

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW101718\
 Data File : VW006139.D
 Acq On : 17 Oct 2018 20:45
 Operator : VA/AP
 Sample : J5515-18
 Misc : 5.96G/5ML/MSVOA W/SOIL
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 AOC1-01A

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\82W101518S.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	4.928	484	496	508	rBV7	5423	23253	1.63%	0.182%
2	6.994	825	835	843	rVB3	9511	27237	1.91%	0.213%
3	7.147	850	860	872	rBV2	6752	20232	1.42%	0.158%
4	7.891	971	982	986	rBV2	19221	50909	3.57%	0.398%
5	7.952	986	992	999	rVV6	34250	94796	6.65%	0.741%
6	8.037	999	1006	1015	rVB2	16911	41855	2.94%	0.327%
7	8.159	1017	1026	1033	rBV2	22309	51336	3.60%	0.401%
8	8.317	1043	1052	1060	rVV2	22384	53562	3.76%	0.418%
9	8.488	1063	1080	1090	rBV3	30347	122574	8.60%	0.958%
10	8.579	1090	1095	1103	rVV2	10126	22885	1.61%	0.179%
11	8.848	1128	1139	1151	rBV2	52871	126506	8.88%	0.988%
12	9.335	1205	1219	1225	rBV4	51791	145804	10.23%	1.139%
13	9.390	1225	1228	1236	rVB3	12404	21285	1.49%	0.166%
14	9.518	1244	1249	1256	rVB2	12351	22941	1.61%	0.179%
15	9.616	1256	1265	1271	rVB3	7087	17109	1.20%	0.134%
16	9.762	1283	1289	1299	rVB	42230	84146	5.91%	0.657%
17	9.896	1301	1311	1317	rBV	58032	127667	8.96%	0.997%
18	9.963	1317	1322	1324	rVV5	9233	17569	1.23%	0.137%
19	9.994	1324	1327	1335	rVB6	11436	24193	1.70%	0.189%
20	10.098	1335	1344	1356	rVB3	21850	60011	4.21%	0.469%
21	10.250	1365	1369	1375	rVB2	13222	24766	1.74%	0.193%
22	10.329	1375	1382	1394	rBV	70855	151569	10.64%	1.184%
23	10.518	1405	1413	1418	rVB9	12736	38021	2.67%	0.297%
24	10.671	1433	1438	1442	rBV	11729	21396	1.50%	0.167%
25	10.750	1446	1451	1461	rVB5	27525	68575	4.81%	0.536%
26	10.939	1474	1482	1488	rBV8	5549	14987	1.05%	0.117%
27	11.024	1488	1496	1497	rBV6	7676	17284	1.21%	0.135%
28	11.183	1519	1522	1527	rVV3	10343	16250	1.14%	0.127%
29	11.445	1563	1565	1572	rVB3	14205	23114	1.62%	0.181%
30	11.518	1572	1577	1581	rBV2	11332	20480	1.44%	0.160%
31	11.634	1590	1596	1602	rBV	67051	117008	8.21%	0.914%
32	11.731	1608	1612	1617	rVB3	10456	15706	1.10%	0.123%
33	11.847	1622	1631	1634	rBV2	10788	24046	1.69%	0.188%
34	12.140	1674	1679	1682	rBV5	7965	16125	1.13%	0.126%

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35	12.341	1702	1712	1716	rBV10	9927	29997	2.11%	0.234%
36	12.469	1728	1733	1739	rBV	43909	75420	5.29%	0.589%
37	12.621	1752	1758	1767	rVB2	66513	125881	8.84%	0.983%
38	12.707	1767	1772	1777	rBV9	6917	14394	1.01%	0.112%
39	12.810	1783	1789	1795	rVB	93050	142496	10.00%	1.113%
40	13.109	1831	1838	1844	rBV	74147	122792	8.62%	0.959%
41	13.310	1867	1871	1874	rBV4	7932	14949	1.05%	0.117%
42	13.384	1874	1883	1890	rVV4	29071	86351	6.06%	0.675%
43	13.505	1898	1903	1908	rBV8	9114	20132	1.41%	0.157%
44	13.566	1908	1913	1923	rVB2	64145	142168	9.98%	1.111%
45	13.658	1925	1928	1933	rVB	13791	20808	1.46%	0.163%
46	13.725	1933	1939	1942	rBV	329262	519741	36.48%	4.061%
47	13.768	1942	1946	1951	rVV	914999	1424667	100.00%	11.131%
48	13.810	1951	1953	1961	rVB2	100006	146319	10.27%	1.143%
49	13.890	1961	1966	1973	rBV	70286	125980	8.84%	0.984%
50	13.957	1974	1977	1980	rBV	10771	14431	1.01%	0.113%
51	14.036	1983	1990	1995	rVB6	44070	109037	7.65%	0.852%
52	14.103	1995	2001	2007	rBV	891229	1362397	95.63%	10.644%
53	14.170	2007	2012	2015	rBV	74244	123723	8.68%	0.967%
54	14.213	2015	2019	2026	rVB2	446326	744890	52.29%	5.820%
55	14.286	2026	2031	2036	rBV5	22910	54542	3.83%	0.426%
56	14.353	2036	2042	2045	rBV5	23565	41848	2.94%	0.327%
57	14.396	2045	2049	2051	rBV3	23873	32084	2.25%	0.251%
58	14.426	2051	2054	2063	rVB2	41900	74133	5.20%	0.579%
59	14.511	2063	2068	2071	rBV	76406	115031	8.07%	0.899%
60	14.548	2071	2074	2077	rBV2	15156	22489	1.58%	0.176%
61	14.652	2087	2091	2094	rVB	21225	25146	1.77%	0.196%
62	14.700	2094	2099	2103	rBV	347389	559286	39.26%	4.370%
63	14.743	2103	2106	2111	rVB	73898	104014	7.30%	0.813%
64	14.816	2111	2118	2123	rBV	818786	1417463	99.49%	11.074%
65	14.865	2123	2126	2133	rVB2	115426	218947	15.37%	1.711%
66	14.969	2139	2143	2149	rVB5	32546	52702	3.70%	0.412%
67	15.072	2150	2160	2164	rVV3	184487	432297	30.34%	3.377%
68	15.115	2164	2167	2173	rVV2	133322	249940	17.54%	1.953%
69	15.188	2173	2179	2187	rVB3	226209	501426	35.20%	3.918%
70	15.340	2200	2204	2212	rVB7	40150	70036	4.92%	0.547%
71	15.432	2214	2219	2222	rBV	29839	55960	3.93%	0.437%

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Sampling : 1 Min Area: 3 % of largest Peak
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72	15.481	2223	2227	2231	rVV2	71189	147797	10.37%	1.155%
73	15.536	2231	2236	2249	rVB2	104593	276881	19.43%	2.163%
74	15.670	2250	2258	2267	rBV	162924	384591	27.00%	3.005%
75	15.840	2279	2286	2296	rVB2	62777	174735	12.26%	1.365%
76	15.962	2301	2306	2311	rBV2	52056	106543	7.48%	0.832%
77	16.188	2333	2343	2355	rBV7	52277	178452	12.53%	1.394%
78	16.395	2372	2377	2382	rVB6	33250	65534	4.60%	0.512%
79	16.657	2412	2420	2434	rVB	144338	345777	24.27%	2.702%

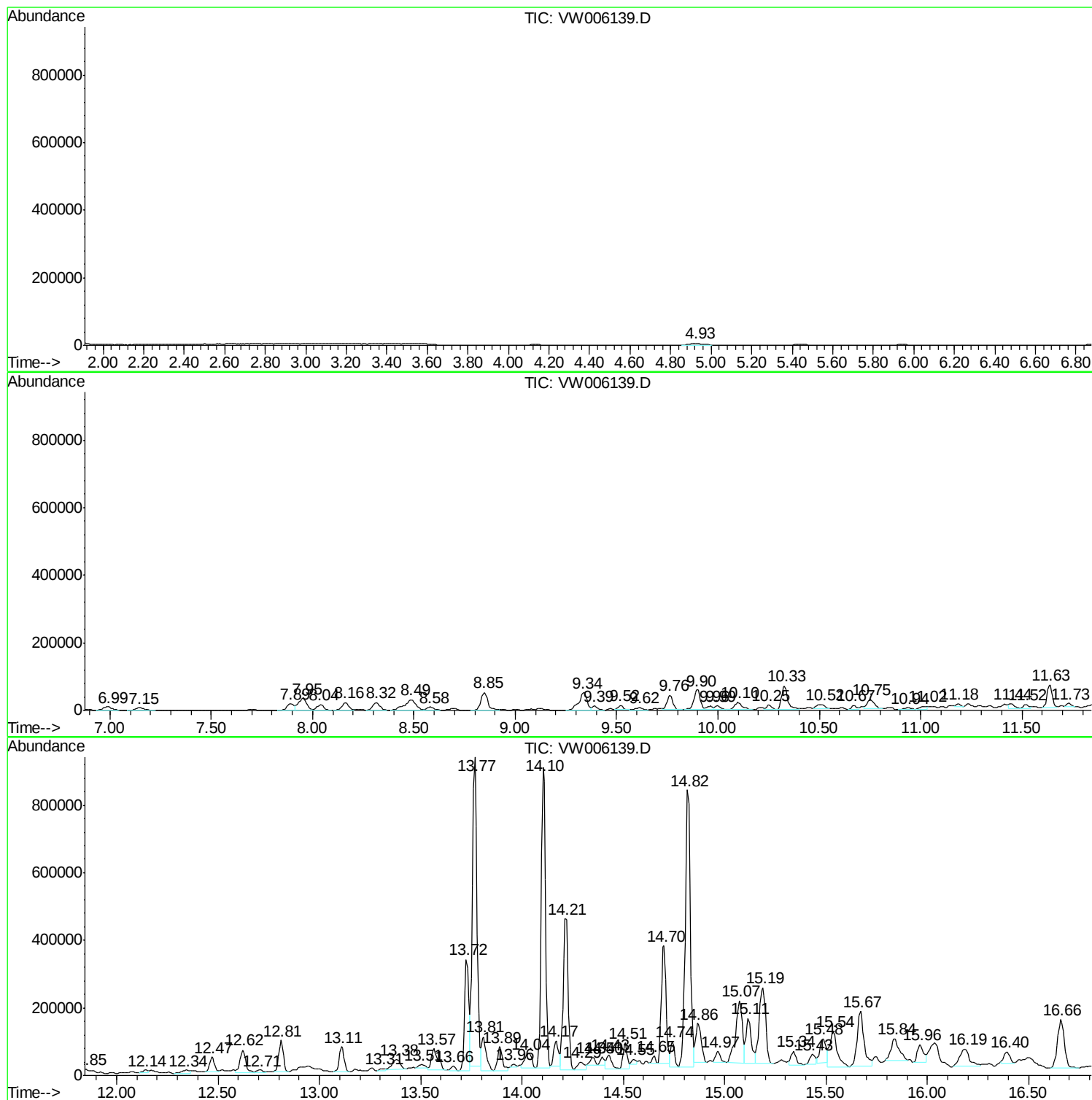
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Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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



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ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
AOC1-01A

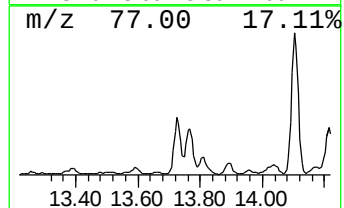
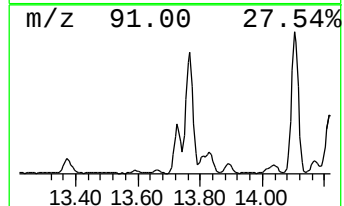
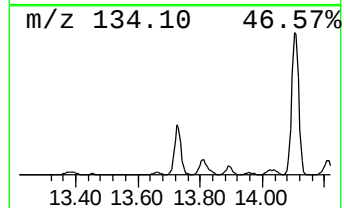
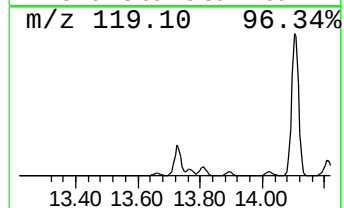
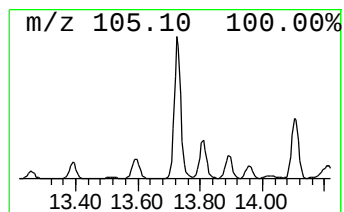
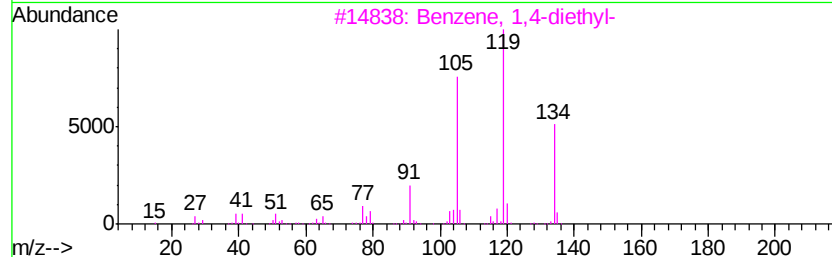
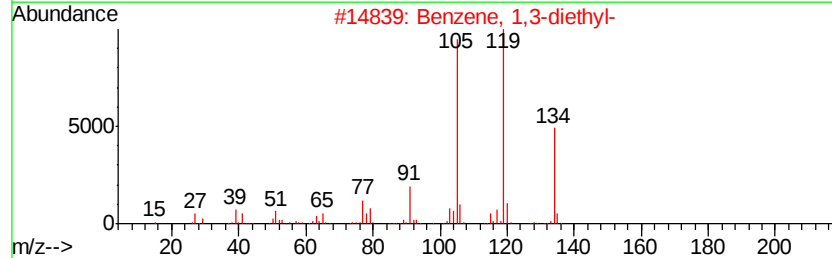
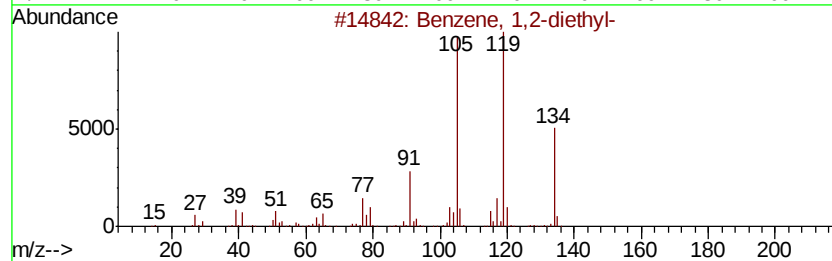
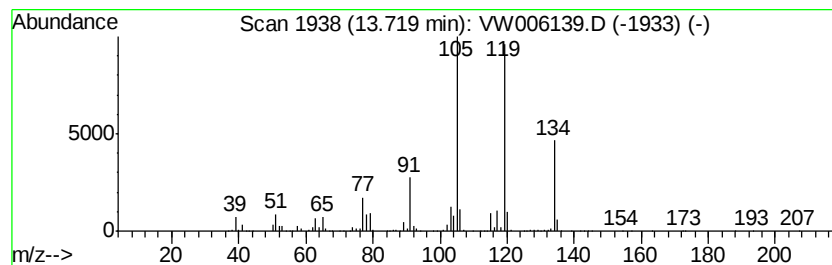
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1,2-diethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	182.79 ug/l	519741	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
2		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
3		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



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Instrument :
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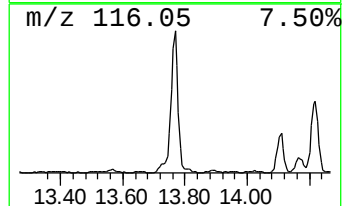
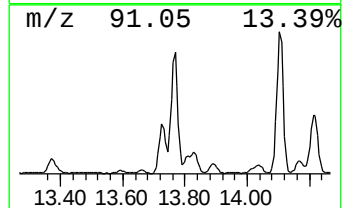
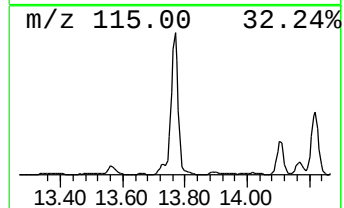
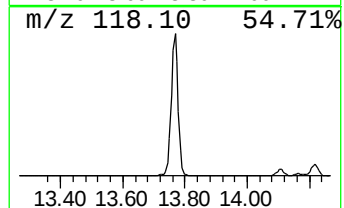
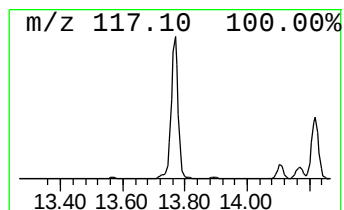
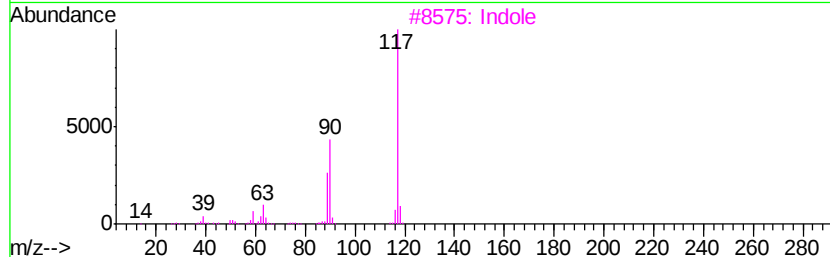
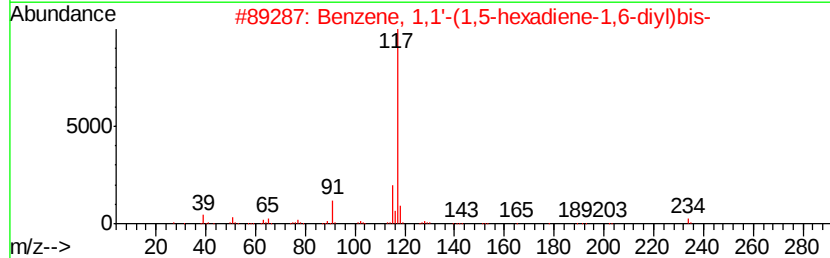
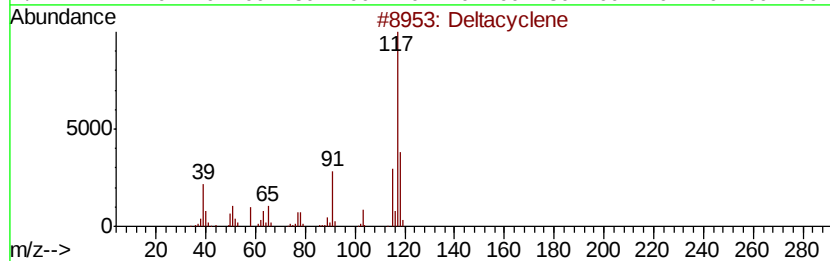
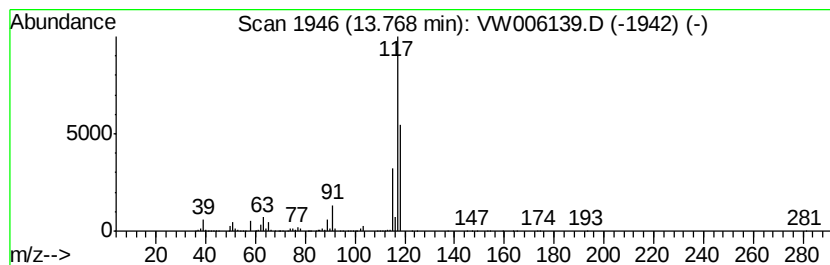
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Deltacyclene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	501.05 ug/l	1424670	1,4-Dichlorobenzene-d4	13.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Deltacyclene	118	C9H10	007785-10-6	59
2		Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	53
3		Indole	117	C8H7N	000120-72-9	46
4		trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	45
5		Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	45



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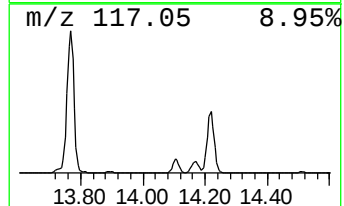
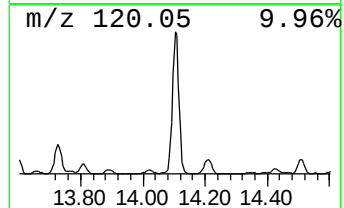
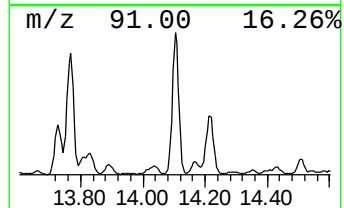
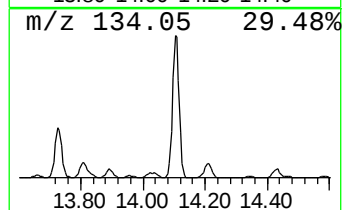
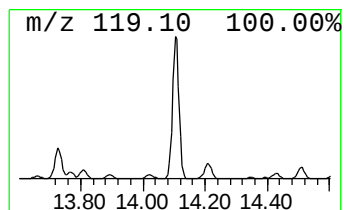
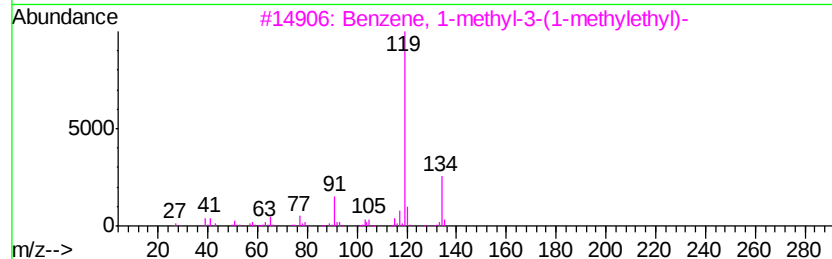
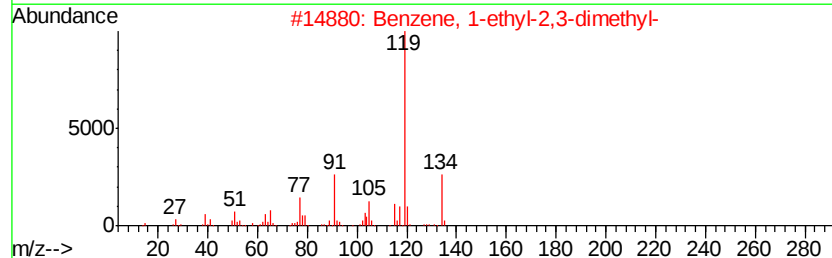
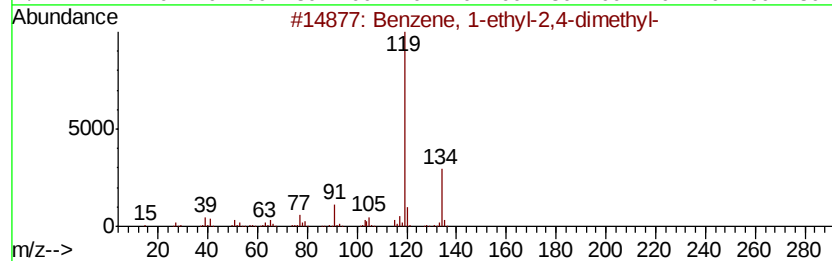
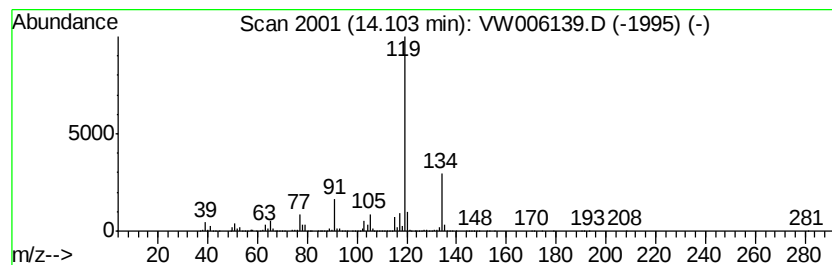
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.10	479.15 ug/l	1362400	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3		Benzene, 1-methyl-3-(1-methylethyl-...)	134	C10H14	000535-77-3	95
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5		o-Cymene	134	C10H14	000527-84-4	95



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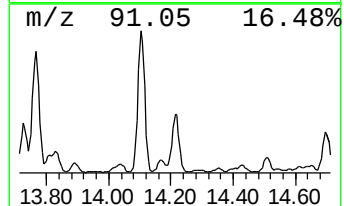
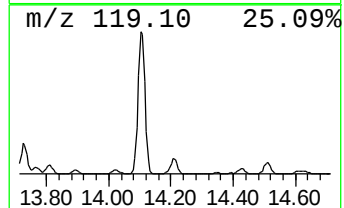
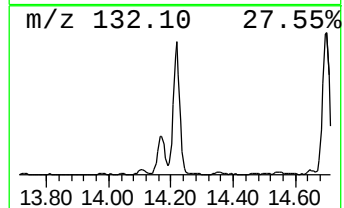
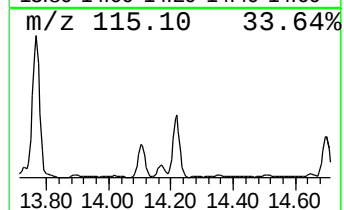
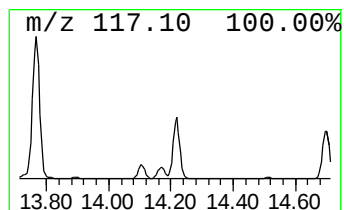
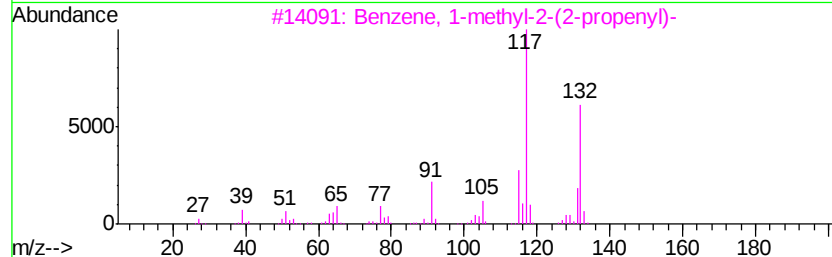
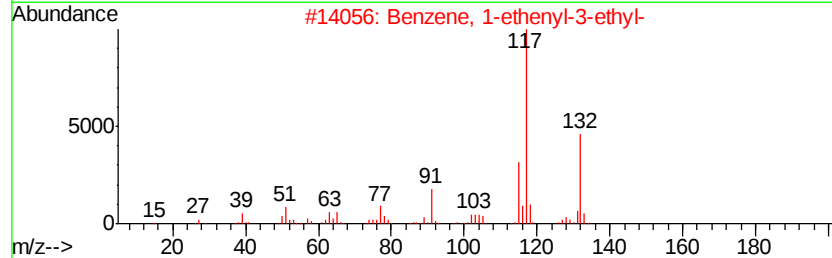
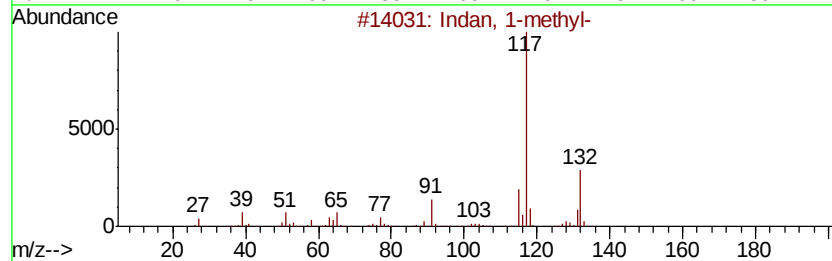
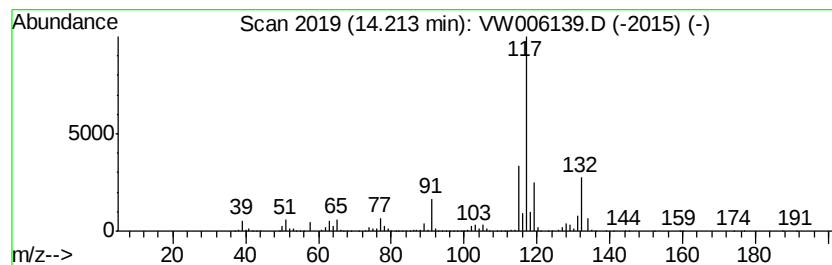
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Quant Title : SW846 8260

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

Peak Number 4 Indan, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	261.98 ug/l	744890	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indan, 1-methyl-	132	C10H12	000767-58-8	76
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	74
3		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	68
4		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	64
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101718\
Data File : VW006139.D
Acq On : 17 Oct 2018 20:45
Operator : VA/AP
Sample : J5515-18
Misc : 5.96G/5ML/MSVOA W/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
AOC1-01A

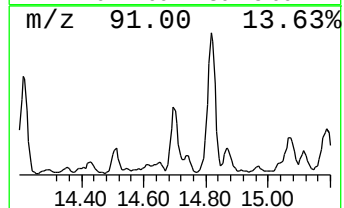
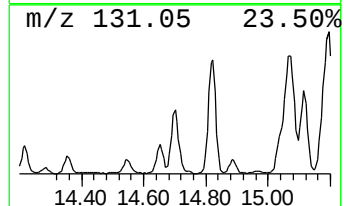
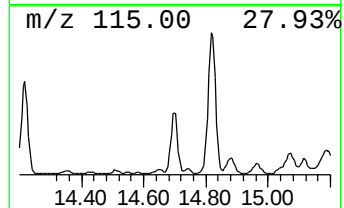
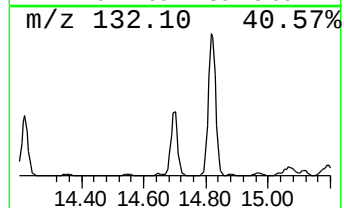
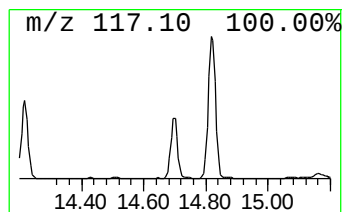
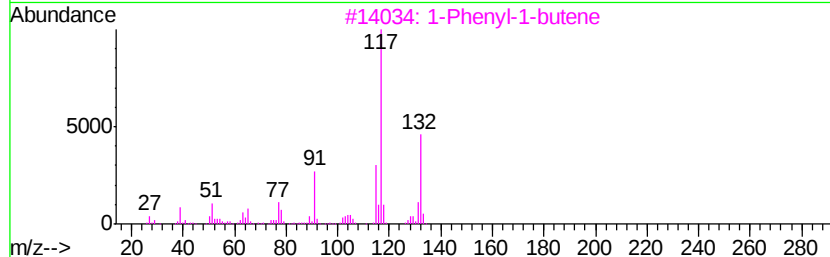
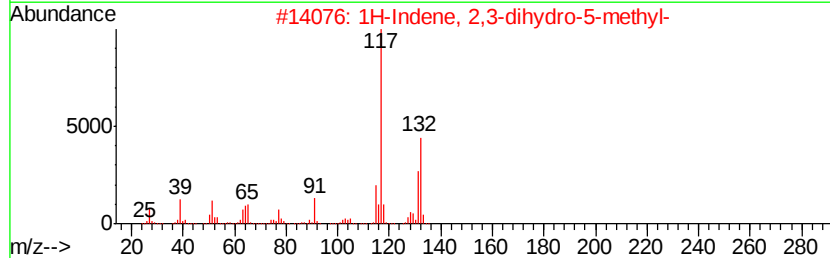
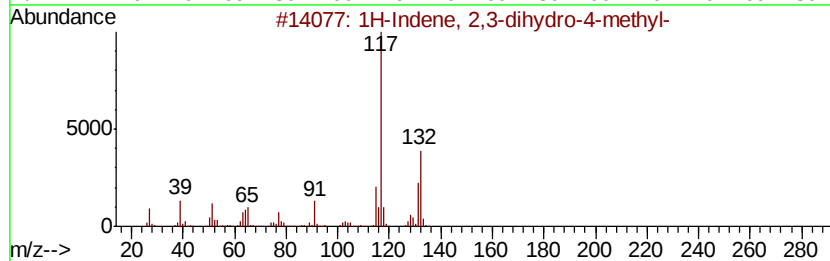
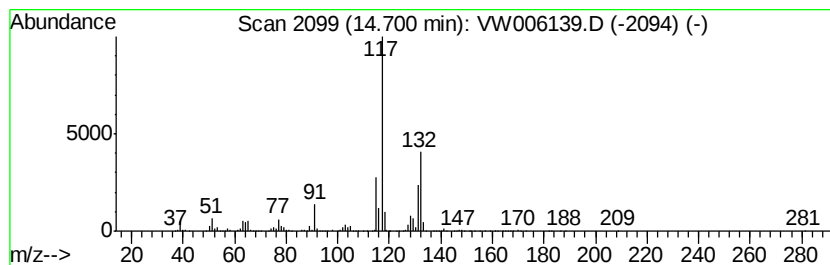
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Quant Title : SW846 8260

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

Peak Number 5 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.70	196.70 ug/l	559286	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	90
4		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	90
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101718\
Data File : VW006139.D
Acq On : 17 Oct 2018 20:45
Operator : VA/AP
Sample : J5515-18
Misc : 5.96G/5ML/MSVOA W/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
AOC1-01A

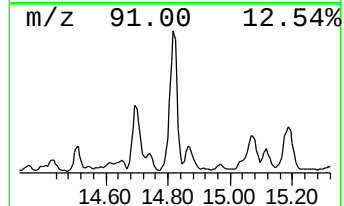
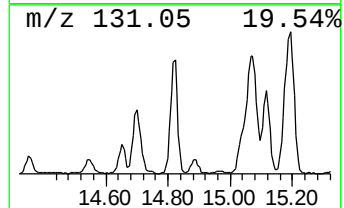
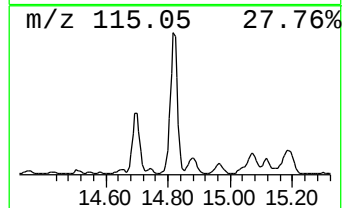
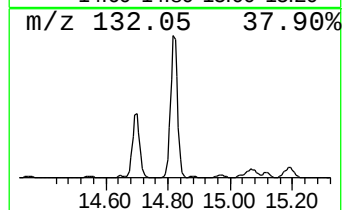
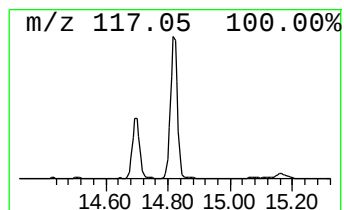
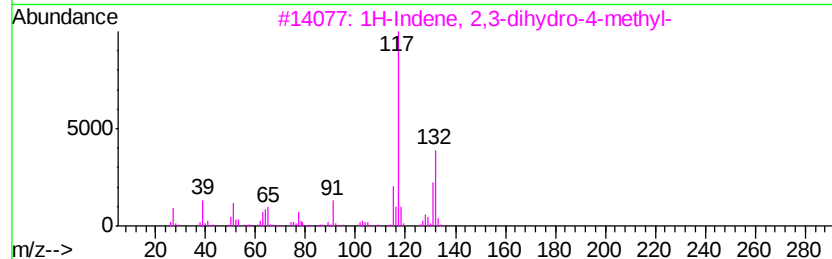
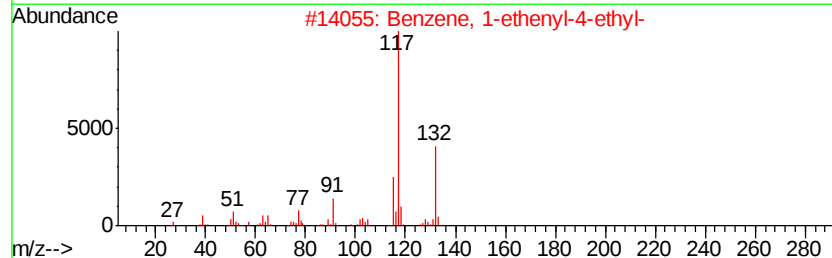
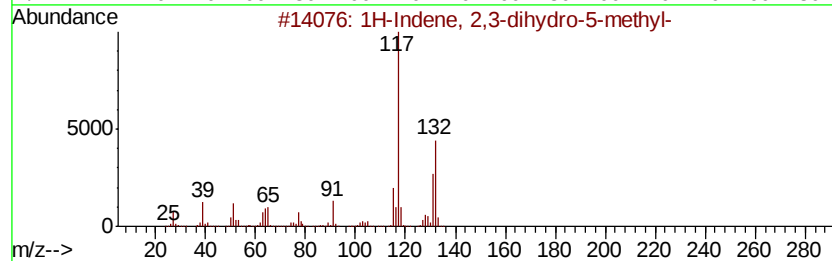
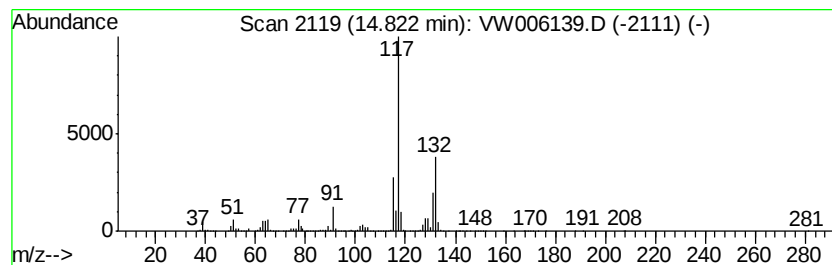
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Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	498.52 ug/l	1417460	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
4		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101718\
Data File : VW006139.D
Acq On : 17 Oct 2018 20:45
Operator : VA/AP
Sample : J5515-18
Misc : 5.96G/5ML/MSVOA W/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
AOC1-01A

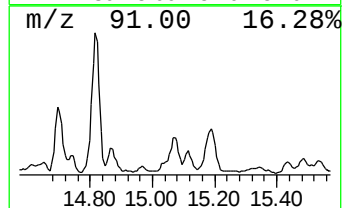
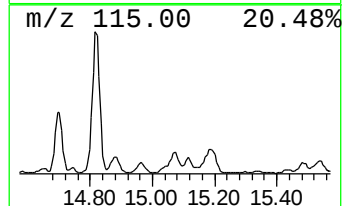
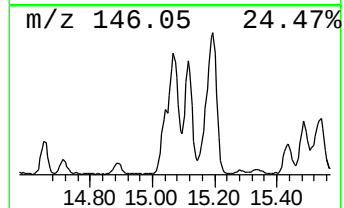
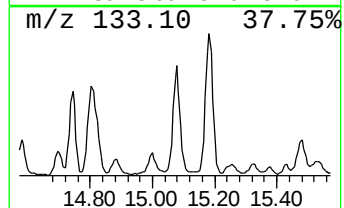
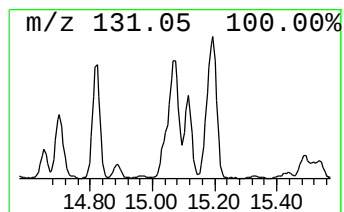
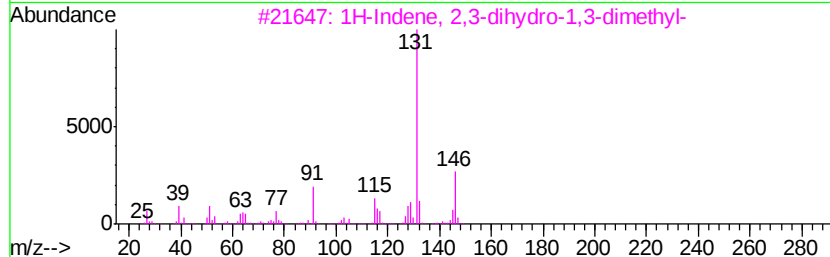
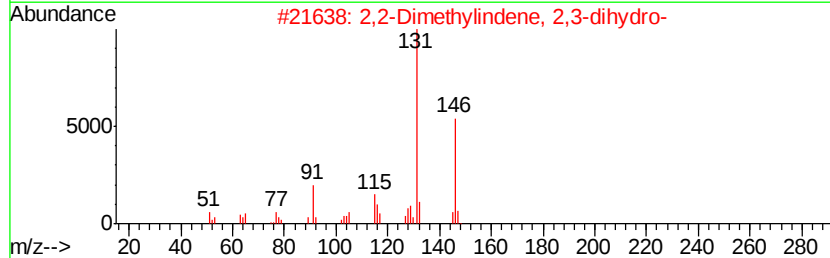
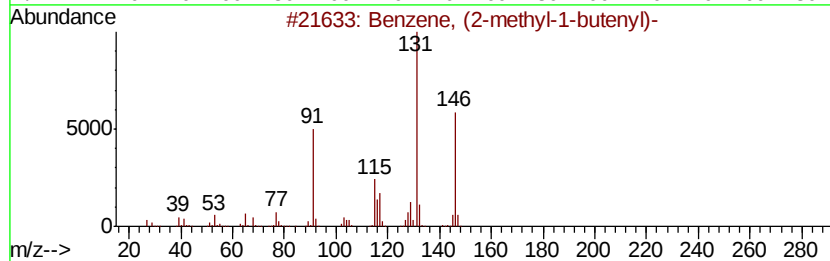
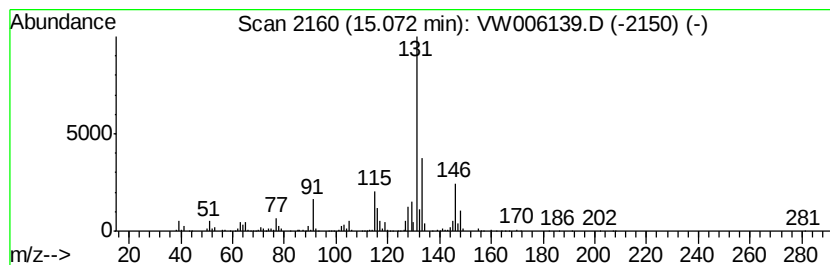
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, (2-methyl-1-butenyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	152.04 ug/l	432297	1,4-Dichlorobenzene-d4	13.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	91
2		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91
3		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90
4		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	87
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101718\
Data File : VW006139.D
Acq On : 17 Oct 2018 20:45
Operator : VA/AP
Sample : J5515-18
Misc : 5.96G/5ML/MSVOA W/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
AOC1-01A

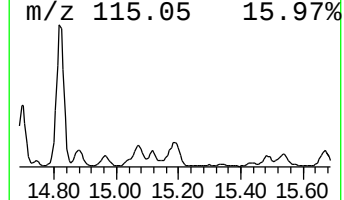
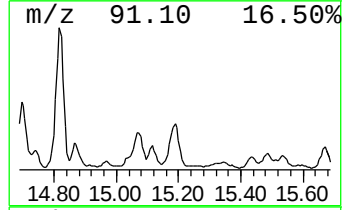
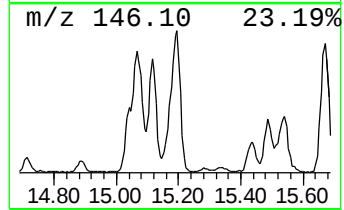
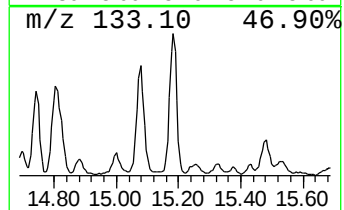
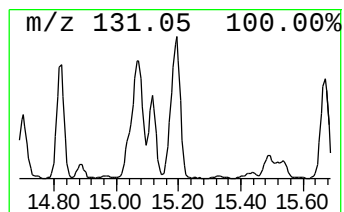
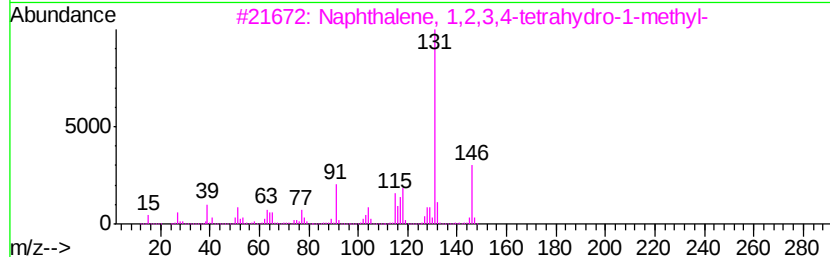
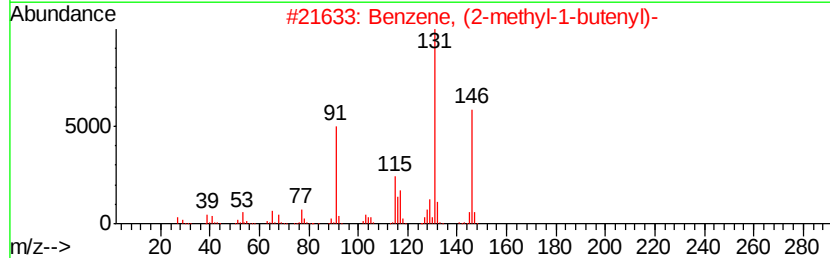
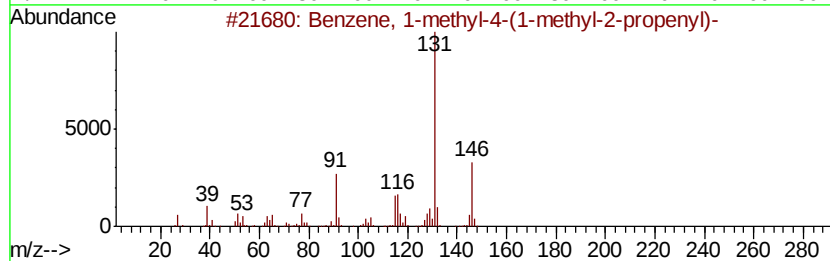
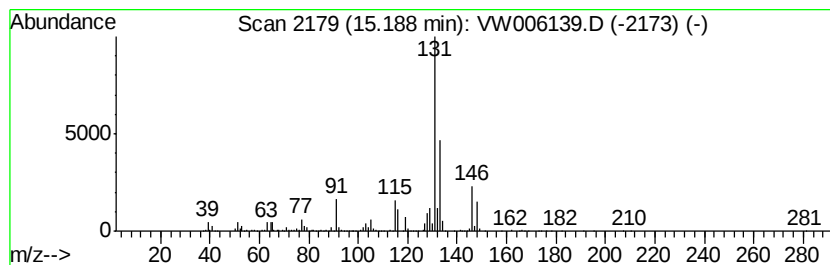
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-methyl-4-(1-meth... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	176.35 ug/l	501426	1,4-Dichlorobenzene-d4	13.57

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	86
2			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	70
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	70
4			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	70
5			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101718\
Data File : VW006139.D
Acq On : 17 Oct 2018 20:45
Operator : VA/AP
Sample : J5515-18
Misc : 5.96G/5ML/MSVOA W/SOIL
ALS Vial : 21 Sample Multiplier: 1

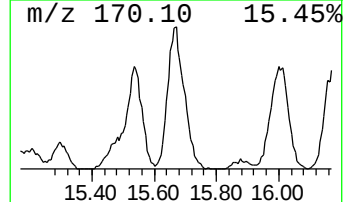
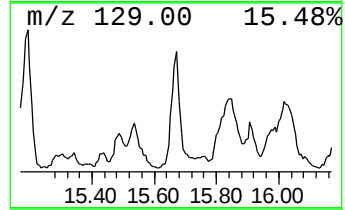
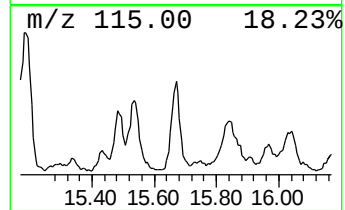
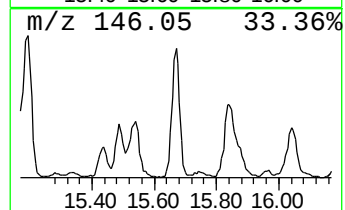
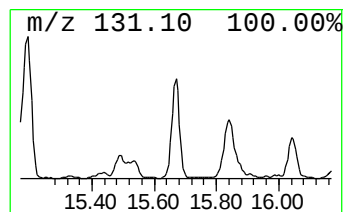
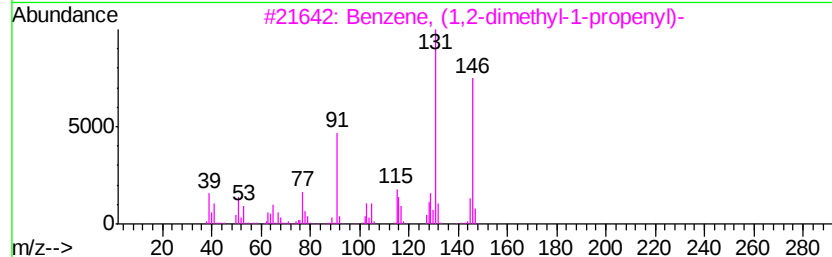
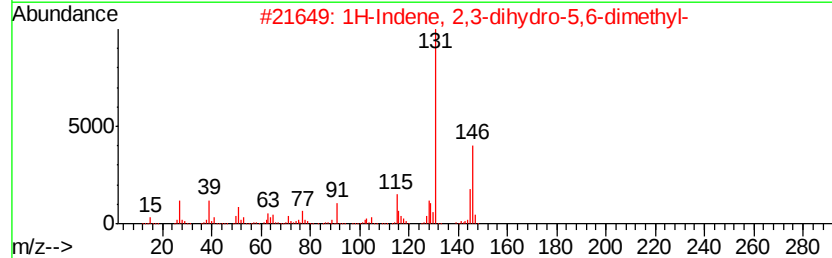
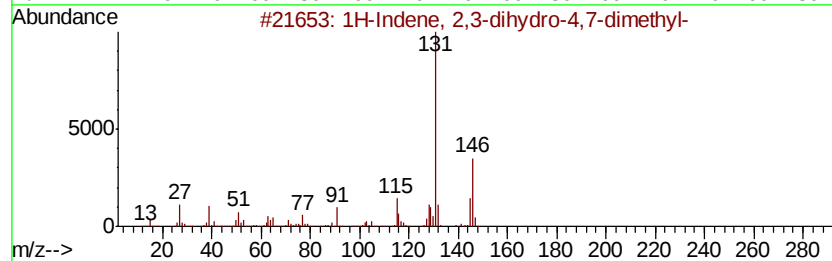
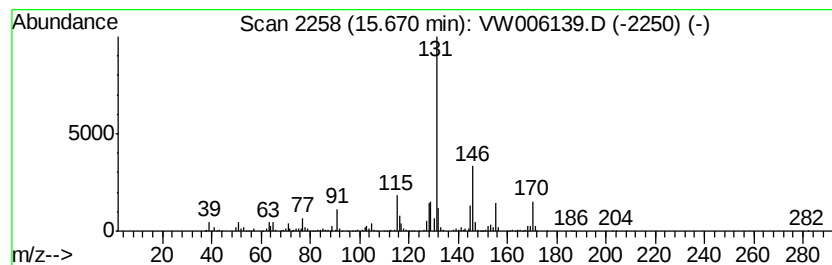
Instrument :
MSVOA_W
ClientSampleId :
AOC1-01A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.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.
15.67	135.26 ug/l	384591	1,4-Dichlorobenzene-d4	13.57
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9	95
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	87
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5	87
5	Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101718\
Data File : VW006139.D
Acq On : 17 Oct 2018 20:45
Operator : VA/AP
Sample : J5515-18
Misc : 5.96G/5ML/MSVOA W/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
AOC1-01A

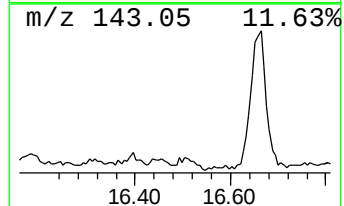
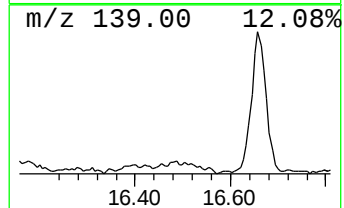
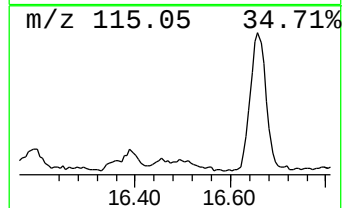
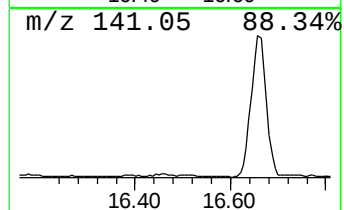
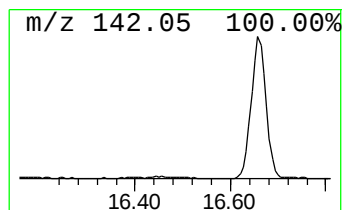
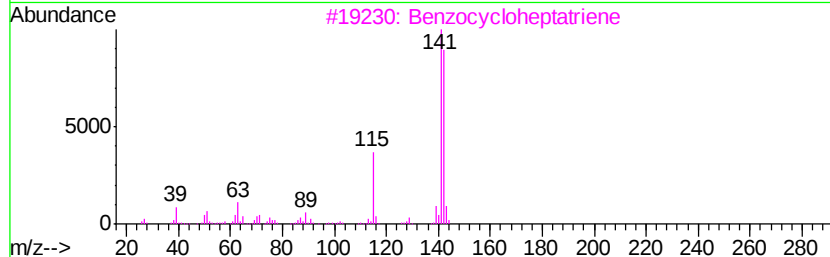
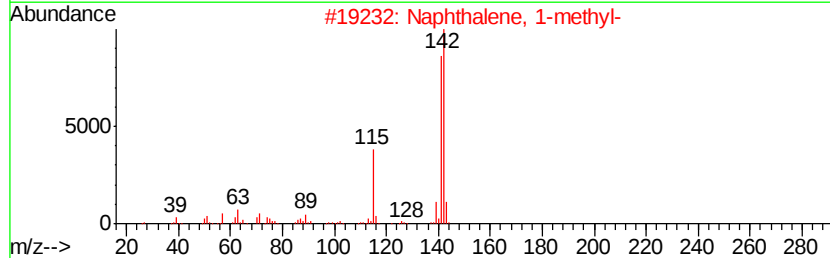
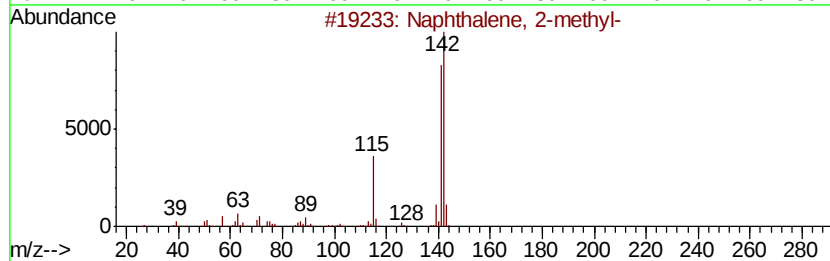
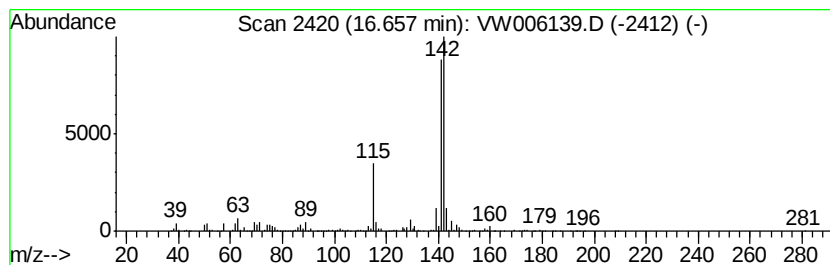
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.66	121.61 ug/l	345777	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	96
3		Benzocycloheptatriene	142	C11H10	000264-09-5	91
4		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	90
5		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	87



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW101718\
Data File : VW006139.D
Acq On : 17 Oct 2018 20:45
Operator : VA/AP
Sample : J5515-18
Misc : 5.96G/5ML/MSVOA_W/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
AOC1-01A

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.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, 1,2-diet...	13.72	182.8	ug/l	519741	4	13.57	142168	50.0
Deltacyclene	13.77	501.1	ug/l	1424670	4	13.57	142168	50.0
Benzene, 1-ethyl-...	14.10	479.1	ug/l	1362400	4	13.57	142168	50.0
Indan, 1-methyl-	14.21	262.0	ug/l	744890	4	13.57	142168	50.0
1H-Indene, 2,3-di...	14.70	196.7	ug/l	559286	4	13.57	142168	50.0
1H-Indene, 2,3-di...	14.82	498.5	ug/l	1417460	4	13.57	142168	50.0
Benzene, (2-methy...	15.07	152.0	ug/l	432297	4	13.57	142168	50.0
Benzene, 1-methyl...	15.19	176.3	ug/l	501426	4	13.57	142168	50.0
1H-Indene, 2,3-di...	15.67	135.3	ug/l	384591	4	13.57	142168	50.0
Naphthalene, 2-me...	16.66	121.6	ug/l	345777	4	13.57	142168	50.0