

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW082220\
 Data File : VW016301.D
 Acq On : 21 Aug 2020 13:13
 Operator : SY/VA
 Sample : L3504-07
 Misc : 7.81G/5ML/MSVOA W/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 EO-01-081920

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\82W081420S.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	4.916	482	494	514	rBV2	44130	148806	1.89%	0.130%
2	7.885	971	981	986	rBV	133192	312208	3.96%	0.272%
3	7.952	986	992	1002	rVB	250083	547194	6.93%	0.476%
4	8.311	1041	1051	1063	rBV	116825	266331	3.37%	0.232%
5	8.848	1130	1139	1151	rBV	358566	719455	9.11%	0.626%
6	9.335	1207	1219	1232	rVB3	43437	95965	1.22%	0.084%
7	10.323	1373	1381	1397	rBV	566613	1083019	13.72%	0.943%
8	10.506	1404	1411	1419	rVB3	60582	119385	1.51%	0.104%
9	11.183	1517	1522	1525	rVV	94163	162902	2.06%	0.142%
10	11.231	1525	1530	1536	rVB	121835	217650	2.76%	0.189%
11	11.341	1542	1548	1551	rBV4	52246	89745	1.14%	0.078%
12	11.414	1551	1560	1571	rVB4	134580	418673	5.30%	0.364%
13	11.512	1571	1576	1581	rBV	150516	245115	3.11%	0.213%
14	11.634	1589	1596	1606	rBV	442443	750783	9.51%	0.654%
15	11.731	1606	1612	1616	rBV2	169762	296179	3.75%	0.258%
16	11.841	1620	1630	1634	rBV	498263	969220	12.28%	0.844%
17	11.914	1639	1642	1648	rVB3	142373	259684	3.29%	0.226%
18	11.975	1648	1652	1659	rBV3	37268	82326	1.04%	0.072%
19	12.061	1659	1666	1671	rVB3	59433	128066	1.62%	0.111%
20	12.164	1671	1683	1691	rBV2	958453	2306504	29.22%	2.008%
21	12.256	1691	1698	1702	rVB	435825	715608	9.07%	0.623%
22	12.372	1714	1717	1724	rVB4	299231	648512	8.22%	0.565%
23	12.463	1727	1732	1736	rBV3	195576	407791	5.17%	0.355%
24	12.512	1737	1740	1741	rBV2	84438	100147	1.27%	0.087%
25	12.542	1742	1745	1748	rBV	239896	273742	3.47%	0.238%
26	12.621	1753	1758	1761	rBV	416348	706330	8.95%	0.615%
27	12.652	1761	1763	1767	rVB	352617	395448	5.01%	0.344%
28	12.701	1767	1771	1775	rVB5	104722	150537	1.91%	0.131%
29	12.786	1779	1785	1792	rVB3	486163	1231220	15.60%	1.072%
30	12.865	1792	1798	1801	rBV3	461127	890477	11.28%	0.775%
31	12.932	1801	1809	1816	rVV2	3322476	7392702	93.66%	6.436%
32	12.993	1816	1819	1824	rVB4	343920	515589	6.53%	0.449%
33	13.042	1824	1827	1829	rBV4	65215	90222	1.14%	0.079%
34	13.103	1829	1837	1843	rBV	1598759	3347716	42.41%	2.914%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	13.164	1843	1847	1854	rVB2	1187388	1874048	23.74%	1.631%
36	13.256	1860	1862	1865	rVB2	164264	177937	2.25%	0.155%
37	13.292	1865	1868	1873	rVB	231697	309367	3.92%	0.269%
38	13.347	1873	1877	1880	rBV	366598	582791	7.38%	0.507%
39	13.384	1880	1883	1887	rVB2	390385	512550	6.49%	0.446%
40	13.445	1887	1893	1897	rBV2	1480388	2697330	34.17%	2.348%
41	13.487	1897	1900	1904	rVV2	775536	1369563	17.35%	1.192%
42	13.554	1908	1911	1912	rVV	695330	764339	9.68%	0.665%
43	13.591	1912	1917	1921	rVV	3287421	5293851	67.07%	4.609%
44	13.627	1921	1923	1930	rVB2	483536	599276	7.59%	0.522%
45	13.719	1930	1938	1940	rBV5	594554	1297949	16.44%	1.130%
46	13.755	1940	1944	1947	rVV2	1788174	3040369	38.52%	2.647%
47	13.792	1947	1950	1959	rVB3	1819162	3368568	42.68%	2.933%
48	13.877	1959	1964	1969	rBV	3895168	5976692	75.72%	5.203%
49	13.951	1969	1976	1983	rVV	1189734	2696224	34.16%	2.347%
50	14.042	1983	1991	1995	rVV3	915340	2399724	30.40%	2.089%
51	14.097	1995	2000	2004	rVV	1784646	2987084	37.84%	2.600%
52	14.140	2004	2007	2013	rVV2	426910	901456	11.42%	0.785%
53	14.207	2013	2018	2023	rVV2	2101282	3518735	44.58%	3.063%
54	14.261	2023	2027	2034	rVV6	369237	953788	12.08%	0.830%
55	14.341	2034	2040	2044	rVV2	1148579	2298415	29.12%	2.001%
56	14.389	2044	2048	2051	rVV3	1136447	2097972	26.58%	1.826%
57	14.426	2051	2054	2056	rVV2	1013111	1538514	19.49%	1.339%
58	14.457	2056	2059	2063	rVV	1859642	2758651	34.95%	2.402%
59	14.499	2063	2066	2070	rVV2	1005901	1387545	17.58%	1.208%
60	14.548	2070	2074	2081	rVB4	721996	1393180	17.65%	1.213%
61	14.639	2086	2089	2094	rVB2	501238	728146	9.22%	0.634%
62	14.694	2094	2098	2099	rBV2	605028	833758	10.56%	0.726%
63	14.725	2099	2103	2109	rVB2	2524150	3913092	49.57%	3.407%
64	14.798	2109	2115	2121	rBV2	3286518	7893343	100.00%	6.872%
65	14.963	2137	2142	2148	rVB	1460000	2462383	31.20%	2.144%
66	15.030	2150	2153	2154	rBV2	164843	187242	2.37%	0.163%
67	15.066	2155	2159	2163	rBV2	778794	1323184	16.76%	1.152%
68	15.176	2171	2177	2185	rBV3	1033243	2420123	30.66%	2.107%
69	15.249	2186	2189	2196	rVB6	485409	1059458	13.42%	0.922%
70	15.328	2196	2202	2211	rVB5	936515	1898877	24.06%	1.653%
71	15.426	2212	2218	2222	rBV	799674	1506779	19.09%	1.312%

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Integration Parameters: RTEINT.P

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72	15.475	2222	2226	2227	rBV2	342332	427484	5.42%	0.372%
73	15.554	2236	2239	2248	rVB2	1075755	1941594	24.60%	1.690%
74	15.658	2248	2256	2263	rBV4	483724	1313108	16.64%	1.143%
75	15.731	2264	2268	2274	rBV3	338297	586639	7.43%	0.511%
76	15.822	2279	2283	2286	rVV3	342213	716107	9.07%	0.623%
77	15.865	2286	2290	2297	rVB	1055549	1843926	23.36%	1.605%
78	16.182	2334	2342	2354	rVB3	706517	1846739	23.40%	1.608%
79	16.285	2356	2359	2365	rVB2	237963	399411	5.06%	0.348%
80	16.377	2368	2374	2380	rVV	550984	1202632	15.24%	1.047%
81	16.444	2380	2385	2389	rVV2	246165	519070	6.58%	0.452%
82	16.493	2390	2393	2398	rVB2	229435	382477	4.85%	0.333%
83	16.651	2411	2419	2424	rBV8	185100	583350	7.39%	0.508%

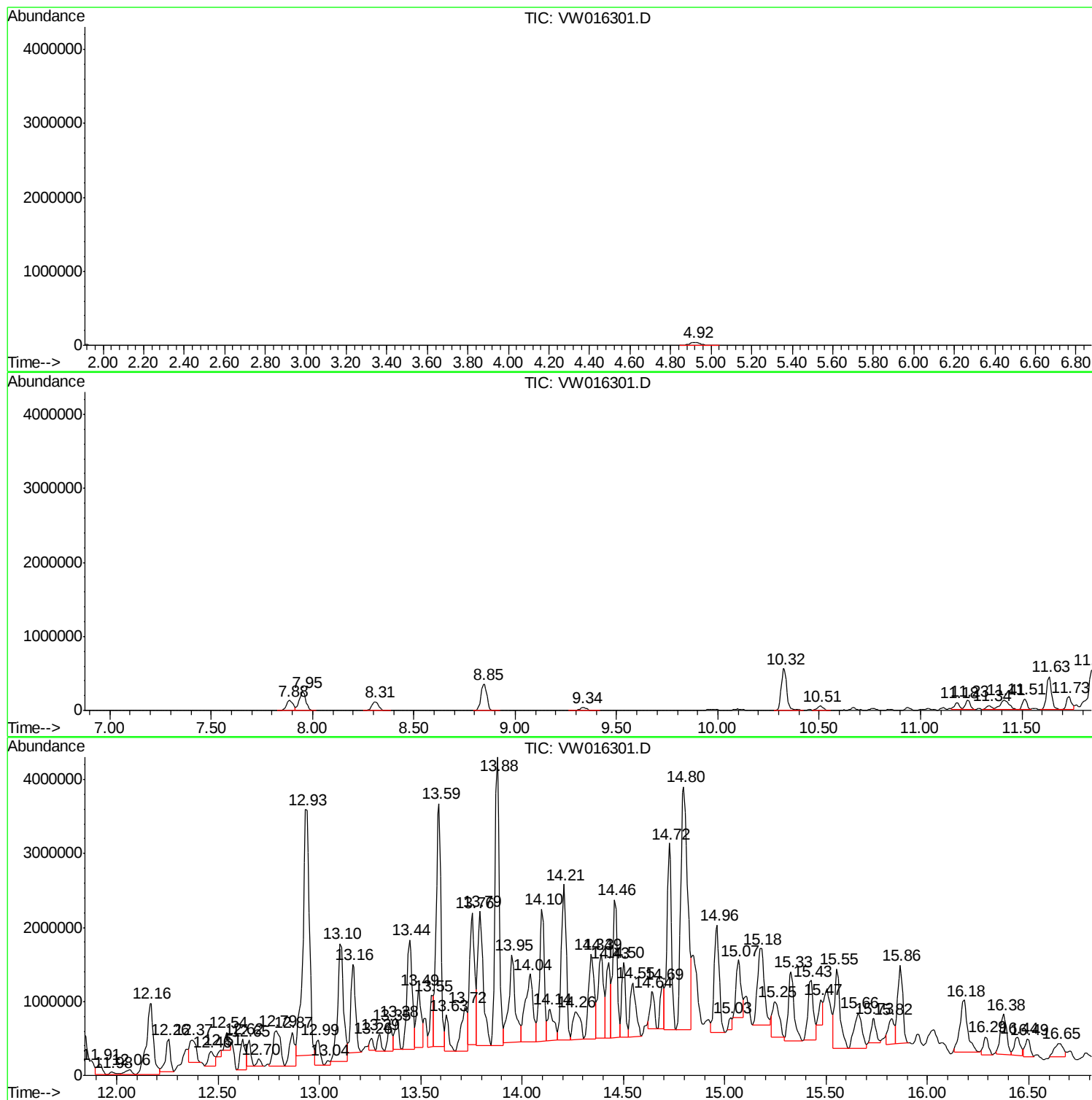
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W081420S.M
Quant Title : SW846 8260

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



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EO-01-081920

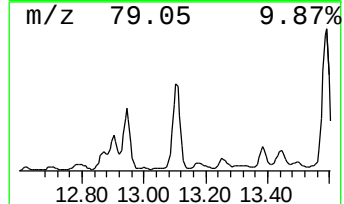
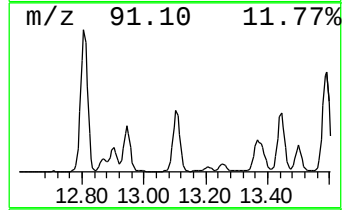
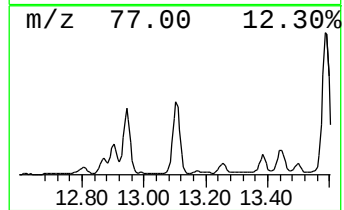
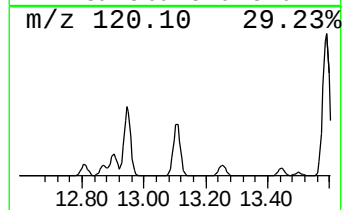
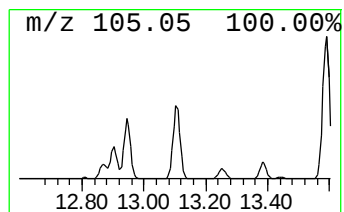
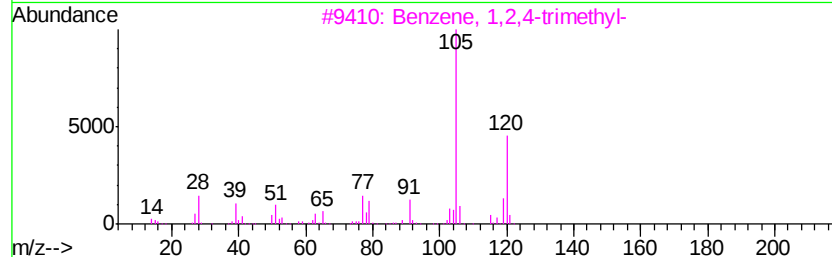
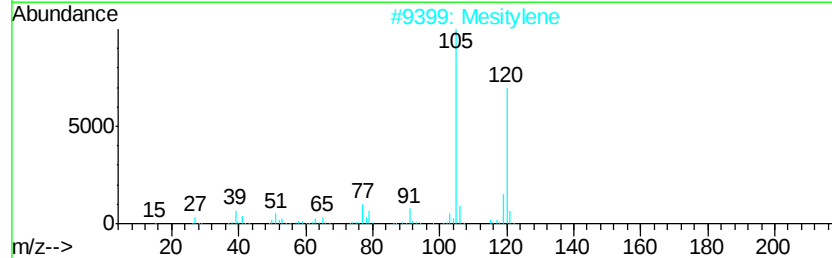
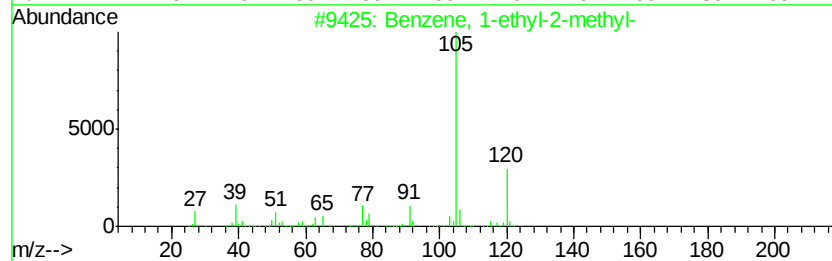
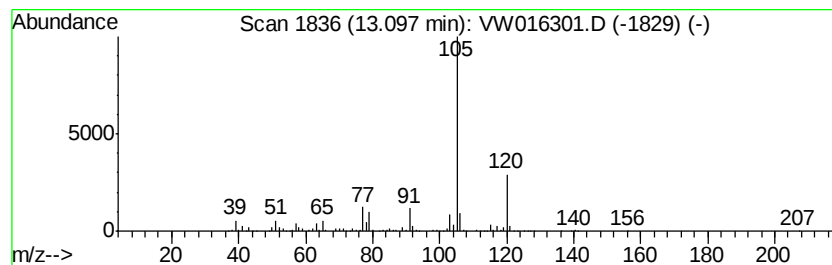
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W081420S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.10	218.99 ug/l	3347720	1,4-Dichlorobenzene-d4	13.56

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



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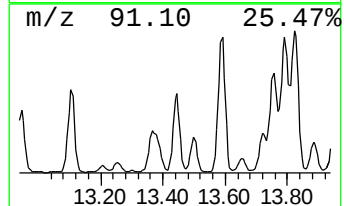
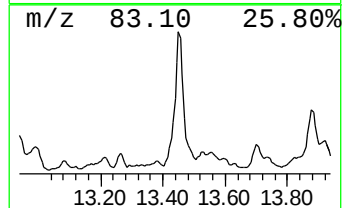
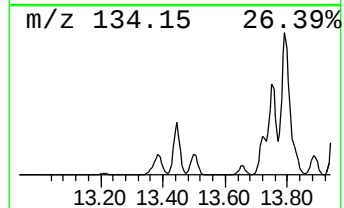
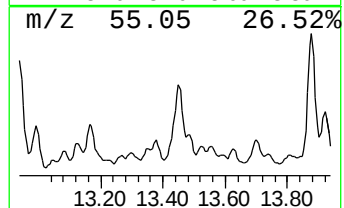
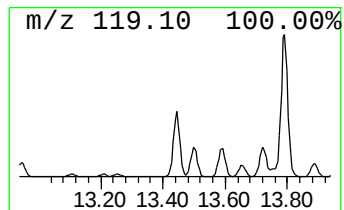
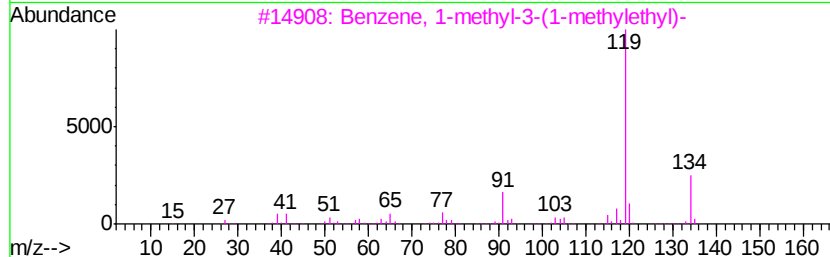
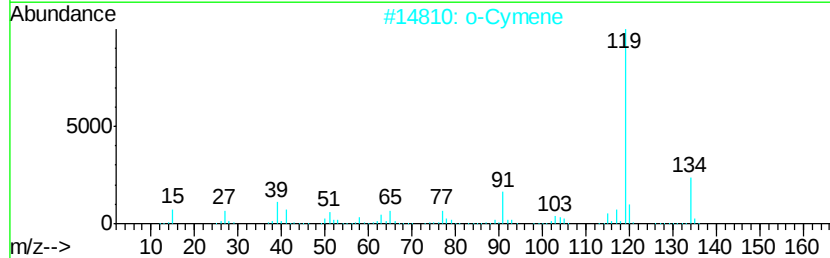
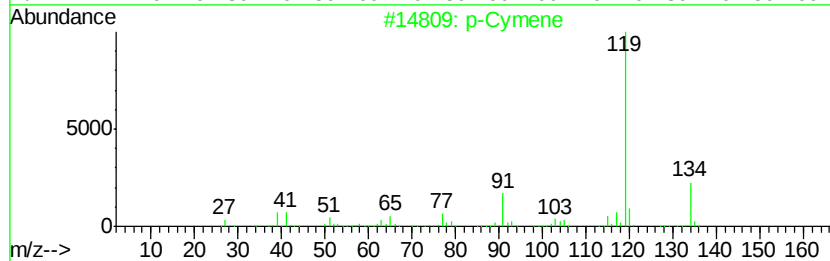
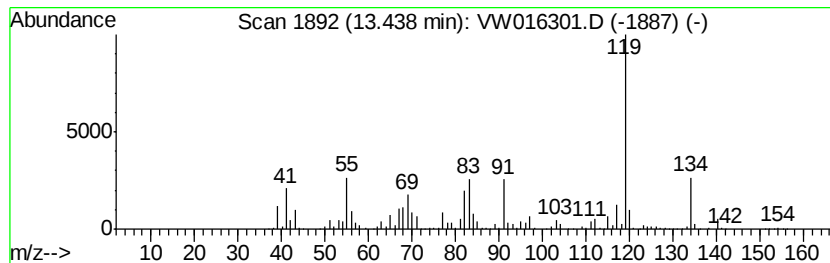
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 o-Cymene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.44	176.45 ug/l	2697330	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	p-Cymene	134	C10H14	000099-87-6	95
2		o-Cymene	134	C10H14	000527-84-4	95
3		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	93
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	70
5		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	70



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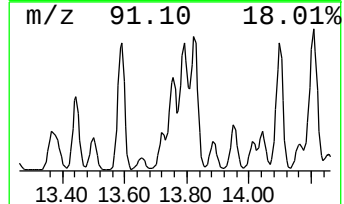
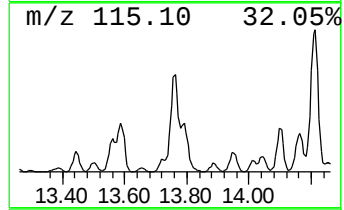
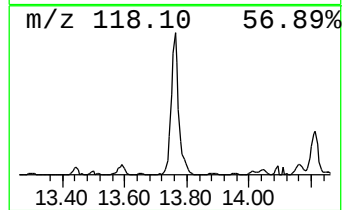
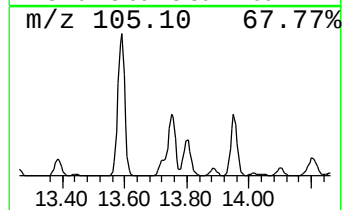
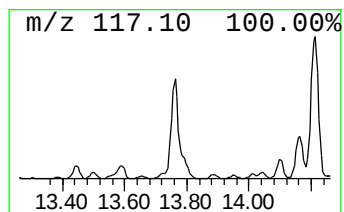
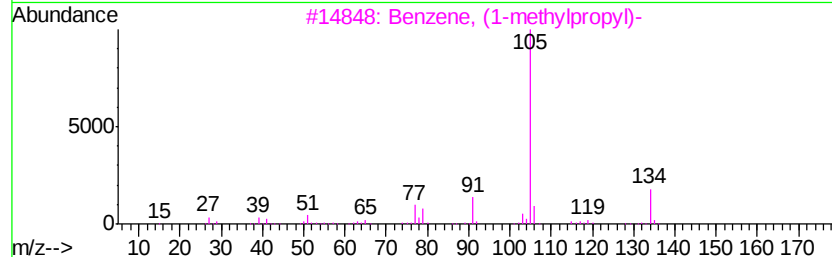
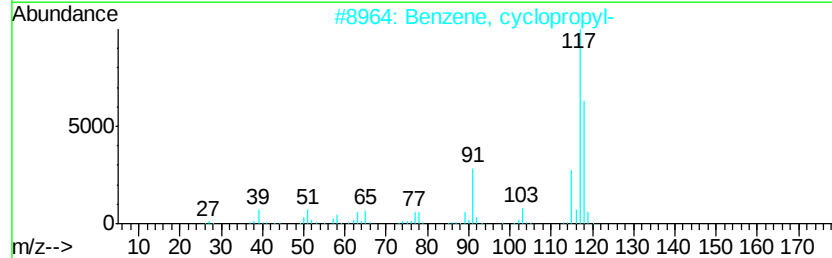
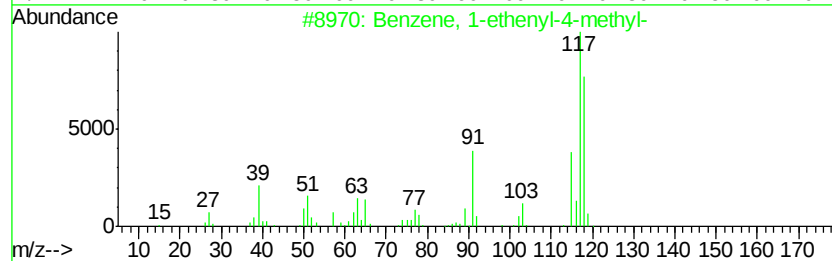
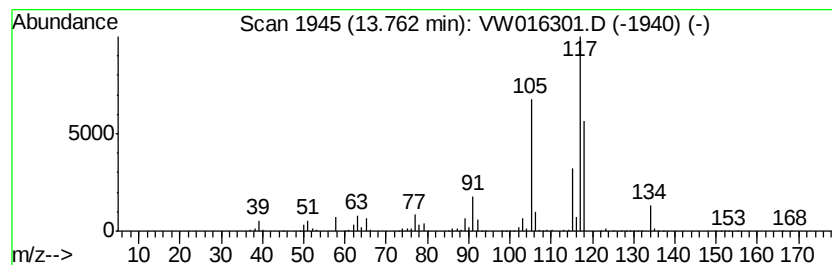
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W081420S.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-ethenyl-4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	198.89 ug/l	3040370	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	50
2		Benzene, cyclopropyl-	118	C9H10	000873-49-4	50
3		Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	46
4		Benzene, 1-propenyl-	118	C9H10	000637-50-3	45
5		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	43



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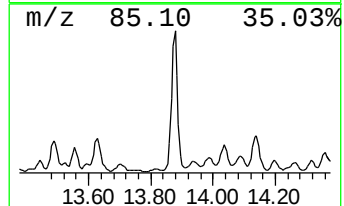
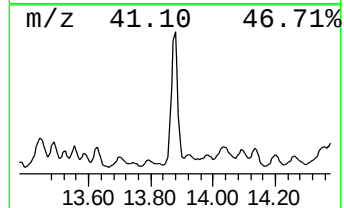
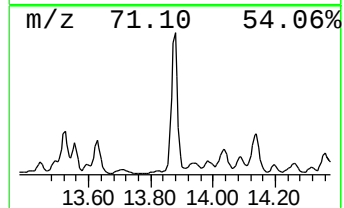
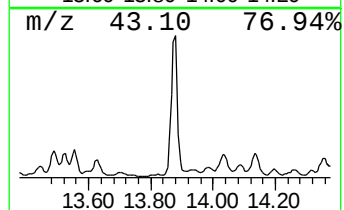
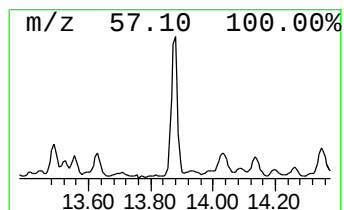
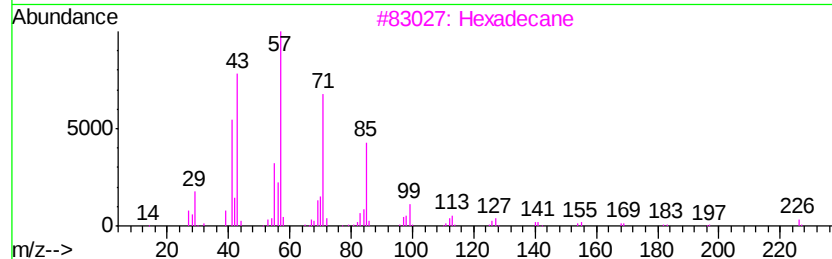
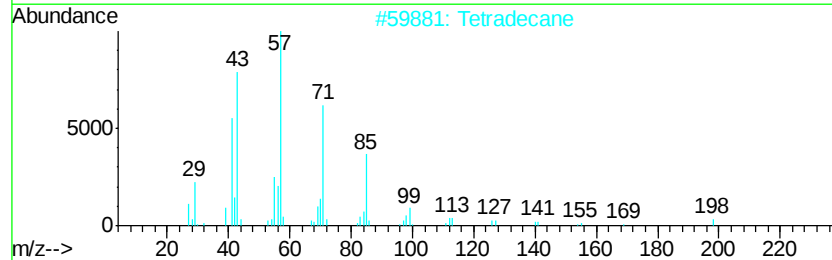
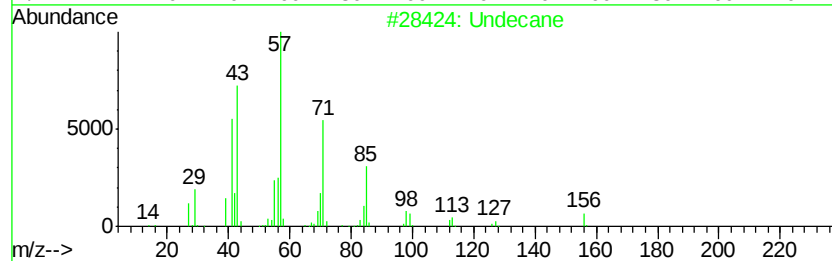
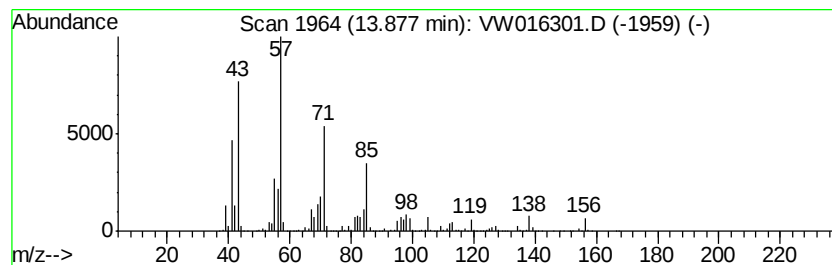
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Quant Title : SW846 8260

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

Peak Number 4 Undecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.88	390.97 ug/l	5976690	1,4-Dichlorobenzene-d4	13.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane		156	C11H24	001120-21-4	96
2	Tetradecane		198	C14H30	000629-59-4	80
3	Hexadecane		226	C16H34	000544-76-3	80
4	Decane		142	C10H22	000124-18-5	74
5	Decane, 2,3,5,8-tetramethyl-		198	C14H30	192823-15-7	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW082220\
Data File : VW016301.D
Acq On : 21 Aug 2020 13:13
Operator : SY/VA
Sample : L3504-07
Misc : 7.81G/5ML/MSVOA W/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EO-01-081920

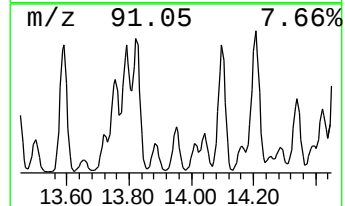
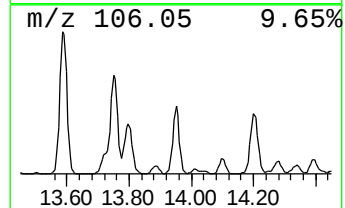
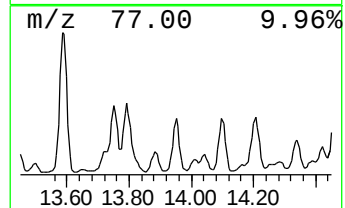
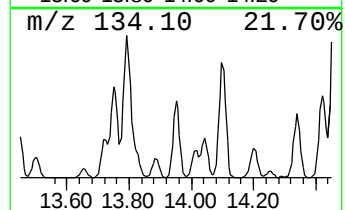
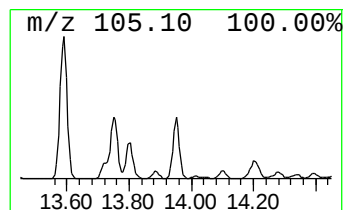
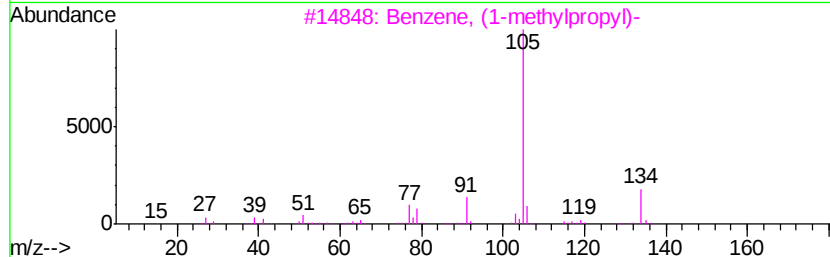
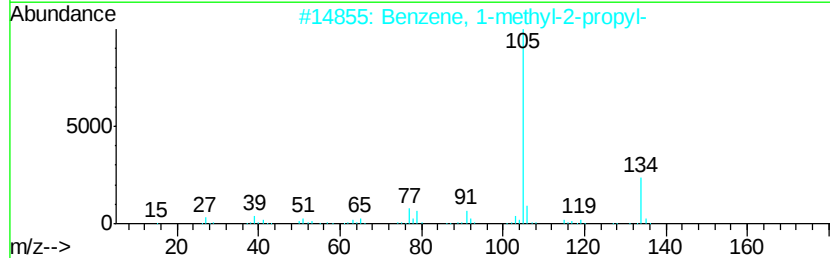
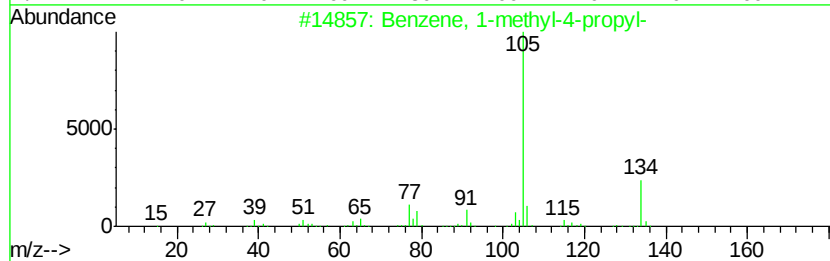
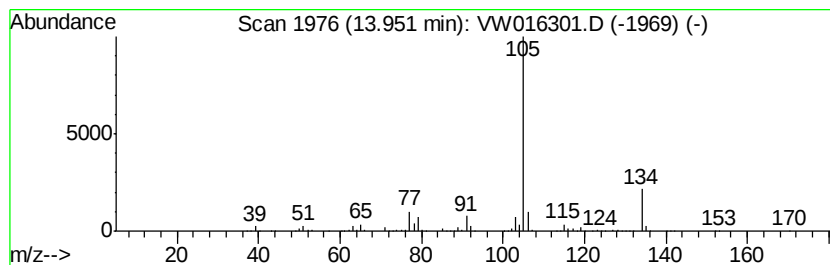
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-methyl-4-propyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	176.38 ug/l	2696220	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	94
2		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	90
3		Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	86
4		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	80
5		Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW082220\
Data File : VW016301.D
Acq On : 21 Aug 2020 13:13
Operator : SY/VA
Sample : L3504-07
Misc : 7.81G/5ML/MSVOA W/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EO-01-081920

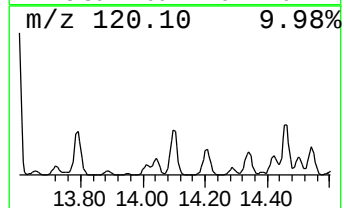
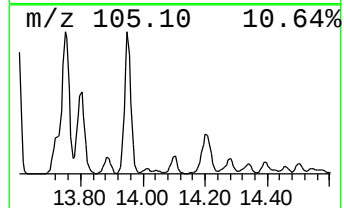
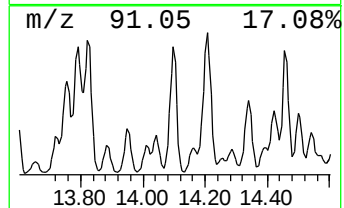
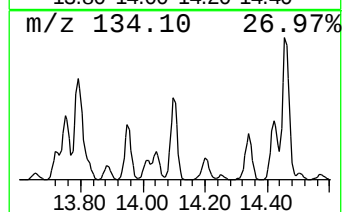
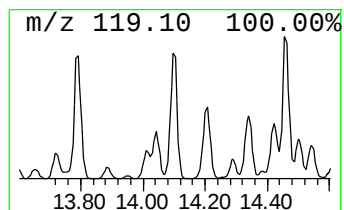
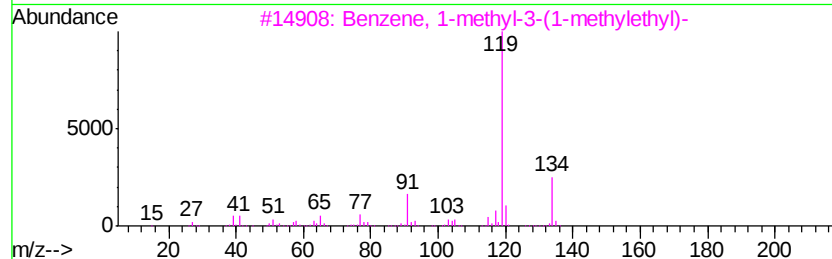
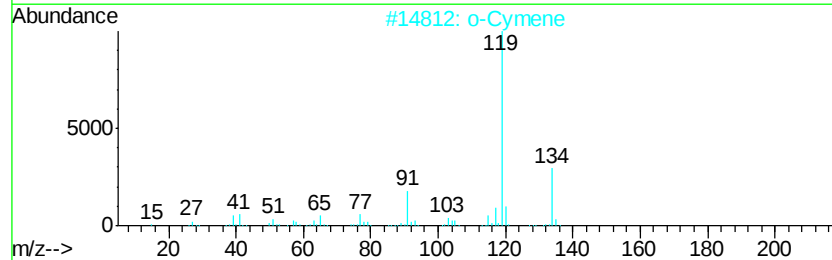
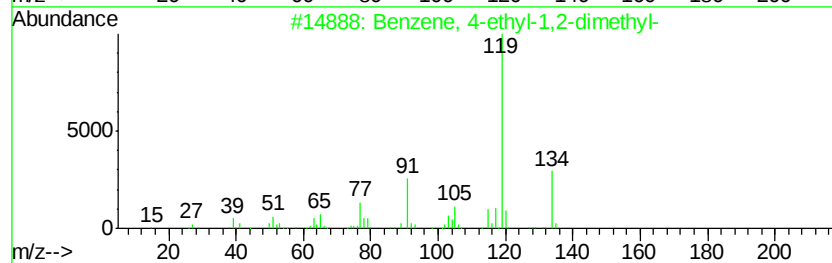
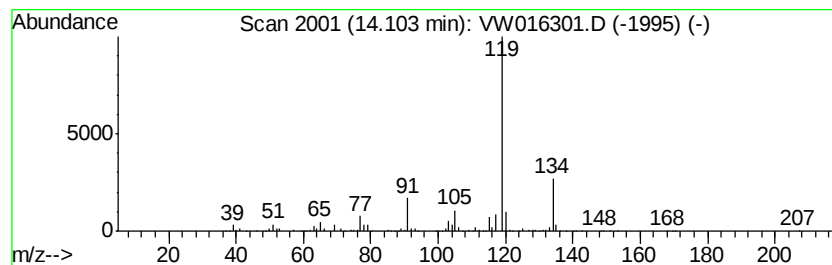
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.10	195.40 ug/l	2987080	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
2		o-Cymene	134	C10H14	000527-84-4	94
3		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	94
4		p-Cymene	134	C10H14	000099-87-6	94
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW082220\
Data File : VW016301.D
Acq On : 21 Aug 2020 13:13
Operator : SY/VA
Sample : L3504-07
Misc : 7.81G/5ML/MSVOA W/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EO-01-081920

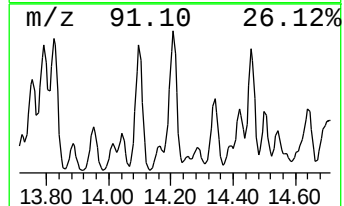
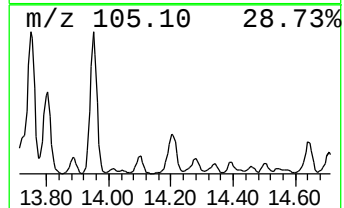
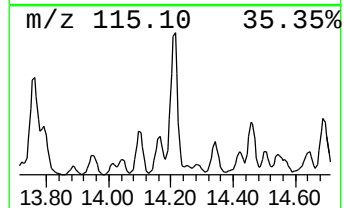
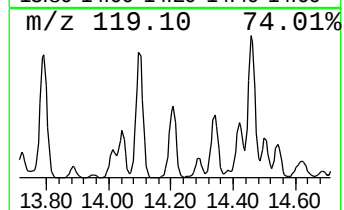
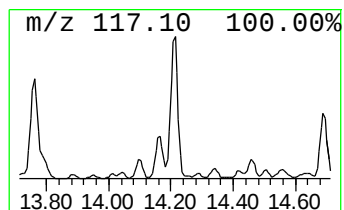
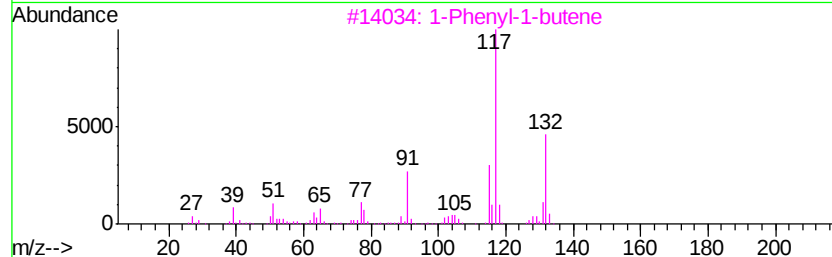
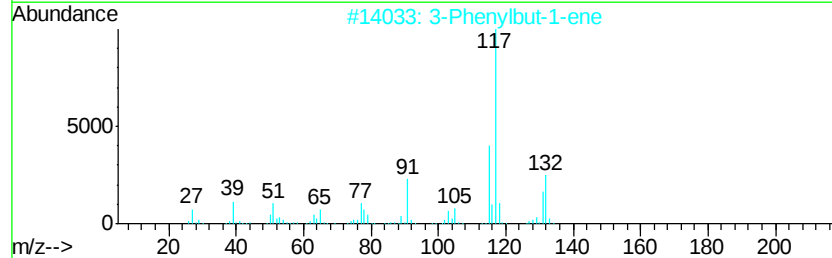
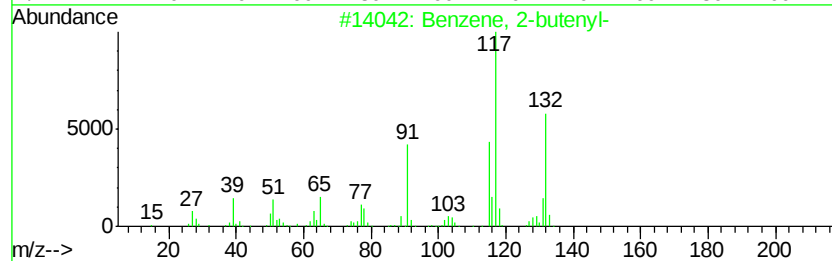
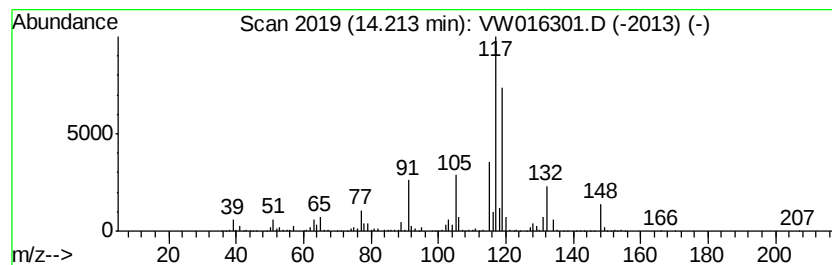
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 2-butenyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	230.18 ug/l	3518740	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	89
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	50
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	50
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	50
5		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	45



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW082220\
Data File : VW016301.D
Acq On : 21 Aug 2020 13:13
Operator : SY/VA
Sample : L3504-07
Misc : 7.81G/5ML/MSVOA W/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EO-01-081920

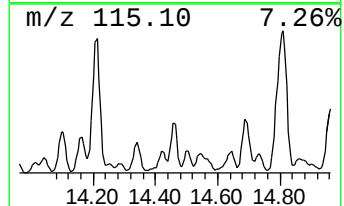
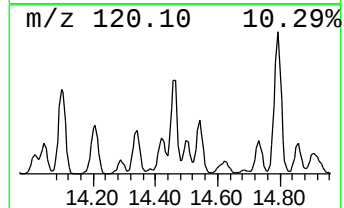
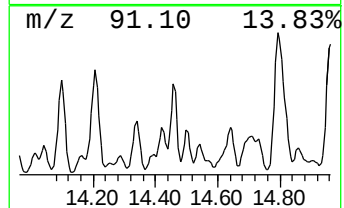
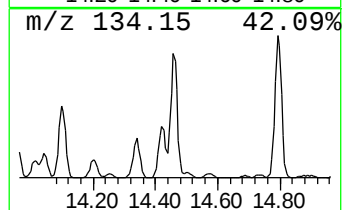
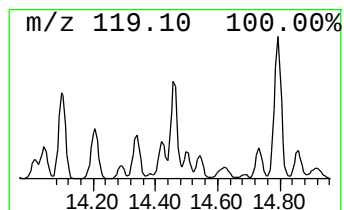
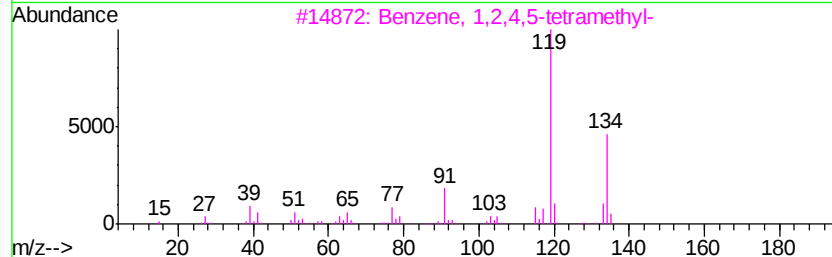
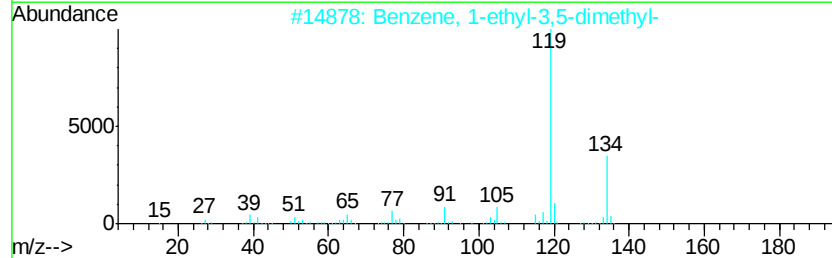
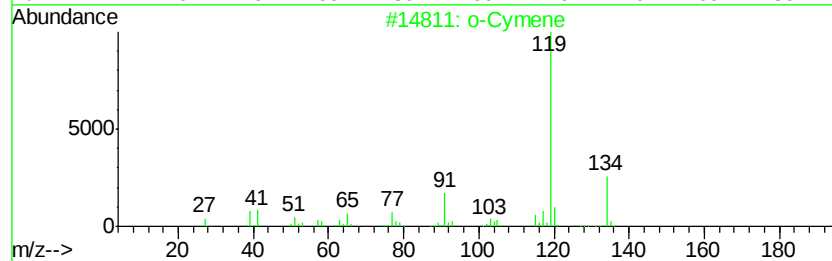
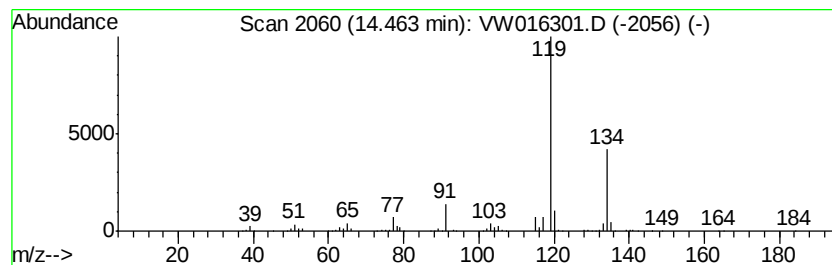
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.46	180.46 ug/l	2758650	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	95
2		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW082220\
Data File : VW016301.D
Acq On : 21 Aug 2020 13:13
Operator : SY/VA
Sample : L3504-07
Misc : 7.81G/5ML/MSVOA W/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EO-01-081920

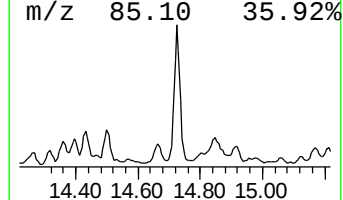
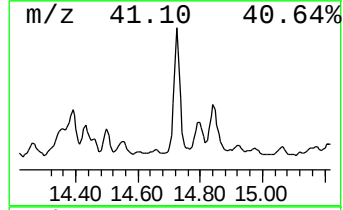
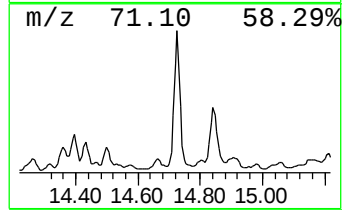
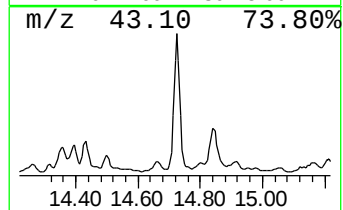
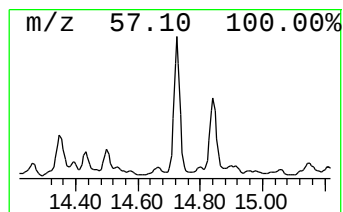
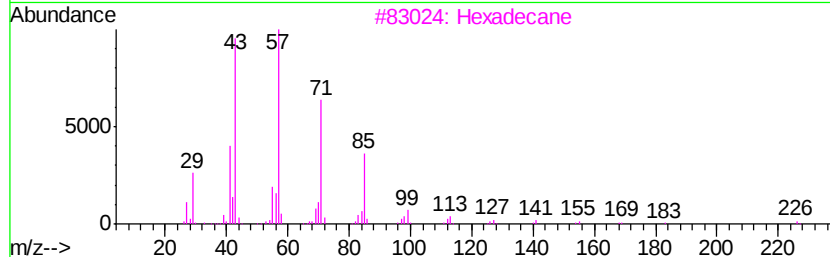
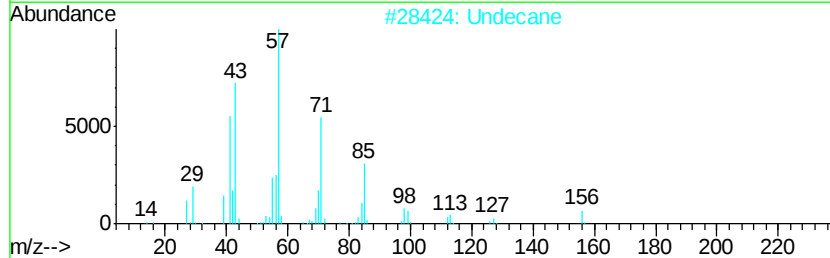
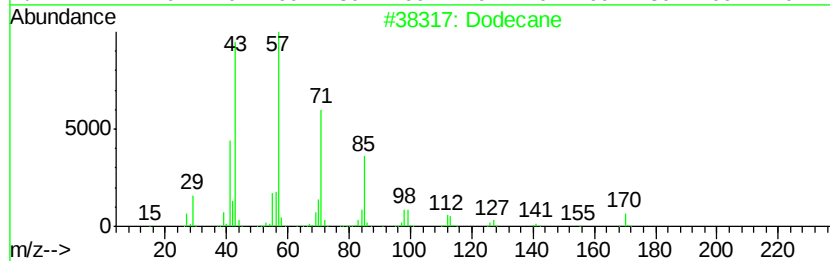
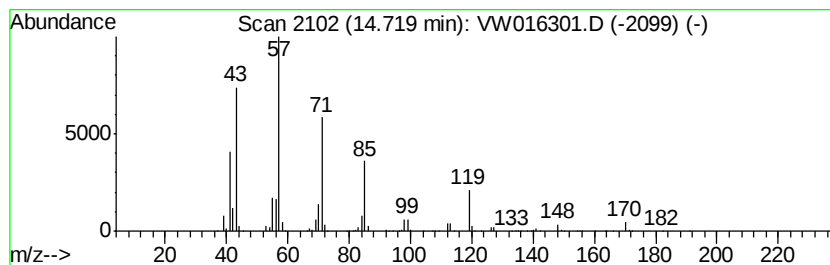
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Quant Title : SW846 8260

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

Peak Number 9 Dodecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.72	255.98 ug/l	3913090	1,4-Dichlorobenzene-d4	13.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane		170	C12H26	000112-40-3	96
2	Undecane		156	C11H24	001120-21-4	70
3	Hexadecane		226	C16H34	000544-76-3	58
4	Decane, 2,3,6-trimethyl-		184	C13H28	062238-12-4	52
5	Decane, 2,3,5-trimethyl-		184	C13H28	062238-11-3	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW082220\
Data File : VW016301.D
Acq On : 21 Aug 2020 13:13
Operator : SY/VA
Sample : L3504-07
Misc : 7.81G/5ML/MSVOA W/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EO-01-081920

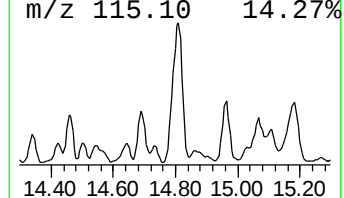
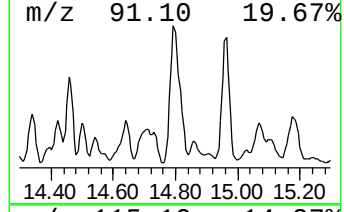
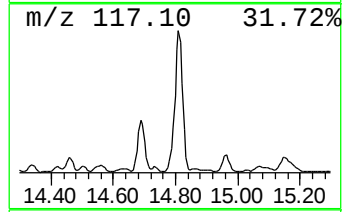
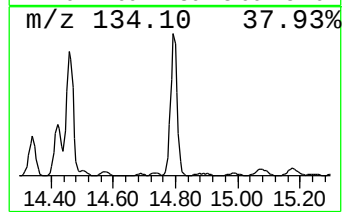
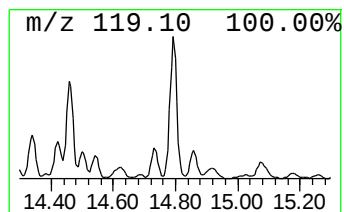
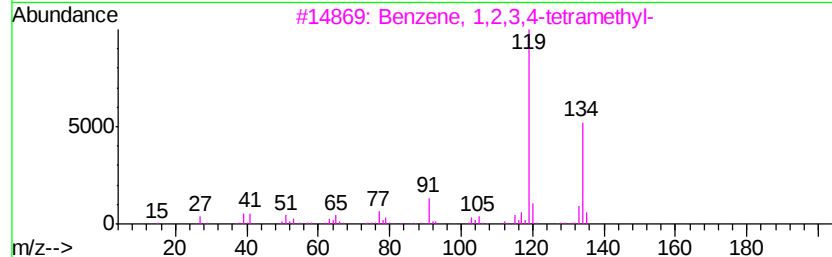
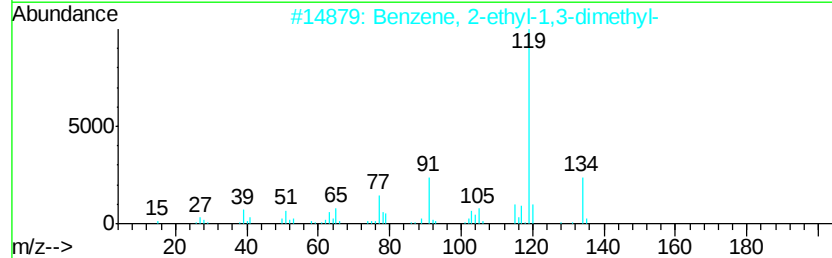
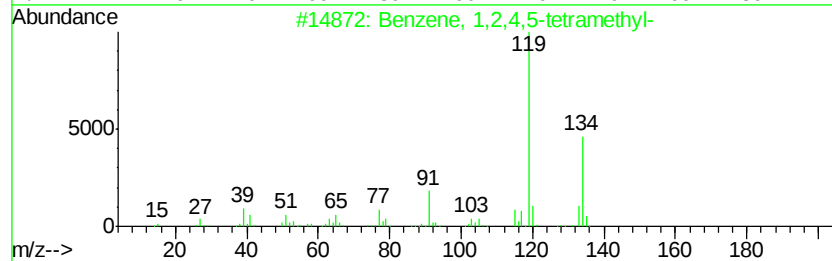
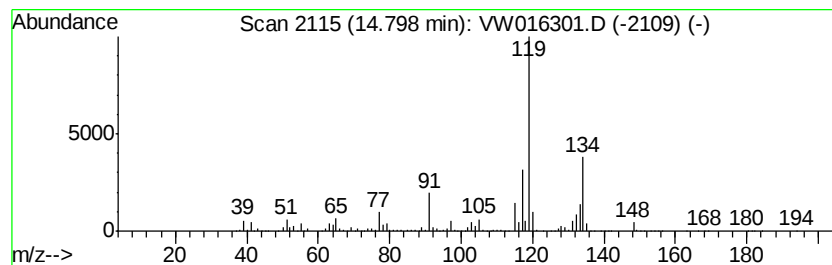
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.80	516.35 ug/l	7893340	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	93
2		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	92
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	81
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	81
5		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW082220\
Data File : VW016301.D
Acq On : 21 Aug 2020 13:13
Operator : SY/VA
Sample : L3504-07
Misc : 7.81G/5ML/MSVOA_W/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EO-01-081920

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W081420S.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-...	13.10	219.0	ug/l	3347720	4	13.56	764339	50.0
o-Cymene	13.44	176.4	ug/l	2697330	4	13.56	764339	50.0
Benzene, 1-etheny...	13.76	198.9	ug/l	3040370	4	13.56	764339	50.0
Undecane	13.88	391.0	ug/l	5976690	4	13.56	764339	50.0
Benzene, 1-methyl...	13.95	176.4	ug/l	2696220	4	13.56	764339	50.0
Benzene, 4-ethyl-...	14.10	195.4	ug/l	2987080	4	13.56	764339	50.0
Benzene, 2-butenyl-	14.21	230.2	ug/l	3518740	4	13.56	764339	50.0
Benzene, 1-ethyl-...	14.46	180.5	ug/l	2758650	4	13.56	764339	50.0
Dodecane	14.72	256.0	ug/l	3913090	4	13.56	764339	50.0
Benzene, 1,2,4,5-...	14.80	516.4	ug/l	7893340	4	13.56	764339	50.0