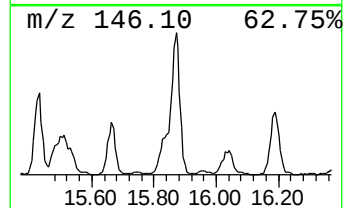
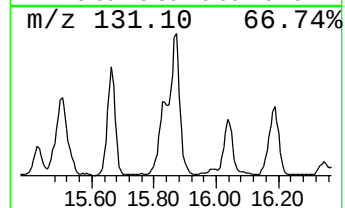
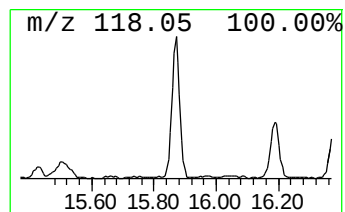
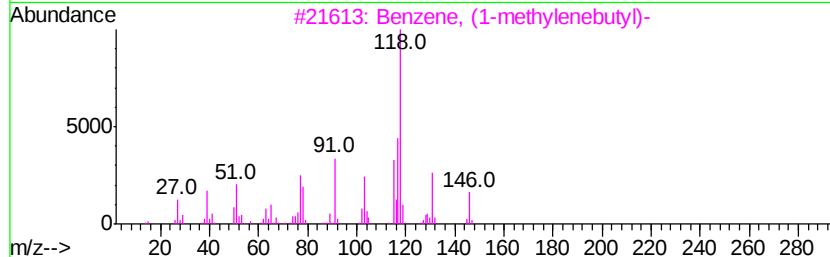
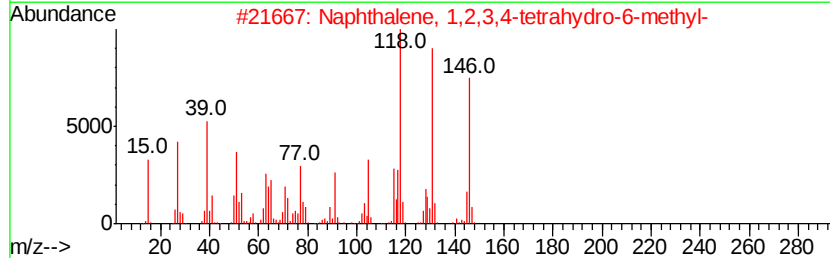
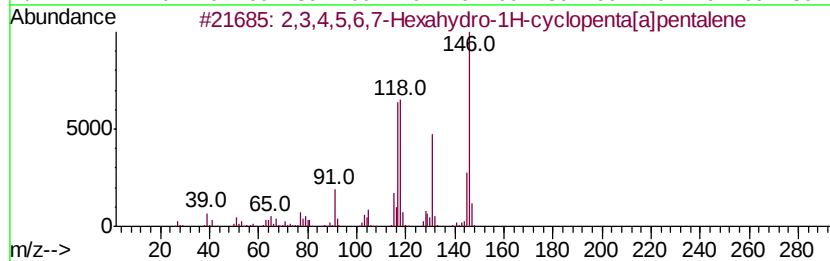
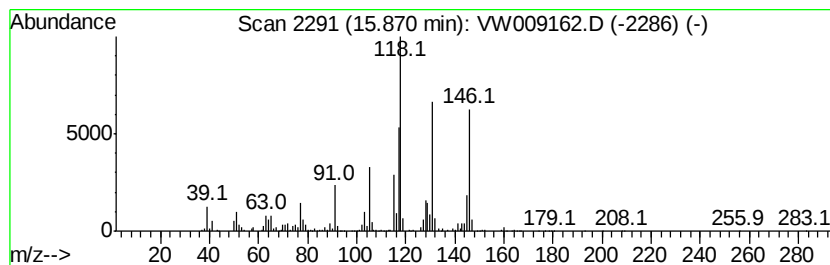
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ClientSampleId :
OR-02-031219-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W030719S.M
Quant Title : SW846 8260

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

Peak Number 9 2,3,4,5,6,7-Hexahydro-1H-cy... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.87	37.73 ug/l	756345	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2,3,4,5,6,7-Hexahydro-1H-cyclope...	146 C11H14	1000189-31-0	89
2	Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9	86
3	Benzene, (1-methylenebutyl)-	146 C11H14	005676-32-4	80
4	Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5	70
5	1H-1,3,2-Benzodiazaborole, 2-eth...	146 C8H11BN2	074663-81-3	49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW031319\
Data File : VW009162.D
Acq On : 13 Mar 2019 18:16
Operator : SY/VA
Sample : K1693-05
Misc : 5.81G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
OR-02-031219-A

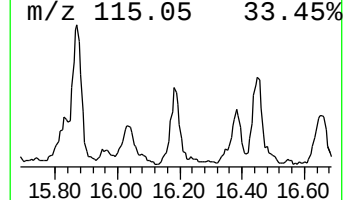
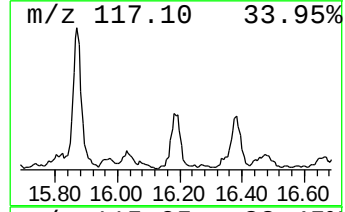
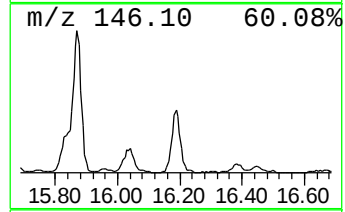
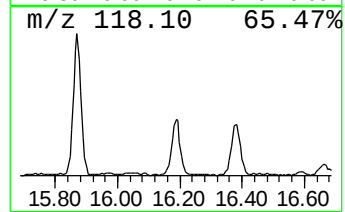
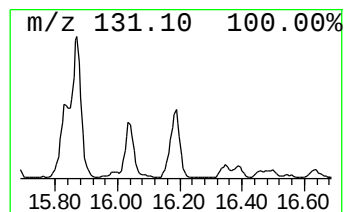
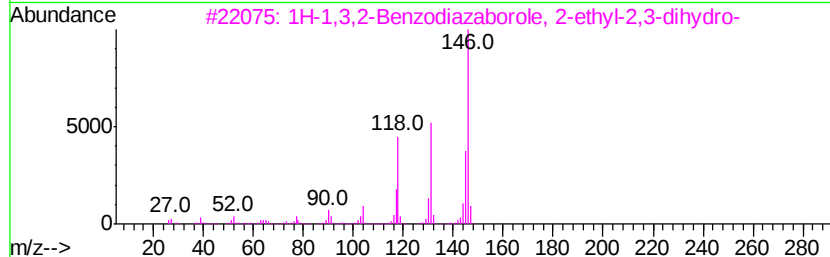
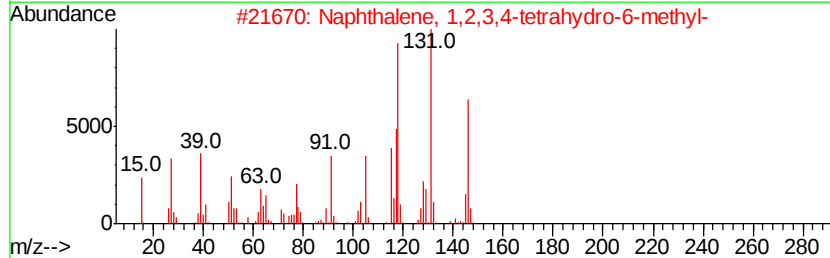
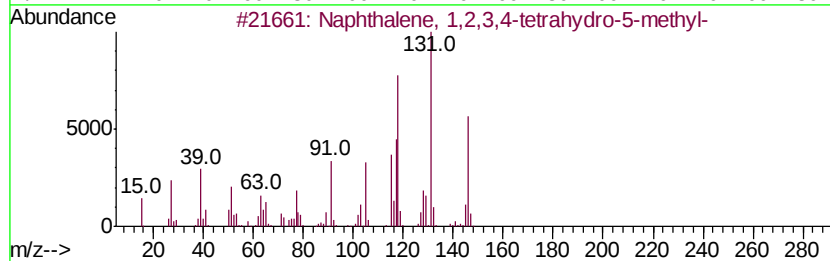
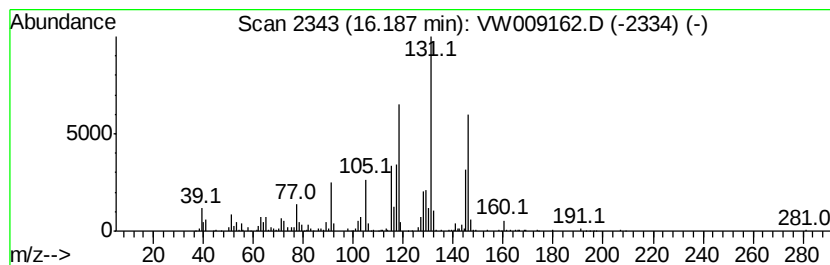
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.19	22.53 ug/l	451610	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	93
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	91
3		1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	90
4		2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	68
5		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW031319\
Data File : VW009162.D
Acq On : 13 Mar 2019 18:16
Operator : SY/VA
Sample : K1693-05
Misc : 5.81G/5ML/MSVOA_W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
OR-02-031219-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W030719S.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-ethyl-...	12.86	33.2	ug/l	666332	4	13.56	1002210	50.0
Bicyclo[4.2.0]oct...	13.76	50.4	ug/l	1010720	4	13.56	1002210	50.0
Benzene, 4-ethyl-...	14.10	23.2	ug/l	464361	4	13.56	1002210	50.0
Benzene, (2-methy...	14.21	30.6	ug/l	614256	4	13.56	1002210	50.0
Benzene, 2-etheny...	14.69	30.2	ug/l	604529	4	13.56	1002210	50.0
Benzene, 1-etheny...	14.81	73.9	ug/l	1482040	4	13.56	1002210	50.0
Naphthalene, 1,2,...	14.96	39.2	ug/l	785047	4	13.56	1002210	50.0
2,2-Dimethylinden...	15.19	25.5	ug/l	511573	4	13.56	1002210	50.0
2,3,4,5,6,7-Hexah...	15.87	37.7	ug/l	756345	4	13.56	1002210	50.0
Naphthalene, 1,2,...	16.19	22.5	ug/l	451610	4	13.56	1002210	50.0