

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW101918\
 Data File : VW006230.D
 Acq On : 19 Oct 2018 17:29
 Operator : VA/AP
 Sample : J5605-01
 Misc : 5.28G/5ML/MSVOA W/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 TP-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\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	2.629	114	119	122	rBV6	2525	4977	1.74%	0.169%
2	3.032	181	185	192	rVB5	1549	3195	1.12%	0.108%
3	3.531	262	267	273	rVB2	3150	7050	2.47%	0.239%
4	4.123	354	364	373	rBV	4964	13887	4.86%	0.471%
5	4.915	485	494	495	rBV3	2645	4880	1.71%	0.165%
6	5.434	571	579	593	rVB5	1598	5745	2.01%	0.195%
7	5.940	654	662	673	rVB4	1921	5262	1.84%	0.178%
8	6.988	824	834	842	rBV3	6344	18620	6.52%	0.631%
9	7.153	851	861	864	rBV2	2020	5424	1.90%	0.184%
10	7.890	973	982	986	rBV2	29727	69638	24.38%	2.361%
11	7.951	986	992	1000	rVV	68417	156918	54.93%	5.319%
12	8.037	1002	1006	1011	rVB4	1497	2991	1.05%	0.101%
13	8.165	1016	1027	1032	rBV5	3924	9561	3.35%	0.324%
14	8.311	1043	1051	1060	rBV	28892	62634	21.92%	2.123%
15	8.439	1063	1072	1078	rVV4	3410	8658	3.03%	0.293%
16	8.512	1078	1084	1089	rVV3	3826	8387	2.94%	0.284%
17	8.579	1089	1095	1101	rVB4	3646	8513	2.98%	0.289%
18	8.848	1128	1139	1151	rBV	83234	171202	59.93%	5.804%
19	9.079	1170	1177	1180	rVV5	1675	3803	1.33%	0.129%
20	9.341	1207	1220	1226	rBV	24722	60038	21.02%	2.035%
21	9.518	1244	1249	1255	rVB4	3942	7313	2.56%	0.248%
22	9.616	1258	1265	1273	rVB6	3242	7467	2.61%	0.253%
23	9.756	1282	1288	1293	rVB3	3835	6940	2.43%	0.235%
24	9.853	1298	1304	1309	rBV3	6120	12166	4.26%	0.412%
25	9.994	1324	1327	1335	rVB6	3626	6406	2.24%	0.217%
26	10.091	1335	1343	1349	rBV5	4559	11879	4.16%	0.403%
27	10.213	1357	1363	1366	rBV4	2465	5266	1.84%	0.179%
28	10.329	1374	1382	1389	rBV	119524	223946	78.39%	7.591%
29	10.451	1397	1402	1404	rBV3	3388	5148	1.80%	0.175%
30	10.512	1405	1412	1420	rVB6	8168	24613	8.62%	0.834%
31	10.609	1420	1428	1433	rBV8	2284	5677	1.99%	0.192%
32	10.670	1433	1438	1443	rBV3	7819	13785	4.83%	0.467%
33	10.756	1445	1452	1459	rVB5	8423	20823	7.29%	0.706%
34	10.853	1463	1468	1473	rVB4	2413	4725	1.65%	0.160%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	10.945	1473	1483	1487	rBV5	3005	6384	2.23%	0.216%
36	11.012	1487	1494	1495	rBV6	3225	5782	2.02%	0.196%
37	11.109	1506	1510	1512	rBV3	3013	4671	1.64%	0.158%
38	11.146	1512	1516	1518	rVV3	5721	9375	3.28%	0.318%
39	11.183	1518	1522	1527	rVV	11942	22225	7.78%	0.753%
40	11.231	1527	1530	1535	rVV2	9379	14598	5.11%	0.495%
41	11.292	1535	1540	1544	rVV4	4063	6762	2.37%	0.229%
42	11.341	1545	1548	1552	rVB5	4557	5979	2.09%	0.203%
43	11.414	1556	1560	1561	rVV3	2099	3402	1.19%	0.115%
44	11.445	1562	1565	1571	rVB4	4634	8726	3.05%	0.296%
45	11.518	1572	1577	1581	rBV3	4214	8176	2.86%	0.277%
46	11.634	1590	1596	1603	rBV	99294	171688	60.10%	5.820%
47	11.731	1608	1612	1616	rVV5	3548	6370	2.23%	0.216%
48	11.768	1616	1618	1622	rVB5	2489	3032	1.06%	0.103%
49	11.847	1622	1631	1641	rBV	47610	95155	33.31%	3.226%
50	12.152	1674	1681	1682	rBV3	4716	9377	3.28%	0.318%
51	12.347	1703	1713	1716	rBV7	6265	14165	4.96%	0.480%
52	12.469	1728	1733	1739	rBV	13841	22299	7.81%	0.756%
53	12.621	1752	1758	1765	rBV	73599	124348	43.53%	4.215%
54	12.707	1765	1772	1780	rVV10	4688	14820	5.19%	0.502%
55	12.810	1780	1789	1794	rVB	29355	51580	18.06%	1.748%
56	12.908	1800	1805	1808	rVV	28641	47813	16.74%	1.621%
57	12.950	1808	1812	1819	rVB	76588	121607	42.57%	4.122%
58	13.109	1832	1838	1845	rVB	59492	96675	33.84%	3.277%
59	13.182	1848	1850	1856	rVB7	2000	3497	1.22%	0.119%
60	13.255	1856	1862	1870	rBV	179872	285675	100.00%	9.684%
61	13.389	1877	1884	1890	rVB2	7048	16894	5.91%	0.573%
62	13.505	1899	1903	1907	rVB	8000	11039	3.86%	0.374%
63	13.566	1907	1913	1916	rBV	81065	145814	51.04%	4.943%
64	13.725	1934	1939	1942	rBV	17625	28644	10.03%	0.971%
65	13.767	1942	1946	1948	rVV	37239	62478	21.87%	2.118%
66	13.798	1948	1951	1961	rVB2	45512	83640	29.28%	2.835%
67	13.889	1961	1966	1970	rVB4	7848	11936	4.18%	0.405%
68	13.956	1972	1977	1982	rBV	14450	22809	7.98%	0.773%
69	14.017	1982	1987	1990	rVV	24202	39275	13.75%	1.331%
70	14.048	1990	1992	1997	rVV	18592	25998	9.10%	0.881%
71	14.103	1997	2001	2007	rVB	35622	55591	19.46%	1.884%

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Integration Parameters: RTEINT.P

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72	14.170	2007	2012	2015	rBV2	8669	14001	4.90%	0.475%
73	14.212	2015	2019	2025	rVB2	34872	56856	19.90%	1.927%
74	14.347	2036	2041	2045	rBV	9637	13368	4.68%	0.453%
75	14.426	2046	2054	2057	rVV	18361	32622	11.42%	1.106%
76	14.469	2057	2061	2065	rVV	28372	42147	14.75%	1.429%
77	14.511	2065	2068	2071	rVV2	6591	8829	3.09%	0.299%
78	14.548	2071	2074	2080	rVB5	5637	8489	2.97%	0.288%
79	14.651	2087	2091	2094	rVB5	3177	4327	1.51%	0.147%
80	14.700	2094	2099	2102	rVV2	9616	15966	5.59%	0.541%
81	14.743	2102	2106	2110	rVB3	5347	9065	3.17%	0.307%
82	14.816	2111	2118	2124	rBV3	23803	48897	17.12%	1.658%
83	14.865	2124	2126	2132	rVB4	2241	4000	1.40%	0.136%
84	14.968	2139	2143	2147	rBV7	2526	3762	1.32%	0.128%
85	15.078	2154	2161	2165	rBV7	6058	13063	4.57%	0.443%
86	15.182	2174	2178	2186	rVB4	6814	15623	5.47%	0.530%
87	15.566	2239	2241	2247	rVB6	2074	3120	1.09%	0.106%

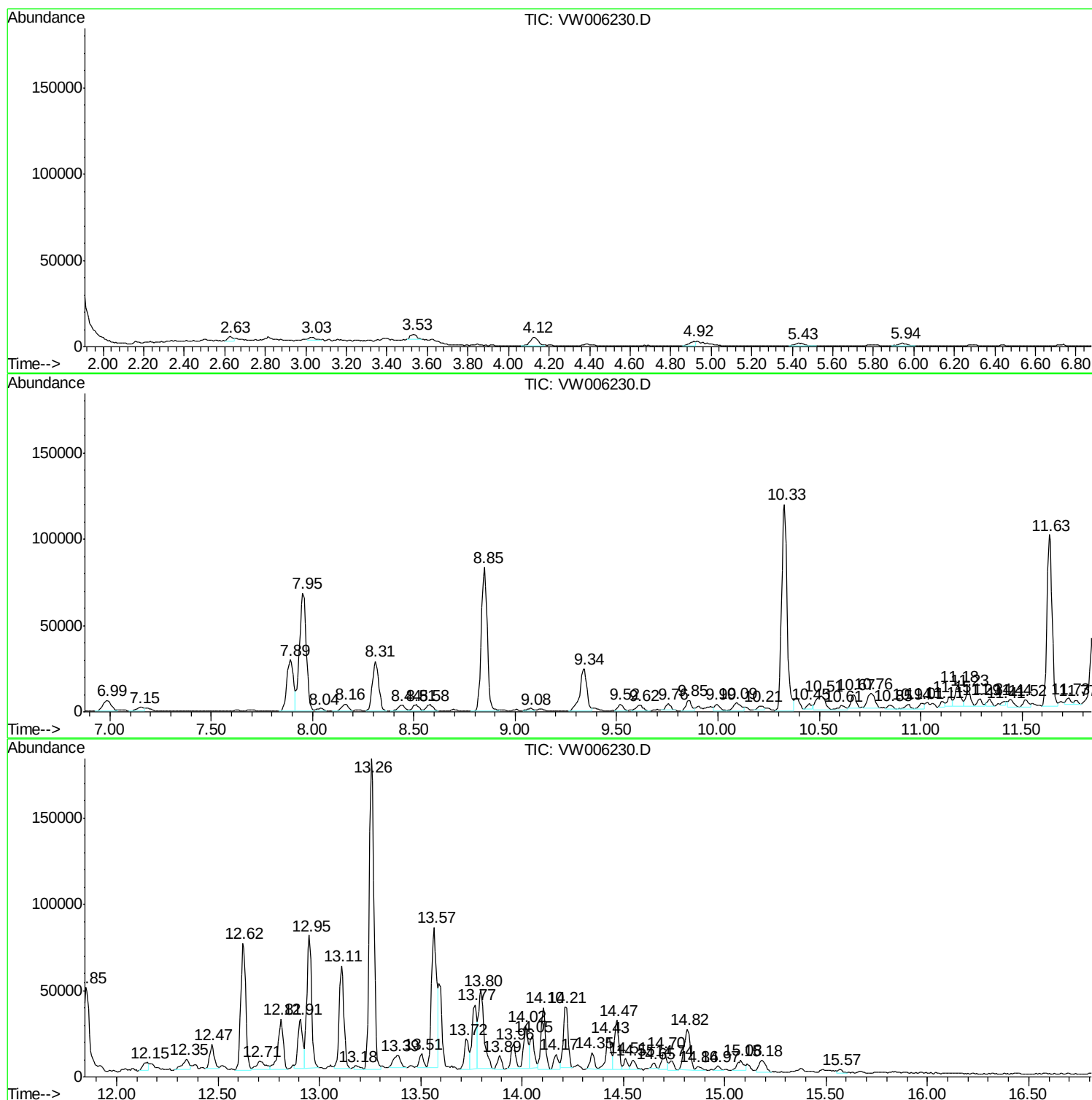
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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



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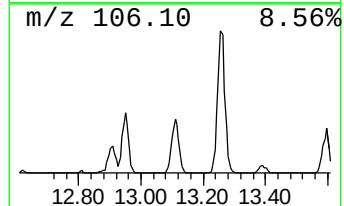
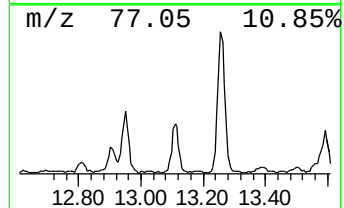
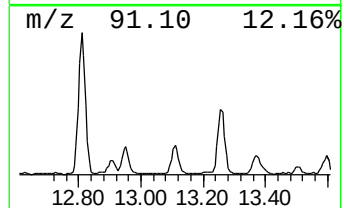
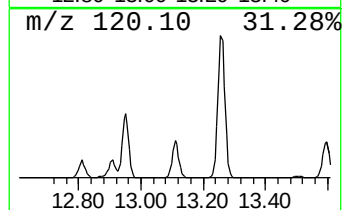
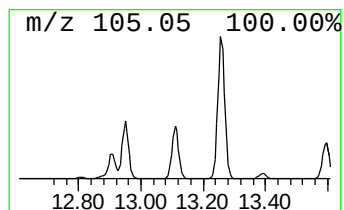
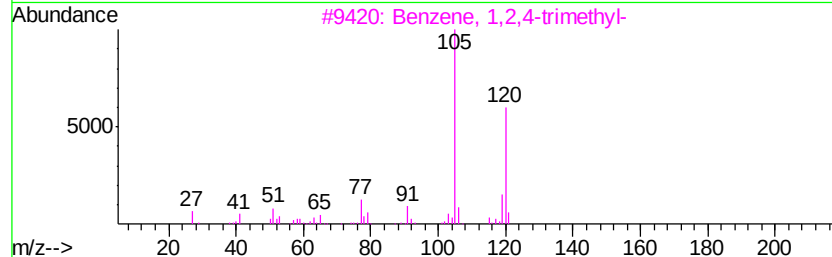
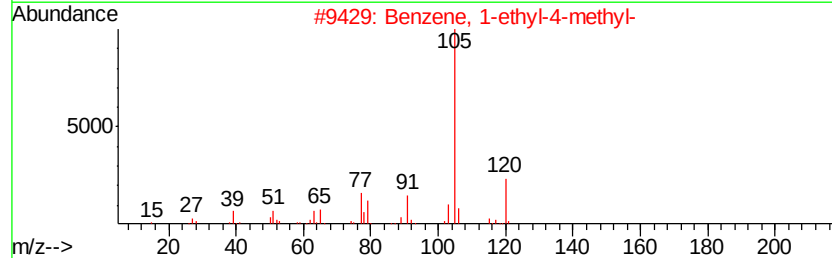
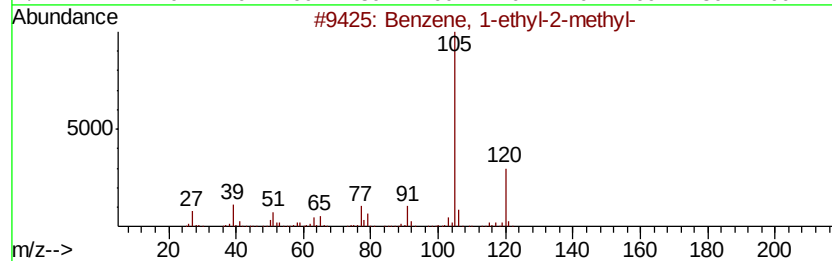
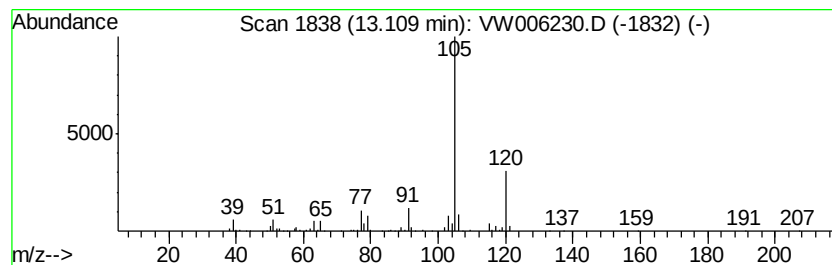
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.11	33.15 ug/l	96675	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Mesitylene	120	C9H12	000108-67-8	91



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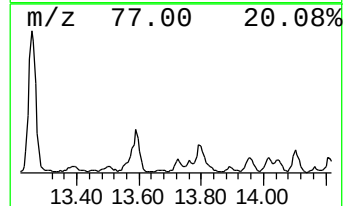
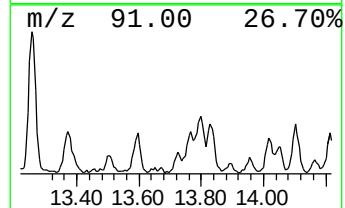
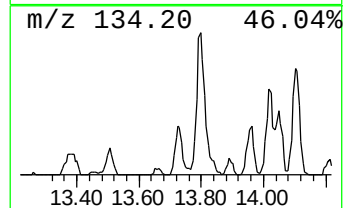
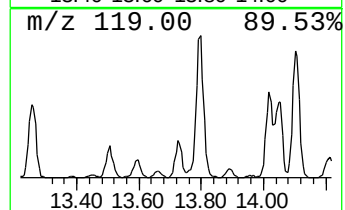
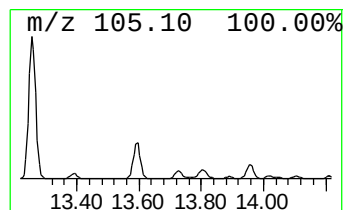
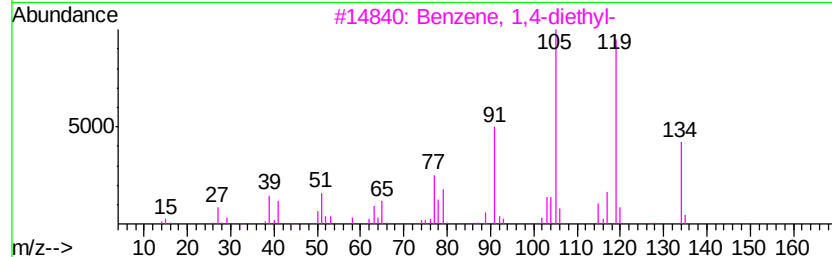
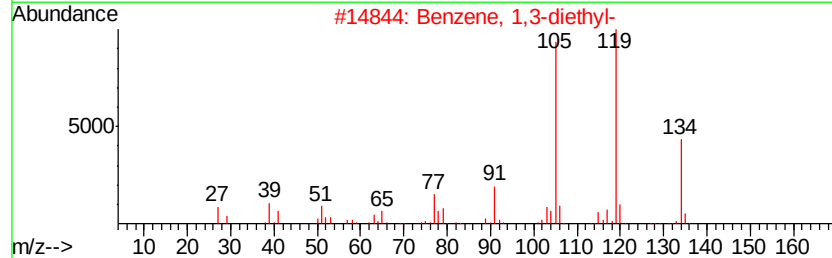
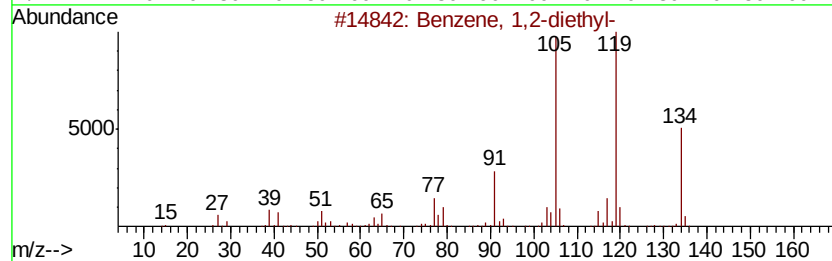
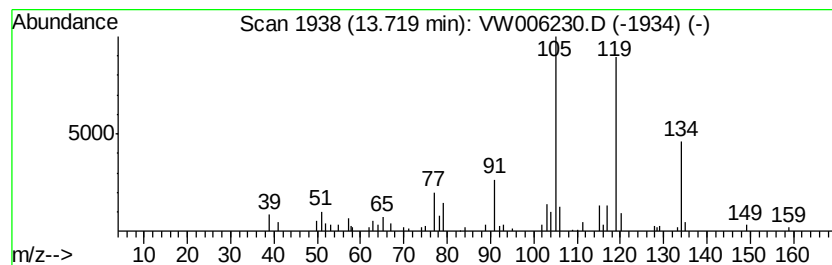
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1,2-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	9.82 ug/l	28644	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	94
2		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94
3		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



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Misc : 5.28G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

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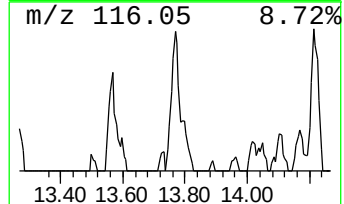
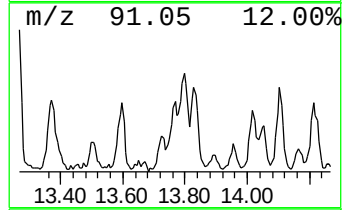
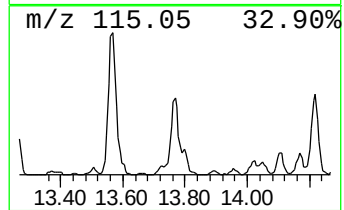
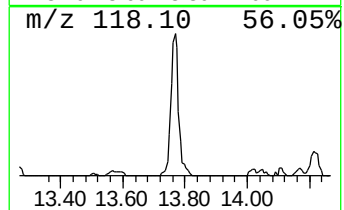
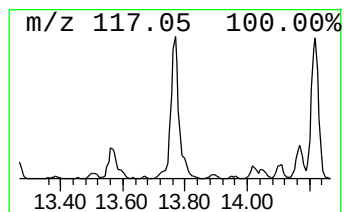
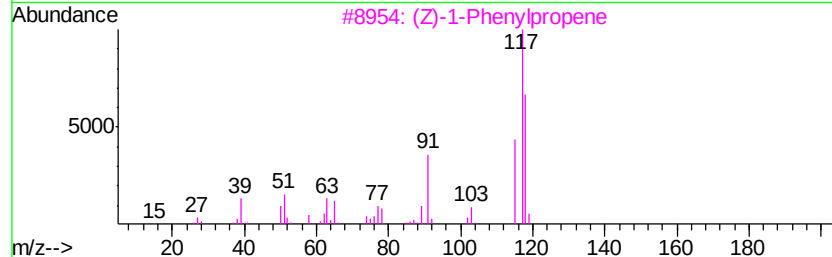
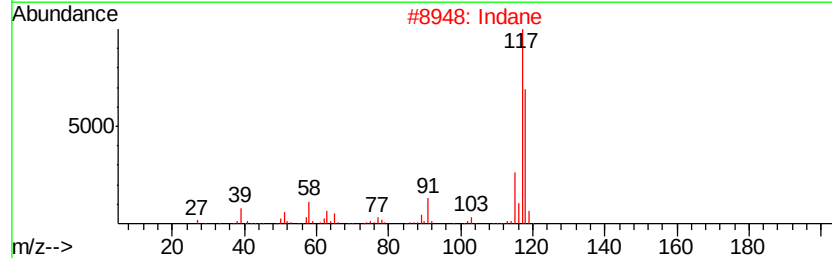
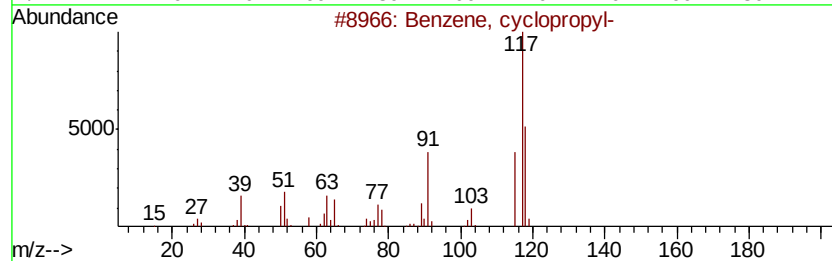
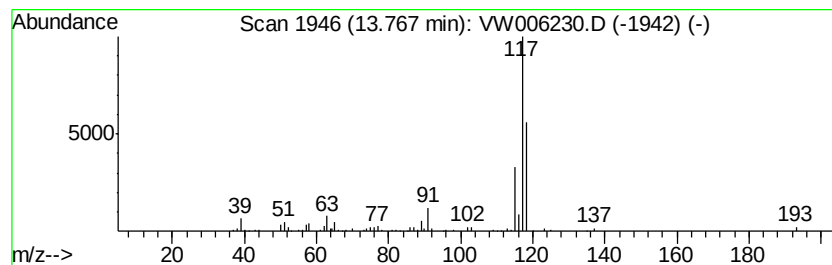
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, cyclopropyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	21.42 ug/l	62478	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, cvclopropyl-	118	C9H10	000873-49-4	74
2		Indane	118	C9H10	000496-11-7	74
3		(Z)-1-Phenylpropene	118	C9H10	000766-90-5	72
4		Benzene, 1-propenyl-	118	C9H10	000637-50-3	72
5		Benzene, 2-propenyl-	118	C9H10	000300-57-2	64



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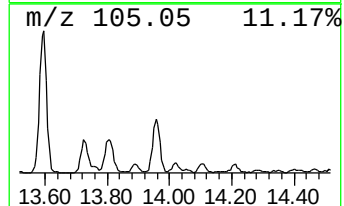
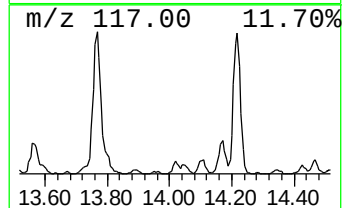
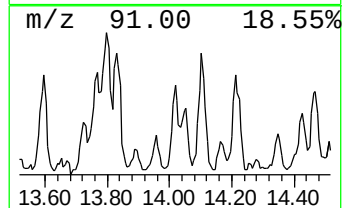
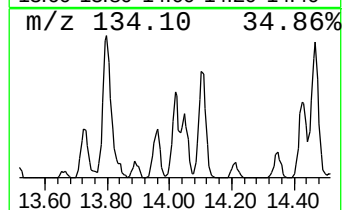
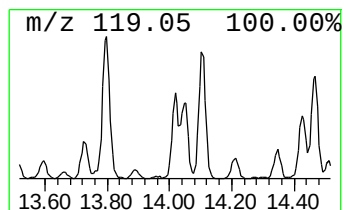
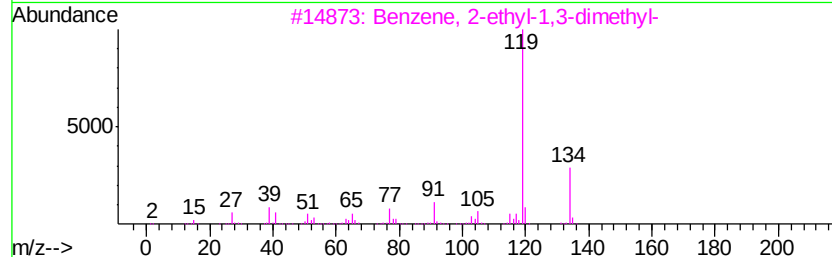
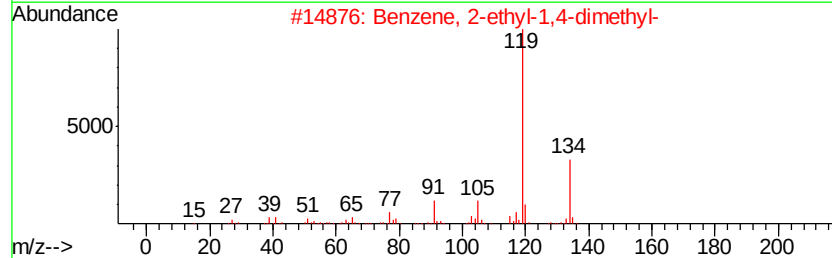
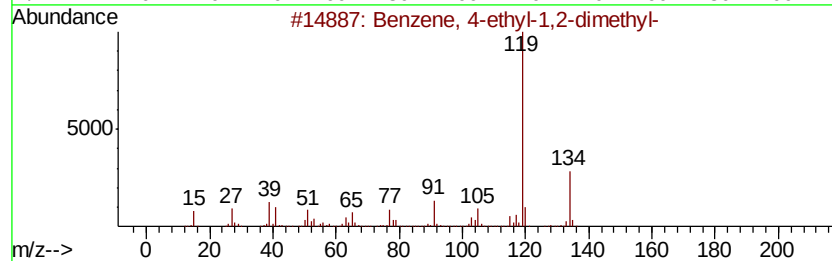
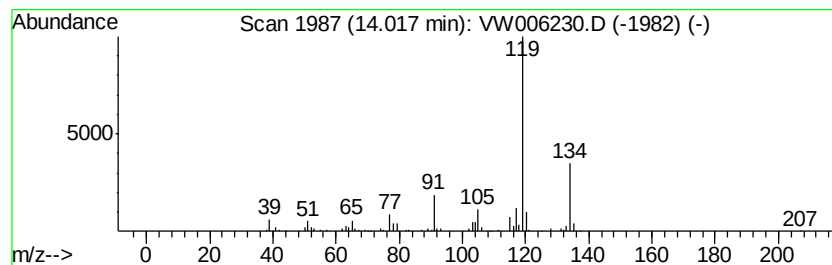
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.02	13.47 ug/l	39275	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
4		o-Cymene	134	C10H14	000527-84-4	94
5		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101918\
Data File : VW006230.D
Acq On : 19 Oct 2018 17:29
Operator : VA/AP
Sample : J5605-01
Misc : 5.28G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
TP-A

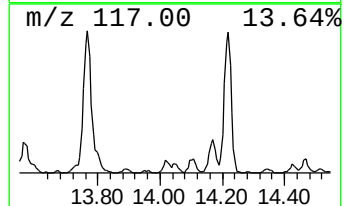
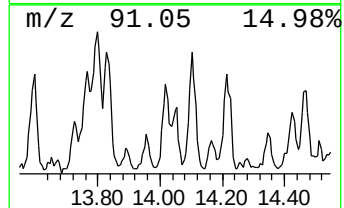
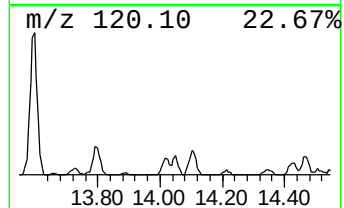
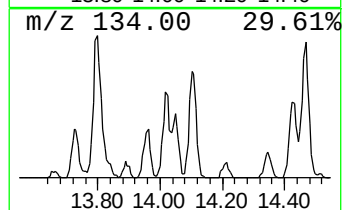
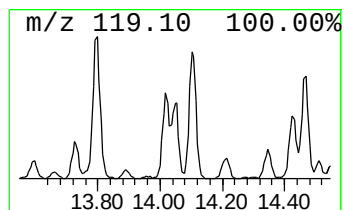
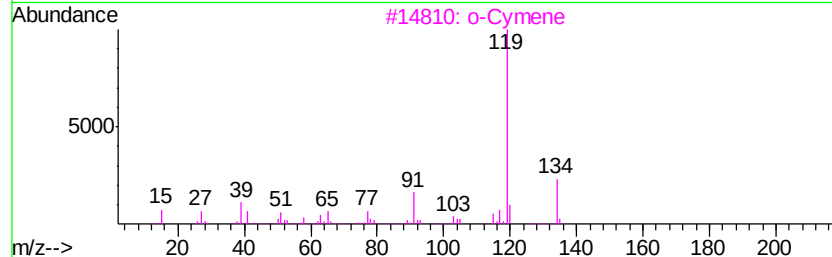
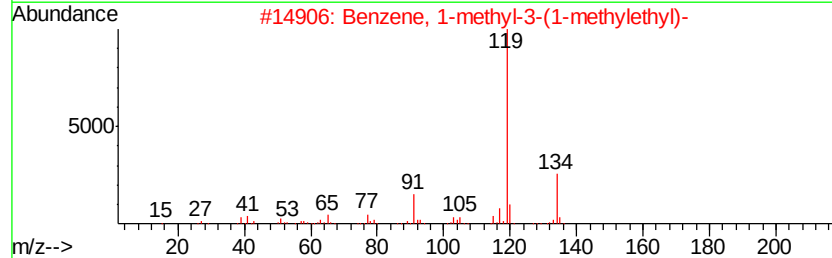
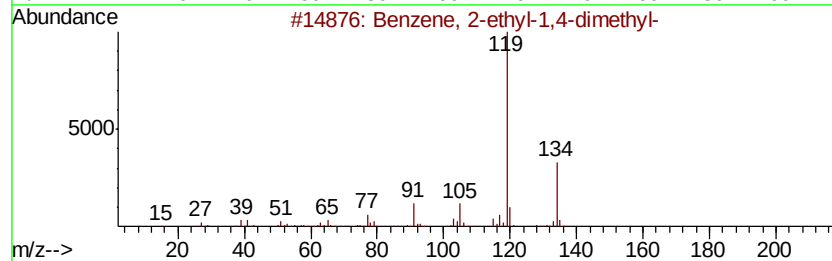
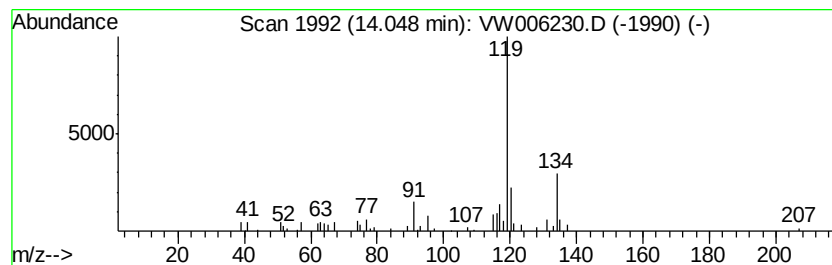
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.05	8.91 ug/l	25998	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	64
2		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	64
3		o-Cymene	134	C10H14	000527-84-4	64
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	64
5		p-Cymene	134	C10H14	000099-87-6	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101918\
Data File : VW006230.D
Acq On : 19 Oct 2018 17:29
Operator : VA/AP
Sample : J5605-01
Misc : 5.28G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
TP-A

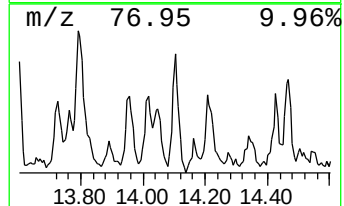
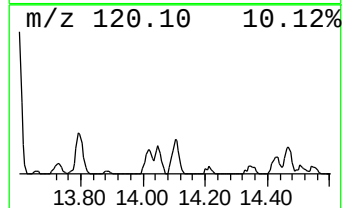
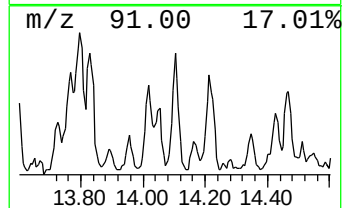
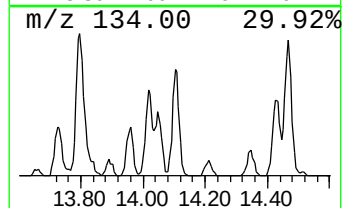
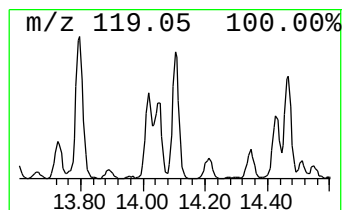
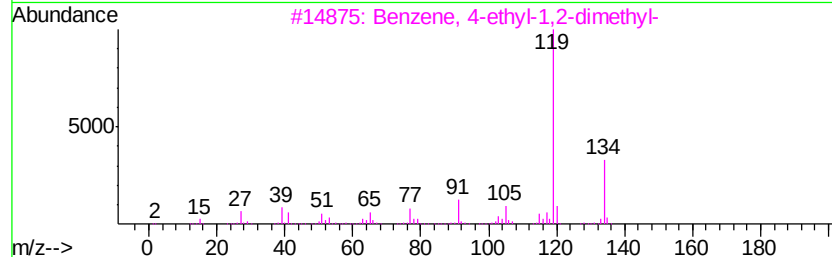
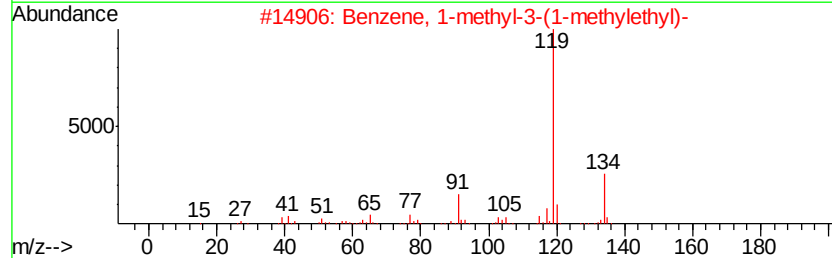
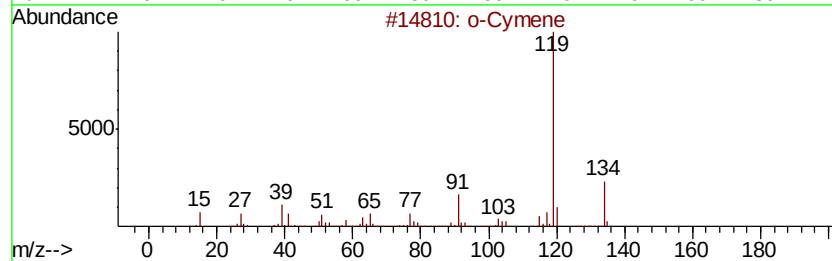
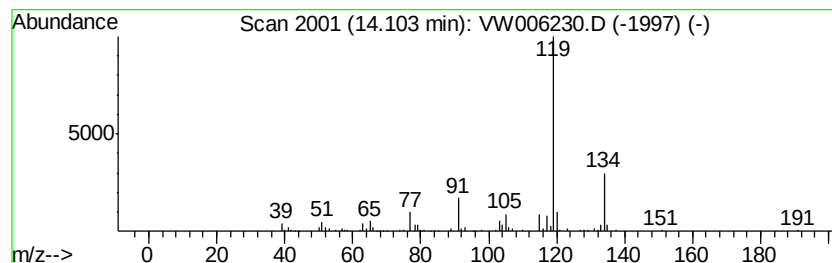
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Quant Title : SW846 8260

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

Peak Number 6 o-Cymene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.10	19.06 ug/l	55591	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	97
2		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101918\
Data File : VW006230.D
Acq On : 19 Oct 2018 17:29
Operator : VA/AP
Sample : J5605-01
Misc : 5.28G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

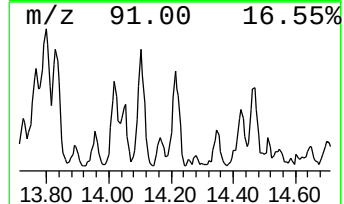
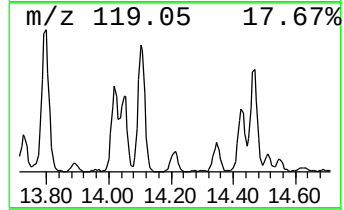
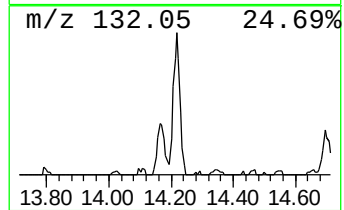
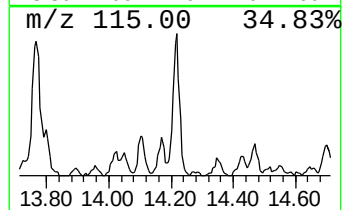
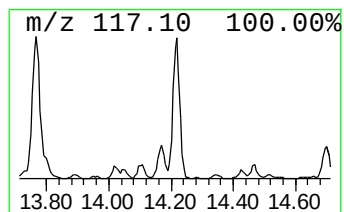
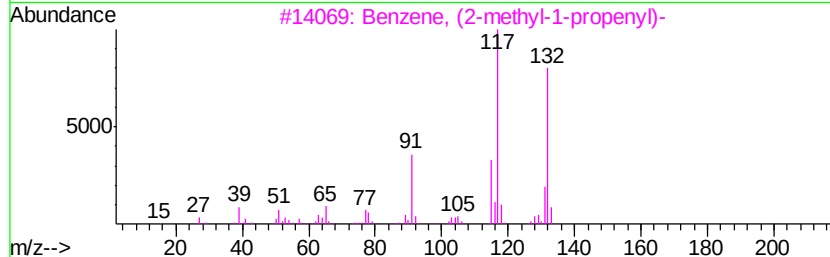
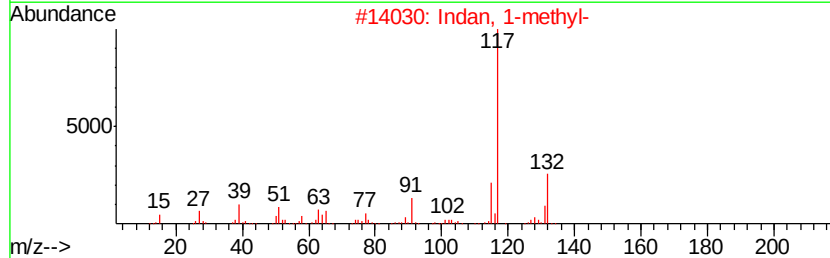
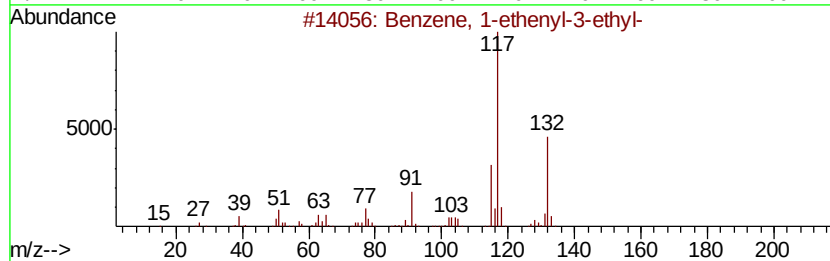
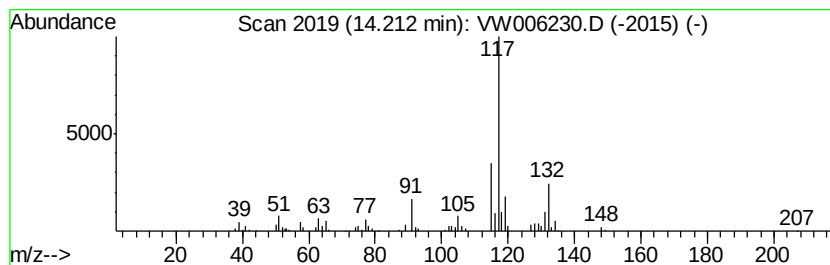
Instrument :
MSVOA_W
ClientSampleId :
TP-A

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 7 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	19.50 ug/l	56856	1,4-Dichlorobenzene-d4	13.57
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4 83
2		Indan, 1-methyl-	132 C10H12	000767-58-8 81
3		Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0 80
4		Benzene, (2-methyl-2-propenyl)-	132 C10H12	003290-53-7 80
5		3-Phenylbut-1-ene	132 C10H12	000934-10-1 72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101918\
Data File : VW006230.D
Acq On : 19 Oct 2018 17:29
Operator : VA/AP
Sample : J5605-01
Misc : 5.28G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
TP-A

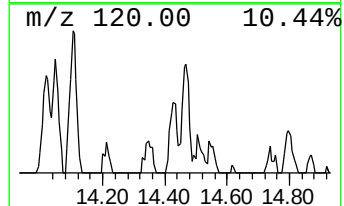
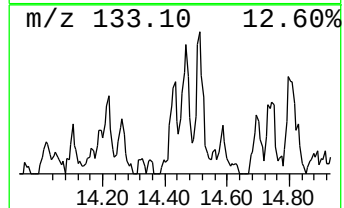
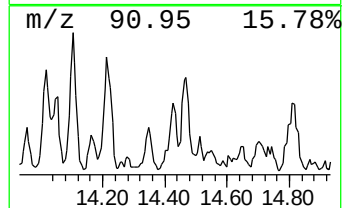
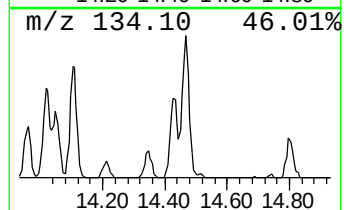
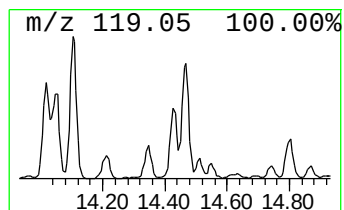
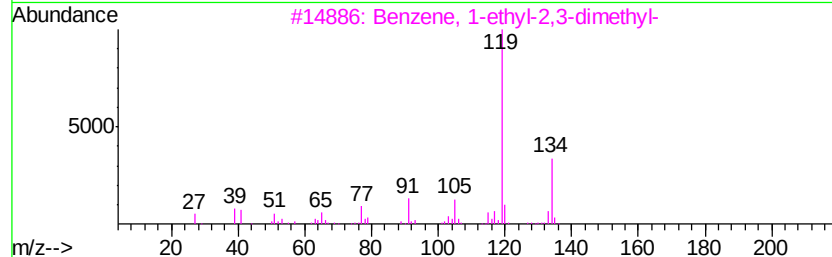
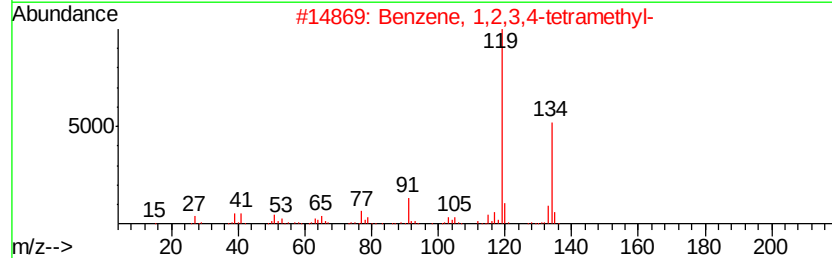
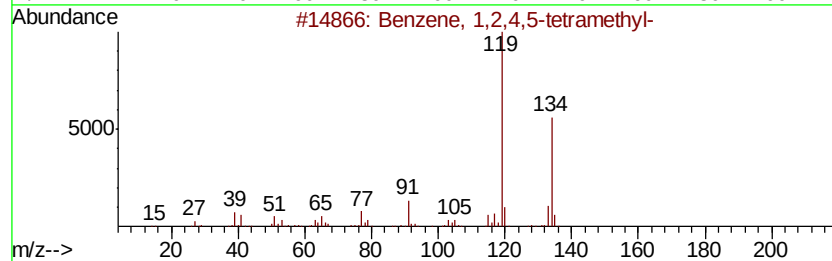
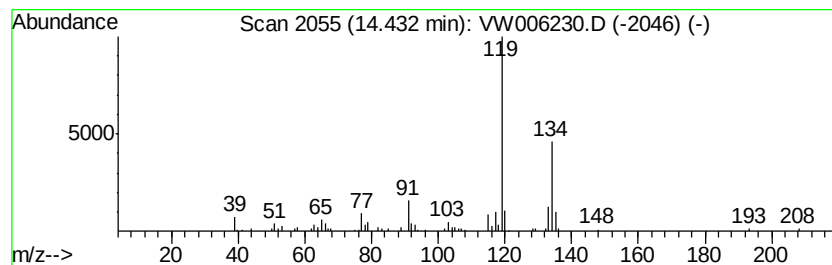
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.43	11.19 ug/l	32622	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	90
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101918\
Data File : VW006230.D
Acq On : 19 Oct 2018 17:29
Operator : VA/AP
Sample : J5605-01
Misc : 5.28G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
TP-A

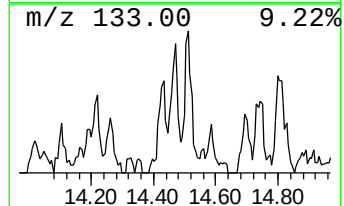
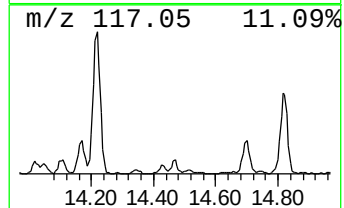
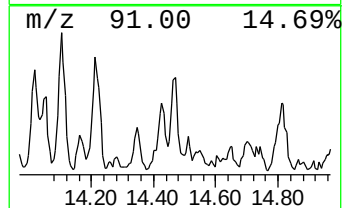
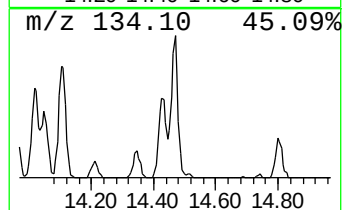
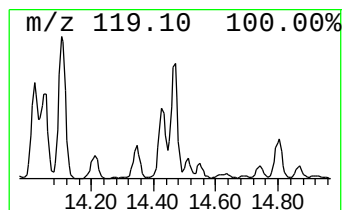
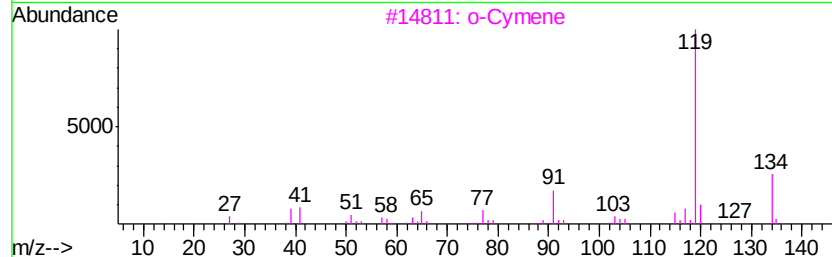
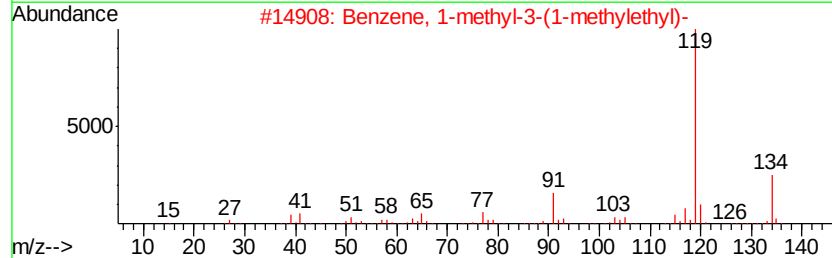
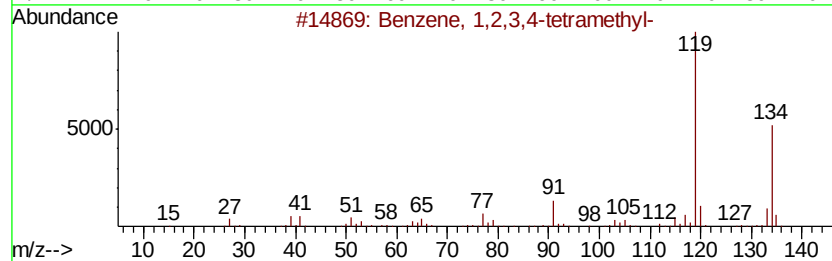
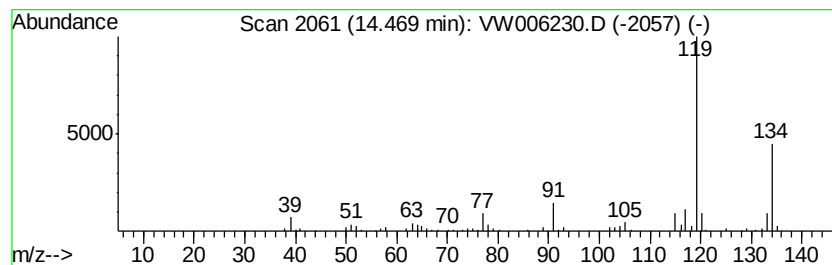
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.47	14.45 ug/l	42147	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91
2		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	90
3		o-Cymene	134	C10H14	000527-84-4	90
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	90
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101918\
Data File : VW006230.D
Acq On : 19 Oct 2018 17:29
Operator : VA/AP
Sample : J5605-01
Misc : 5.28G/5ML/MSVOA W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
TP-A

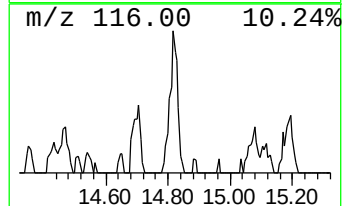
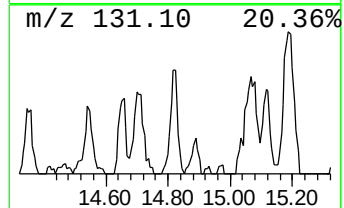
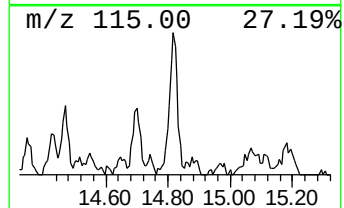
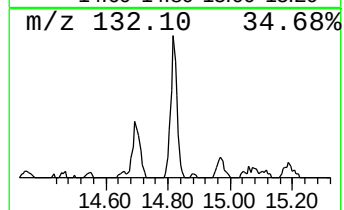
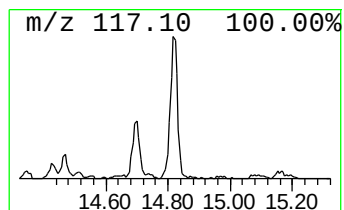
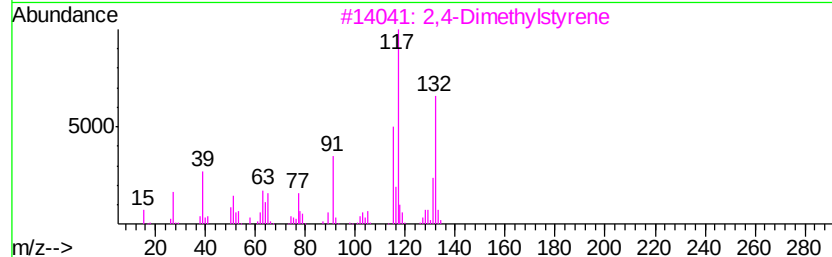
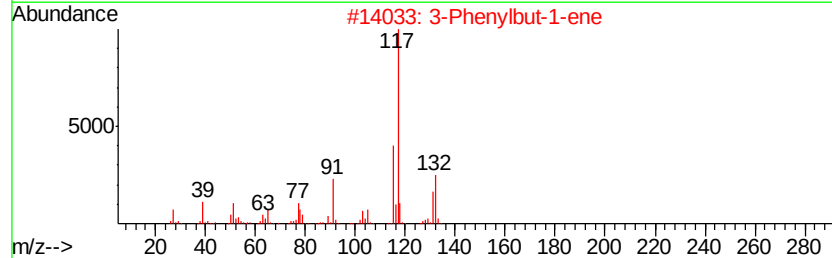
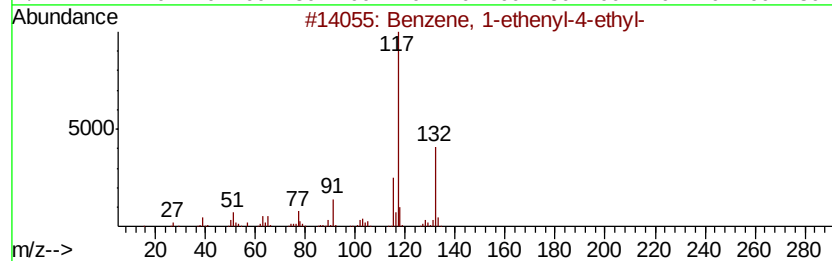
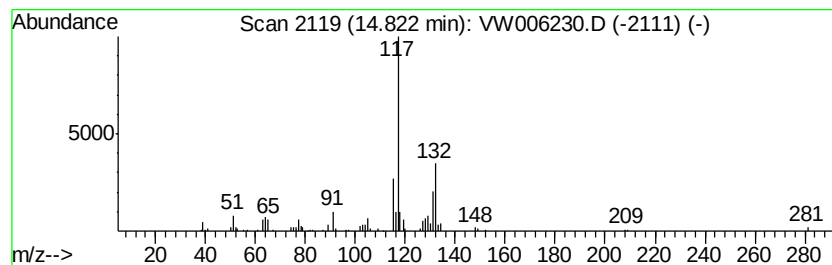
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	16.77 ug/l	48897	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
3		2,4-Dimethylstyrene	132	C10H12	002234-20-0	86
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	83
5		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW101918\
Data File : VW006230.D
Acq On : 19 Oct 2018 17:29
Operator : VA/AP
Sample : J5605-01
Misc : 5.28G/5ML/MSVOA_W/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
TP-A

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-ethyl-...	13.11	33.1	ug/l	96675	4	13.57	145814	50.0
Benzene, 1,2-diet...	13.72	9.8	ug/l	28644	4	13.57	145814	50.0
Benzene, cyclopro...	13.77	21.4	ug/l	62478	4	13.57	145814	50.0
Benzene, 4-ethyl-...	14.02	13.5	ug/l	39275	4	13.57	145814	50.0
Benzene, 2-ethyl-...	14.05	8.9	ug/l	25998	4	13.57	145814	50.0
o-Cymene	14.10	19.1	ug/l	55591	4	13.57	145814	50.0
Benzene, 1-etheny...	14.21	19.5	ug/l	56856	4	13.57	145814	50.0
Benzene, 1,2,4,5-...	14.43	11.2	ug/l	32622	4	13.57	145814	50.0
Benzene, 1,2,3,4-...	14.47	14.4	ug/l	42147	4	13.57	145814	50.0
Benzene, 1-etheny...	14.82	16.8	ug/l	48897	4	13.57	145814	50.0