

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW011119\
 Data File : VW008219.D
 Acq On : 12 Jan 2019 04:28
 Operator : SY/VA
 Sample : K1119-03
 Misc : 5.14G/5ML/MSVOA W/SOIL
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 DS-B

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\82W010819S.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	3.678	287	291	298	rVB6	3978	8493	1.53%	0.082%
2	4.909	483	493	504	rBV2	13562	41812	7.51%	0.403%
3	7.885	973	981	986	rBV	71402	171259	30.76%	1.652%
4	7.945	986	991	1000	rVB	123753	273623	49.14%	2.640%
5	8.305	1039	1050	1060	rBV	69678	153283	27.53%	1.479%
6	8.842	1127	1138	1148	rVB	192942	384836	69.12%	3.713%
7	10.323	1373	1381	1388	rBV	320965	556800	100.00%	5.372%
8	10.390	1388	1392	1397	rVB	13016	23496	4.22%	0.227%
9	10.506	1404	1411	1418	rVB5	5137	9802	1.76%	0.095%
10	11.177	1515	1521	1525	rBV6	3943	7127	1.28%	0.069%
11	11.231	1527	1530	1536	rVB6	3849	6207	1.11%	0.060%
12	11.408	1553	1559	1572	rVB9	6574	20317	3.65%	0.196%
13	11.512	1572	1576	1581	rBV3	6604	10404	1.87%	0.100%
14	11.634	1588	1596	1605	rBV	265200	447739	80.41%	4.320%
15	11.731	1605	1612	1620	rVV	24414	47653	8.56%	0.460%
16	11.841	1621	1630	1639	rVV2	111240	229512	41.22%	2.214%
17	11.914	1639	1642	1649	rVB5	8071	14774	2.65%	0.143%
18	12.055	1662	1665	1672	rVB7	2659	5636	1.01%	0.054%
19	12.164	1672	1683	1692	rBV	61452	142045	25.51%	1.370%
20	12.256	1695	1698	1702	rVV2	7664	11261	2.02%	0.109%
21	12.311	1702	1707	1708	rVV4	5256	8523	1.53%	0.082%
22	12.378	1708	1718	1728	rVV8	21477	95467	17.15%	0.921%
23	12.463	1728	1732	1738	rVV	18923	40803	7.33%	0.394%
24	12.548	1742	1746	1748	rVV4	10284	19129	3.44%	0.185%
25	12.621	1752	1758	1767	rVV	215501	355379	63.83%	3.429%
26	12.701	1767	1771	1775	rVB7	5781	10079	1.81%	0.097%
27	12.804	1780	1788	1793	rVV	43319	87440	15.70%	0.844%
28	12.871	1793	1799	1802	rVV	175356	279591	50.21%	2.698%
29	12.902	1802	1804	1807	rVV	69354	108750	19.53%	1.049%
30	12.945	1807	1811	1817	rVV2	134218	252049	45.27%	2.432%
31	12.993	1817	1819	1824	rVB4	13557	18642	3.35%	0.180%
32	13.109	1830	1838	1844	rBV	90260	159886	28.72%	1.543%
33	13.176	1844	1849	1855	rVB4	22482	45853	8.24%	0.442%
34	13.256	1855	1862	1868	rBV	341929	519152	93.24%	5.009%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	13.384	1878	1883	1887	rVB2	33802	50711	9.11%	0.489%
36	13.451	1887	1894	1898	rBV3	63326	110998	19.93%	1.071%
37	13.499	1898	1902	1907	rBV3	18368	29541	5.31%	0.285%
38	13.560	1907	1912	1916	rBV	271231	480280	86.26%	4.634%
39	13.627	1922	1923	1928	rVB3	8482	9014	1.62%	0.087%
40	13.688	1930	1933	1934	rVV2	8777	8452	1.52%	0.082%
41	13.725	1935	1939	1940	rVV	43721	56241	10.10%	0.543%
42	13.755	1940	1944	1948	rVV2	190737	331510	59.54%	3.199%
43	13.798	1948	1951	1960	rVB3	144050	291068	52.28%	2.808%
44	13.883	1960	1965	1971	rBV2	104926	173946	31.24%	1.678%
45	13.957	1972	1977	1982	rVB	62984	91414	16.42%	0.882%
46	14.018	1982	1987	1989	rBV	84814	126144	22.66%	1.217%
47	14.042	1989	1991	1996	rVB	81781	112773	20.25%	1.088%
48	14.103	1996	2001	2006	rVB	131422	194869	35.00%	1.880%
49	14.164	2006	2011	2014	rBV2	24466	40974	7.36%	0.395%
50	14.213	2014	2019	2024	rVB2	153969	246827	44.33%	2.381%
51	14.292	2024	2032	2036	rBV5	33175	73075	13.12%	0.705%
52	14.341	2036	2040	2044	rBV2	42059	66145	11.88%	0.638%
53	14.389	2044	2048	2051	rBV3	57751	87622	15.74%	0.845%
54	14.426	2051	2054	2057	rVV	50711	60749	10.91%	0.586%
55	14.463	2057	2060	2064	rVB	93588	117139	21.04%	1.130%
56	14.505	2064	2067	2071	rBV	65806	79024	14.19%	0.762%
57	14.548	2071	2074	2080	rVB4	65506	98571	17.70%	0.951%
58	14.609	2080	2084	2087	rBV4	24988	44031	7.91%	0.425%
59	14.646	2087	2090	2094	rVB	63436	84437	15.16%	0.815%
60	14.694	2094	2098	2102	rBV2	98348	179982	32.32%	1.737%
61	14.737	2102	2105	2110	rVB2	77142	118985	21.37%	1.148%
62	14.816	2110	2118	2123	rBV3	193039	471343	84.65%	4.548%
63	14.865	2123	2126	2133	rVB3	47820	75632	13.58%	0.730%
64	14.963	2137	2142	2149	rVB	121286	198092	35.58%	1.911%
65	15.072	2150	2160	2164	rBV4	84221	200094	35.94%	1.931%
66	15.188	2172	2179	2184	rVB2	84577	194287	34.89%	1.875%
67	15.261	2188	2191	2199	rVB7	14063	29990	5.39%	0.289%
68	15.340	2199	2204	2206	rBV4	25275	42167	7.57%	0.407%
69	15.426	2213	2218	2223	rBV	78937	141630	25.44%	1.366%
70	15.481	2223	2227	2229	rBV2	40606	61635	11.07%	0.595%
71	15.664	2250	2257	2265	rBV3	53820	127370	22.88%	1.229%

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Integration Parameters: RTEINT.P

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72	15.743	2265	2270	2275	rVV3	12822	26593	4.78%	0.257%
73	15.834	2277	2285	2286	rVV6	34716	76361	13.71%	0.737%
74	15.871	2286	2291	2300	rVB	124639	241616	43.39%	2.331%
75	15.962	2301	2306	2310	rBV4	13173	25342	4.55%	0.245%
76	16.035	2312	2318	2324	rBV4	21822	52024	9.34%	0.502%
77	16.188	2335	2343	2350	rBV3	60307	140991	25.32%	1.360%
78	16.389	2366	2376	2381	rBV3	69584	164005	29.45%	1.582%
79	16.450	2382	2386	2398	rVB7	30330	84587	15.19%	0.816%
80	16.663	2412	2421	2425	rBV6	19023	54163	9.73%	0.523%
81	16.712	2425	2429	2435	rVB6	11987	25211	4.53%	0.243%

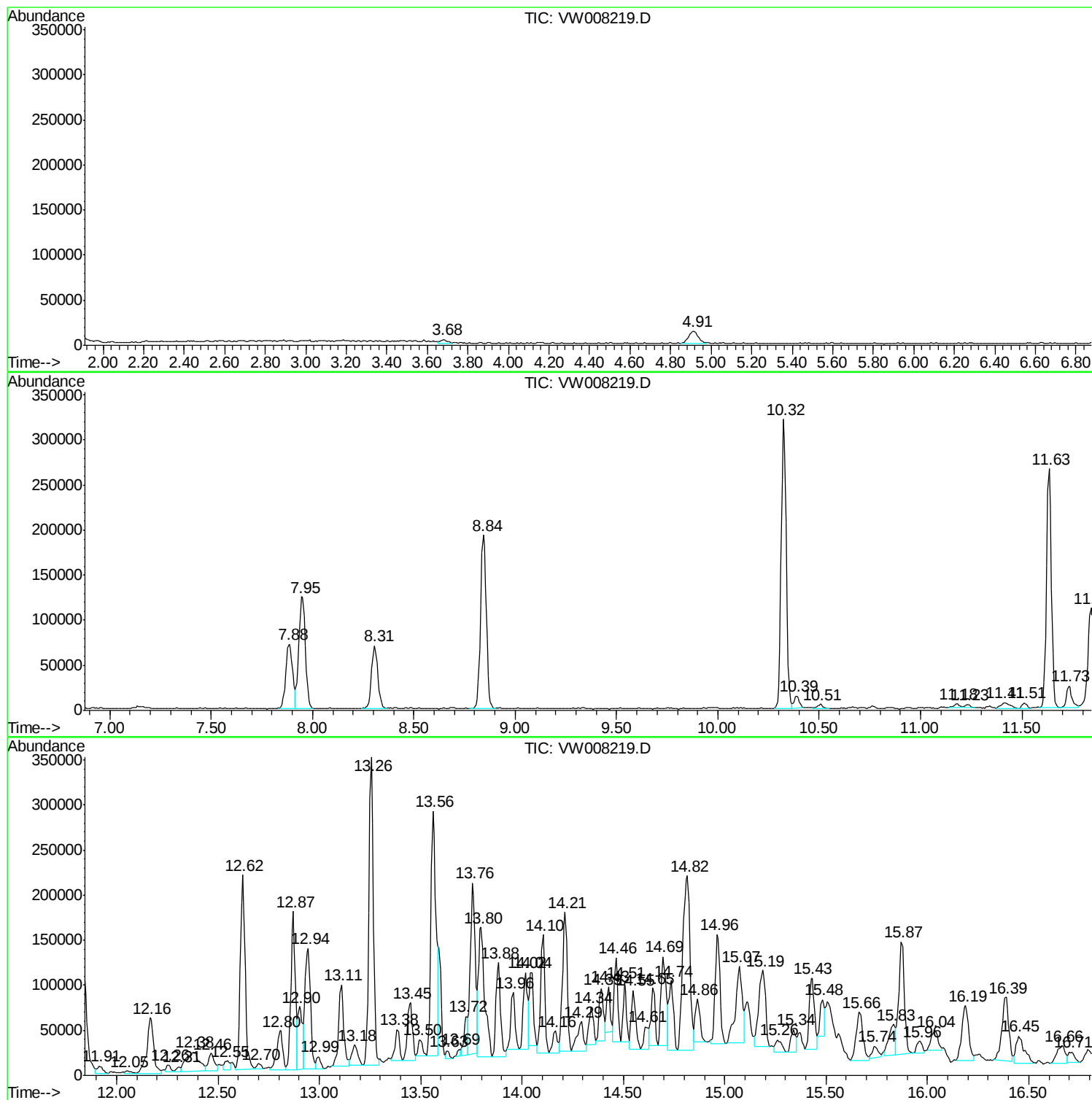
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Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

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



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Operator : SY/VA
Sample : K1119-03
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ALS Vial : 34 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampled :
DS-B

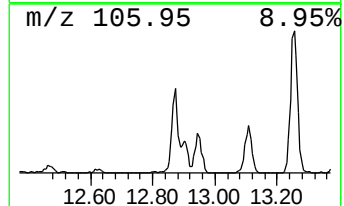
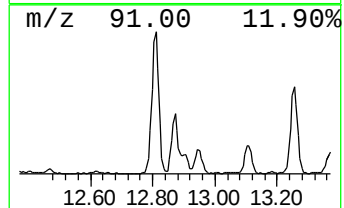
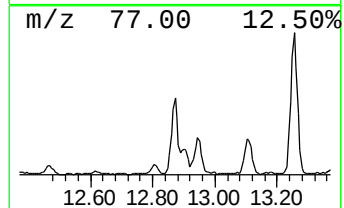
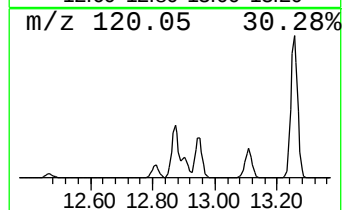
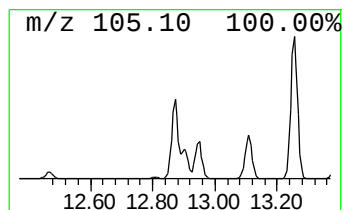
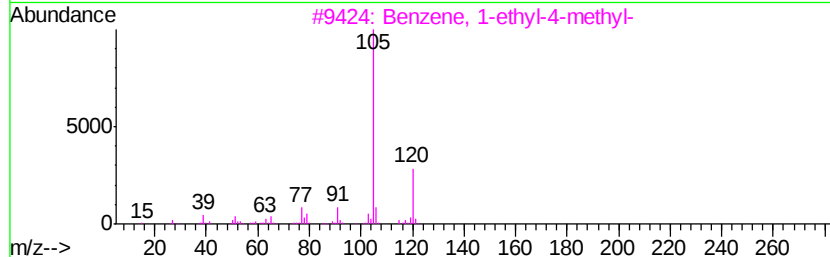
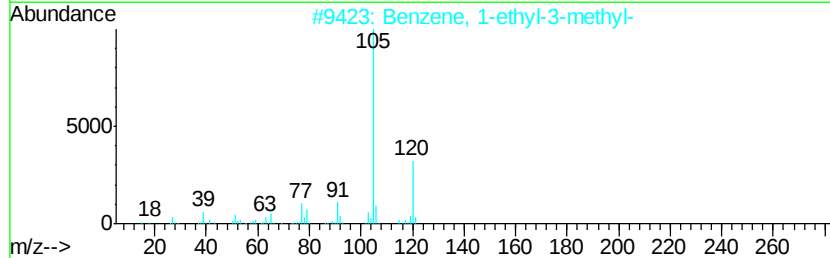
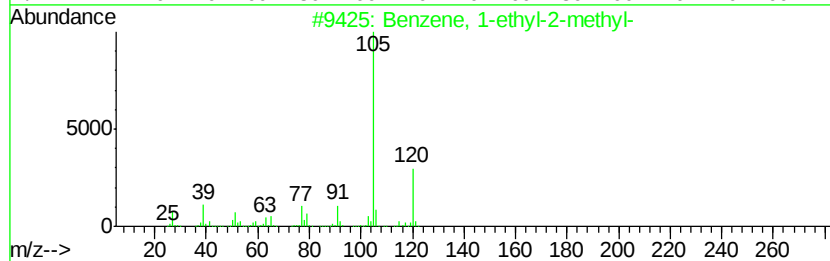
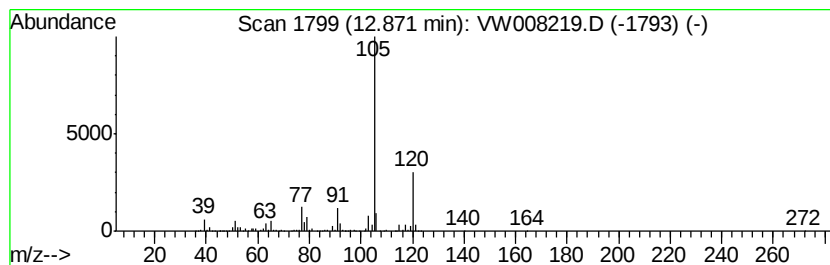
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.87	29.11 ug/l	279591	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		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Mesitylene	120	C9H12	000108-67-8	91



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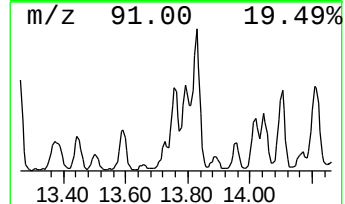
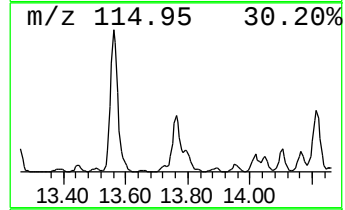
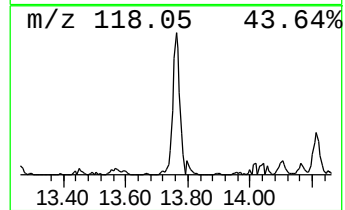
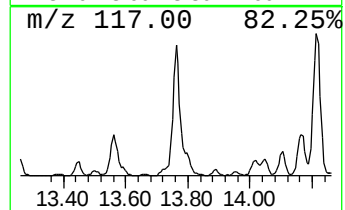
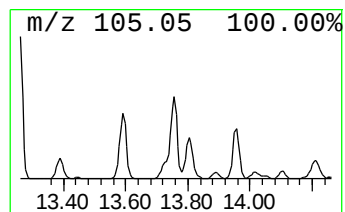
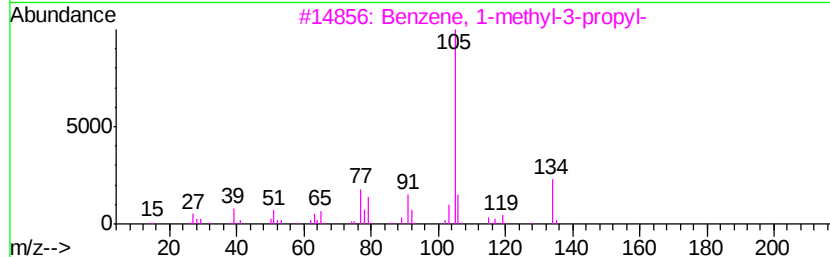
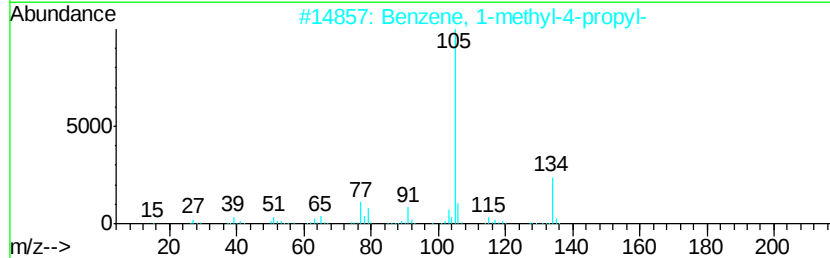
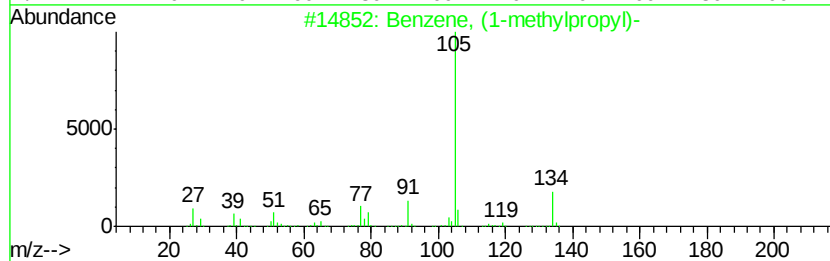
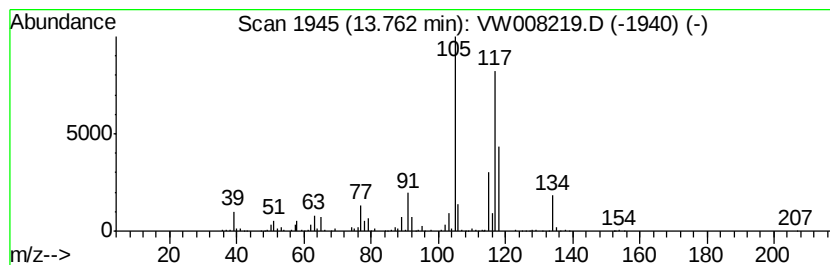
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown13.76 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	34.51 ug/l	331510	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	43
2		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	43
3		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	38
4		Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	35
5		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	35



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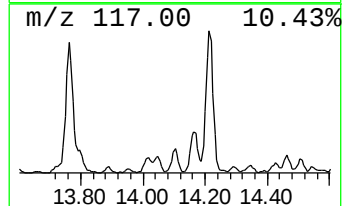
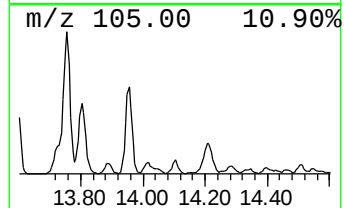
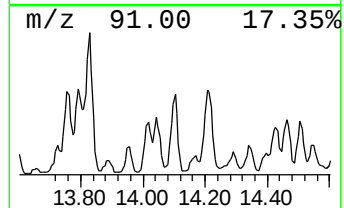
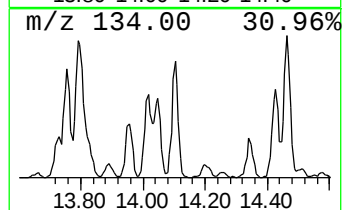
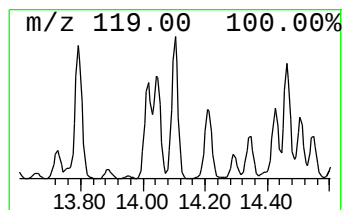
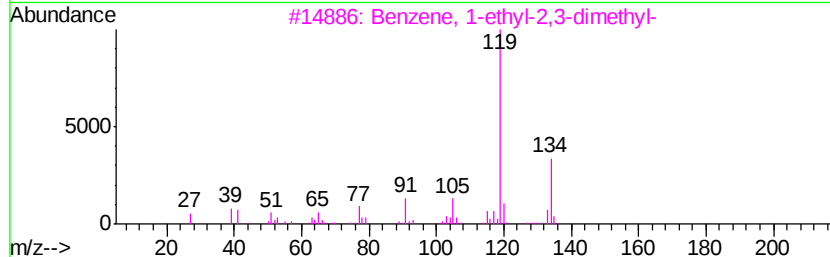
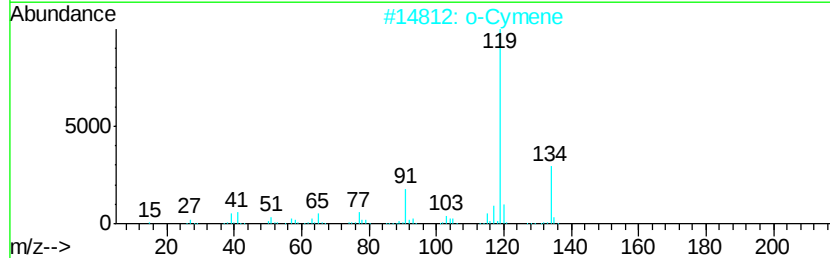
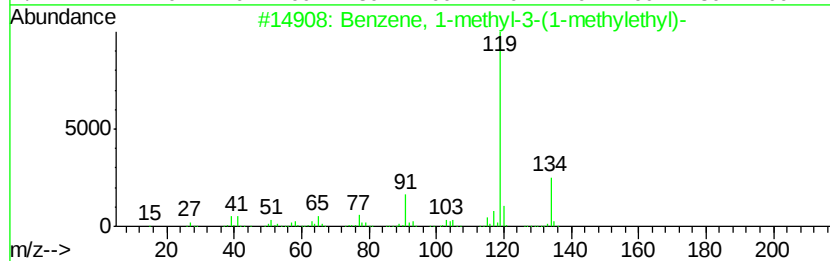
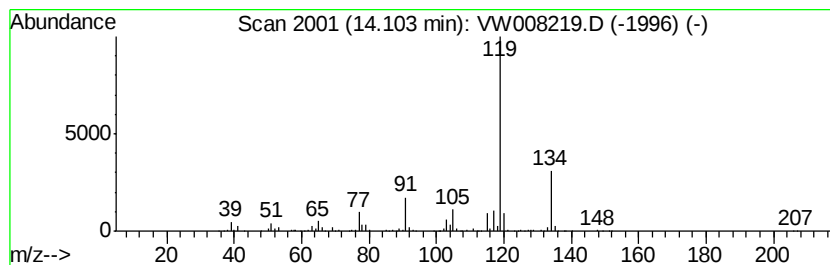
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-methyl-3-(1-meth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.10	20.29 ug/l	194869	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	94
2		o-Cymene	134	C10H14	000527-84-4	93
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91



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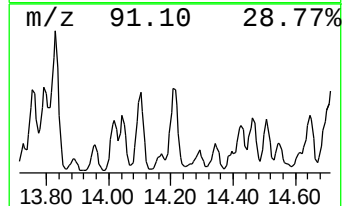
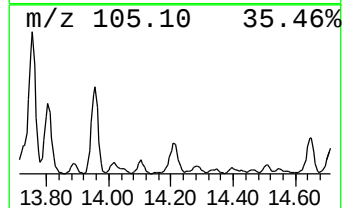
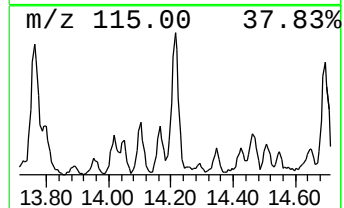
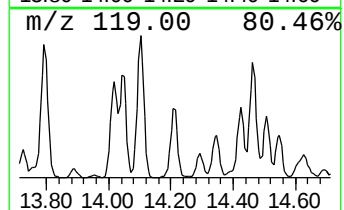
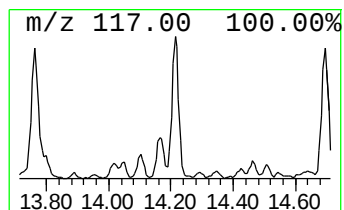
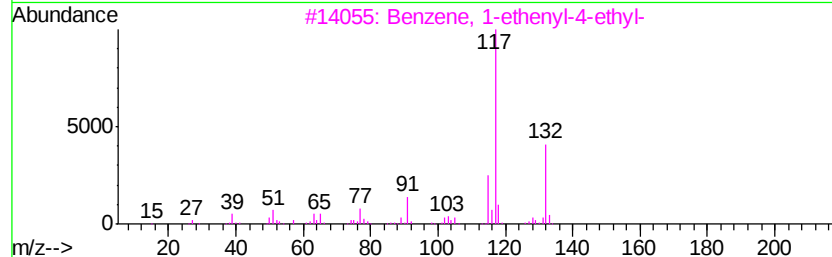
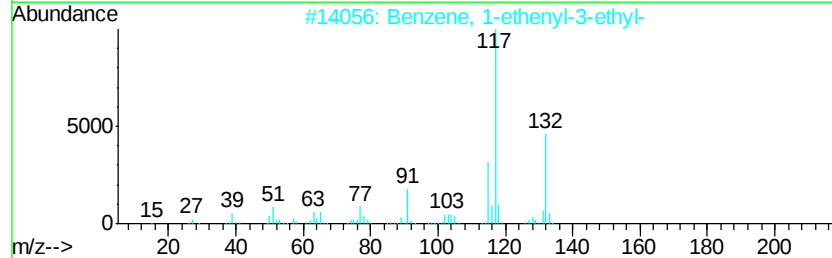
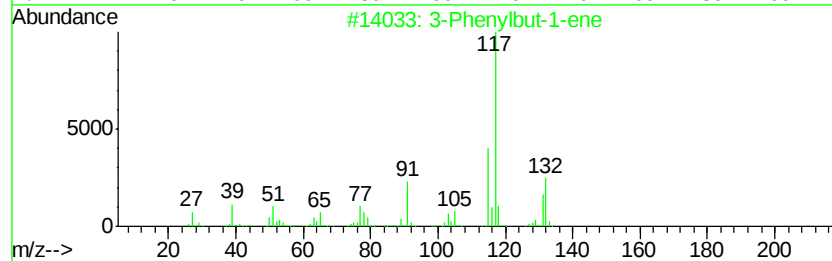
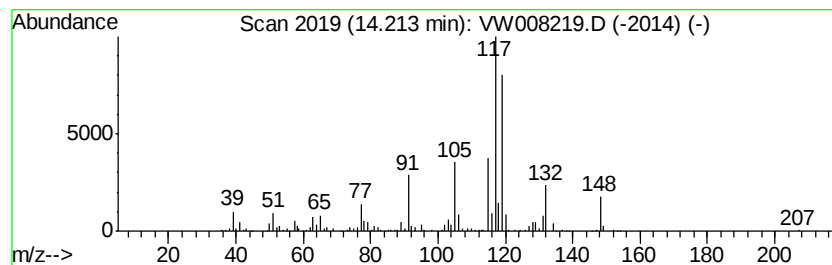
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Quant Title : SW846 8260

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

Peak Number 4 3-Phenylbut-1-ene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	25.70 ug/l	246827	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	55
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	49
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	46
4		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	45
5		Indan, 1-methyl-	132	C10H12	000767-58-8	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW011119\
Data File : VW008219.D
Acq On : 12 Jan 2019 04:28
Operator : SY/VA
Sample : K1119-03
Misc : 5.14G/5ML/MSVOA W/SOIL
ALS Vial : 34 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DS-B

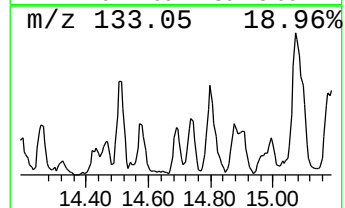
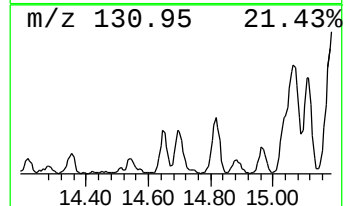
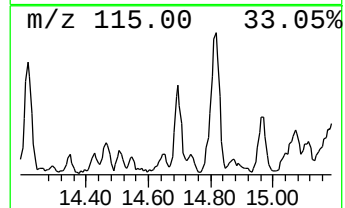
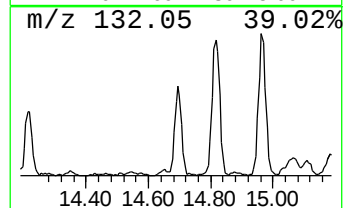
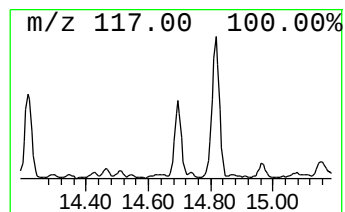
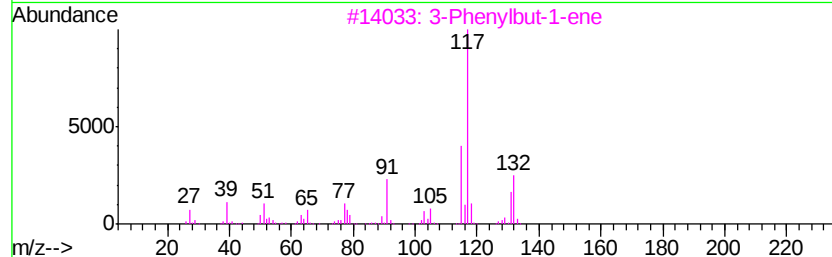
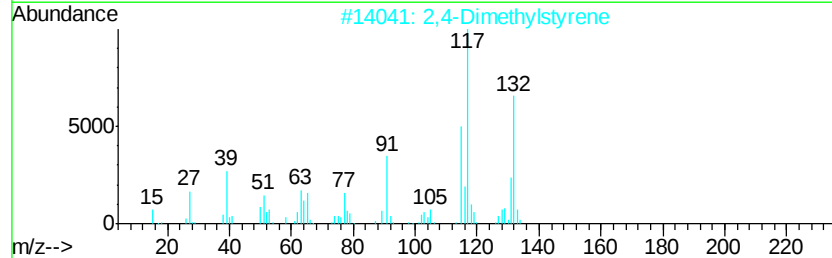
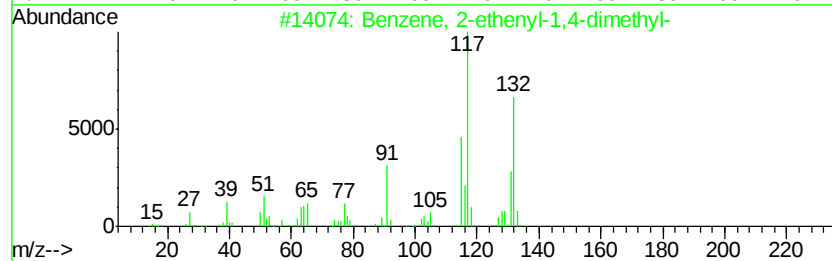
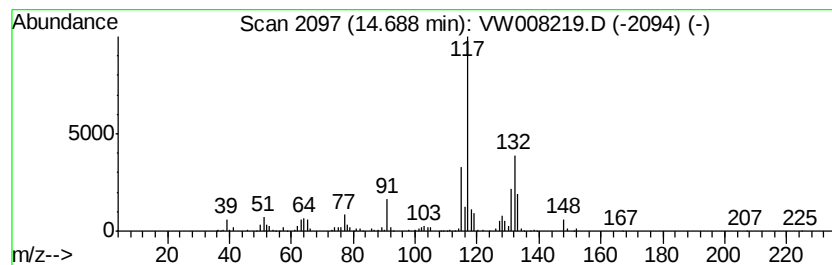
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	18.74 ug/l	179982	1,4-Dichlorobenzene-d4	13.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2		2,4-Dimethylstyrene	132	C10H12	002234-20-0	91
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
5		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW011119\
Data File : VW008219.D
Acq On : 12 Jan 2019 04:28
Operator : SY/VA
Sample : K1119-03
Misc : 5.14G/5ML/MSVOA W/SOIL
ALS Vial : 34 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DS-B

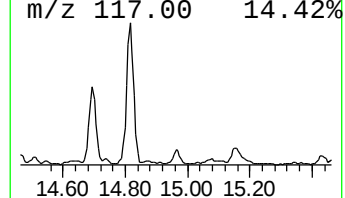
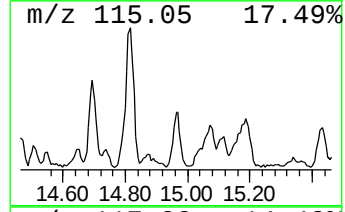
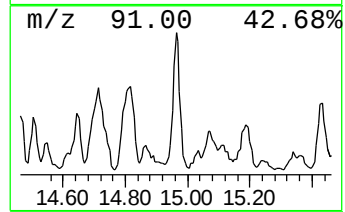
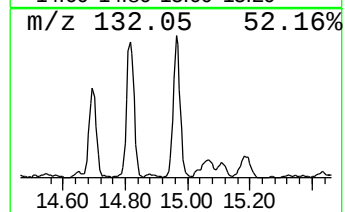
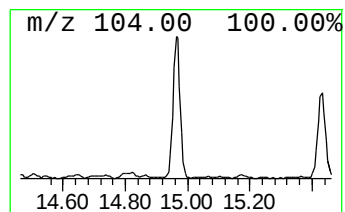
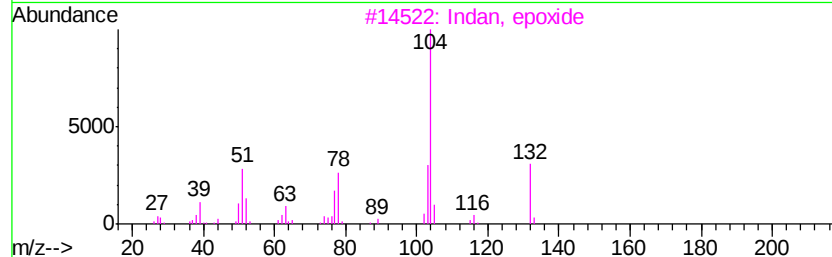
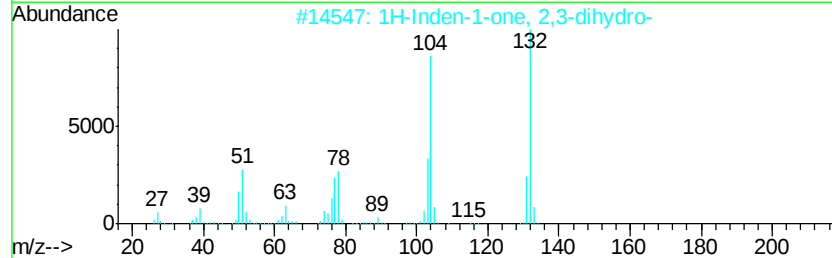
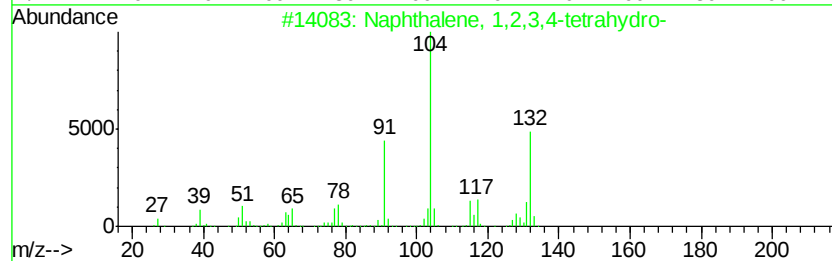
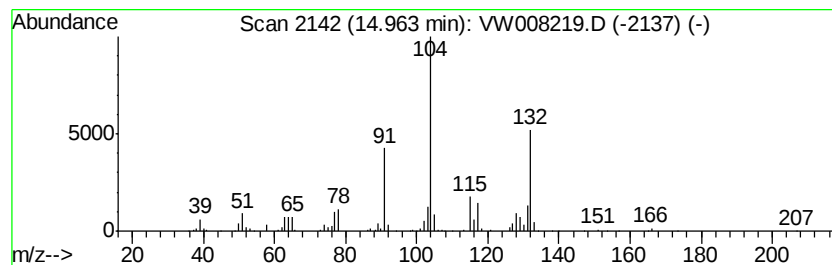
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	20.62 ug/l	198092	1,4-Dichlorobenzene-d4	13.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2		1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	59
3		Indan, epoxide	132	C9H8O	000768-22-9	58
4		2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	53
5		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW011119\
Data File : VW008219.D
Acq On : 12 Jan 2019 04:28
Operator : SY/VA
Sample : K1119-03
Misc : 5.14G/5ML/MSVOA W/SOIL
ALS Vial : 34 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DS-B

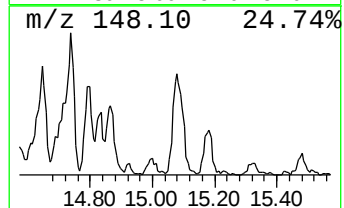
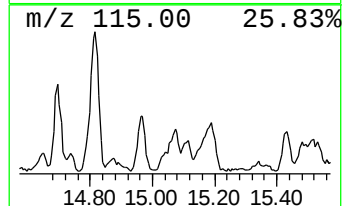
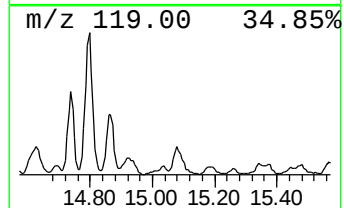
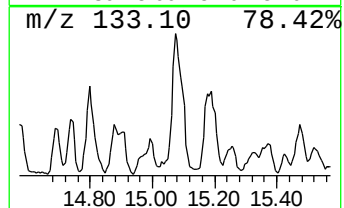
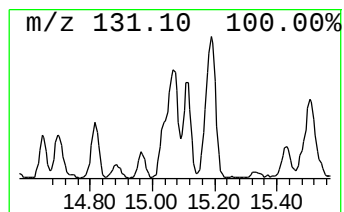
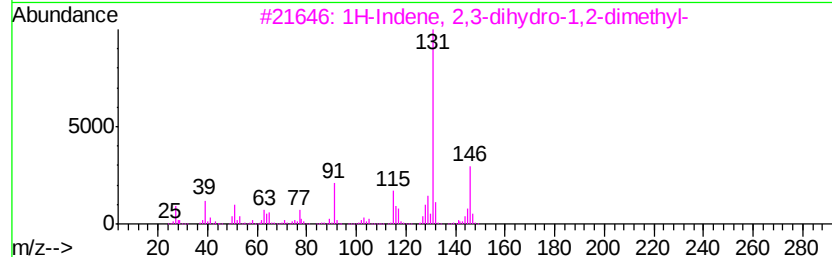
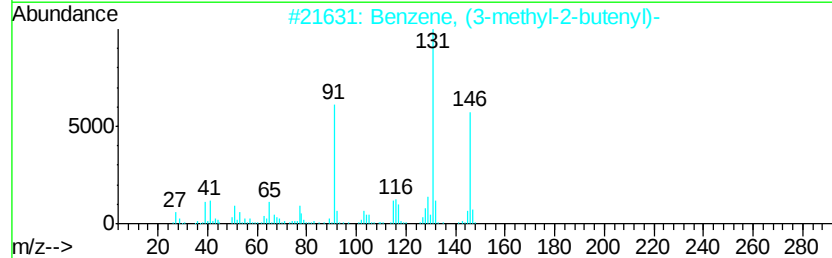
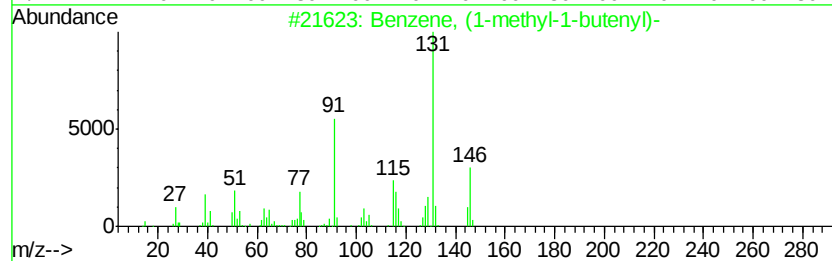
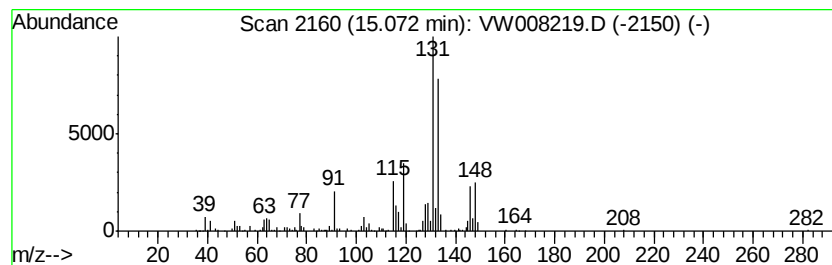
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, (1-methyl-1-butenyl)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	20.83 ug/l	200094	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	90
2		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70
3		1H-Indene, 2,3-dihydro-1,2-dimethyl-	146	C11H14	017057-82-8	70
4		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	64
5		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	55



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW011119\
Data File : VW008219.D
Acq On : 12 Jan 2019 04:28
Operator : SY/VA
Sample : K1119-03
Misc : 5.14G/5ML/MSVOA W/SOIL
ALS Vial : 34 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DS-B

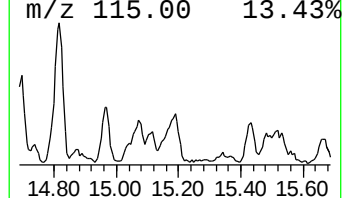
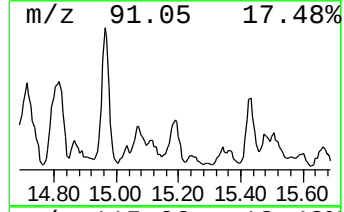
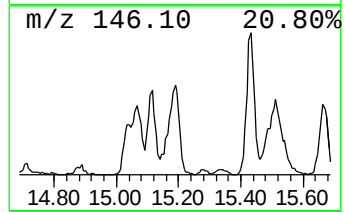
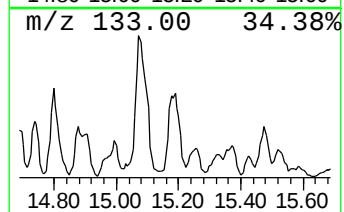
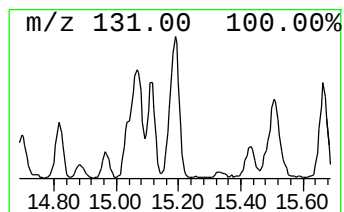
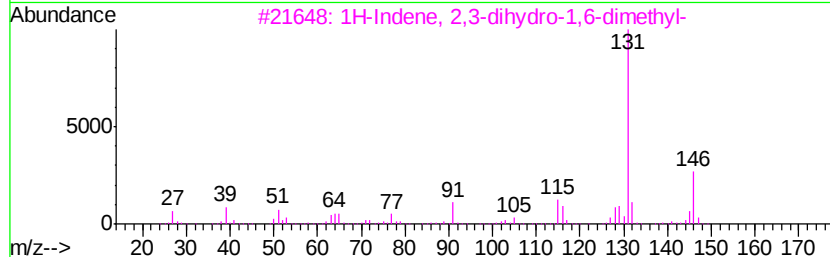
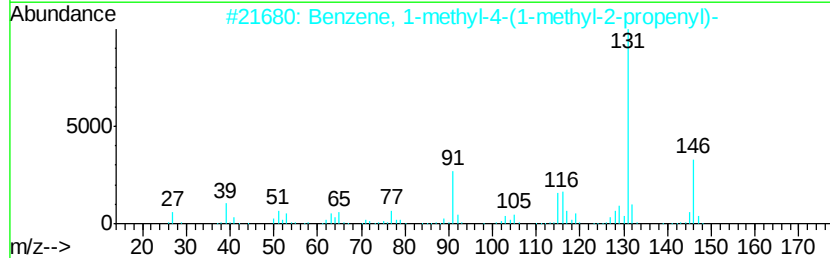
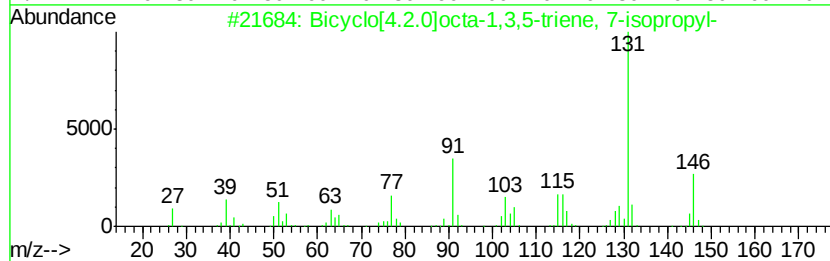
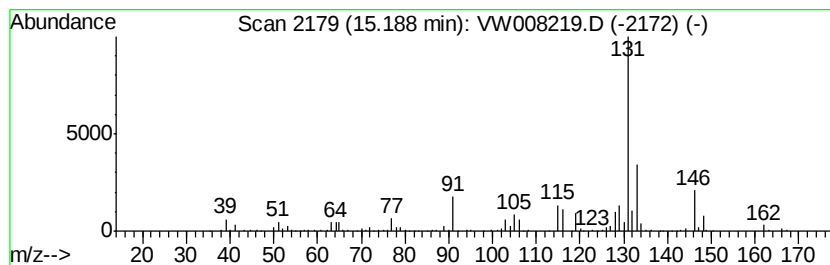
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

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

Peak Number 9 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	20.23 ug/l	194287	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	93
2		Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	81
3		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	76
4		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	76
5		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW011119\
Data File : VW008219.D
Acq On : 12 Jan 2019 04:28
Operator : SY/VA
Sample : K1119-03
Misc : 5.14G/5ML/MSVOA W/SOIL
ALS Vial : 34 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DS-B

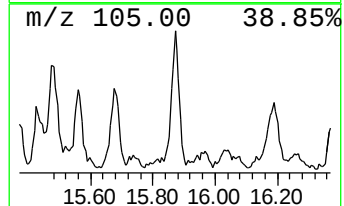
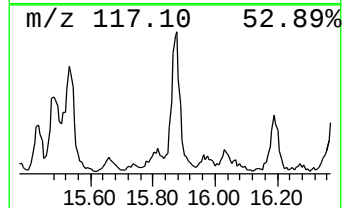
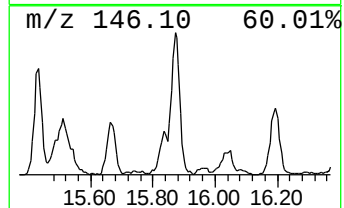
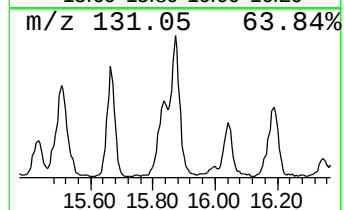
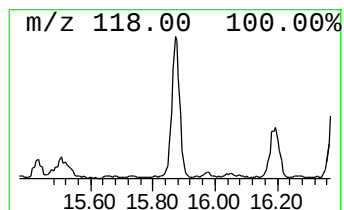
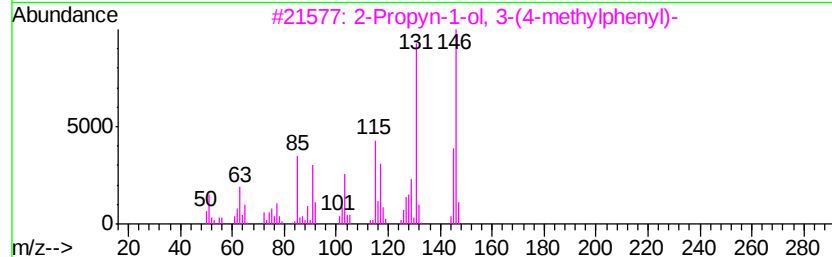
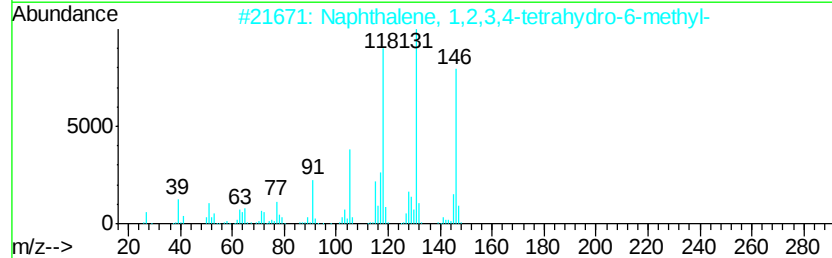
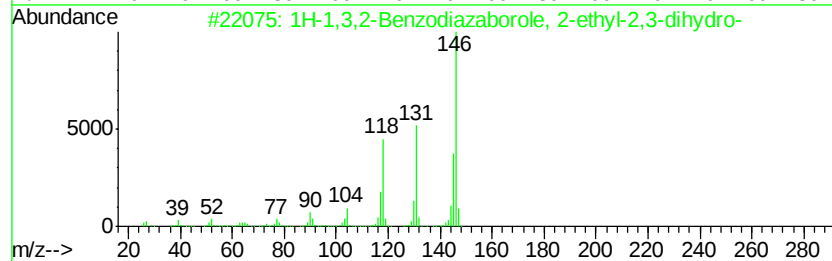
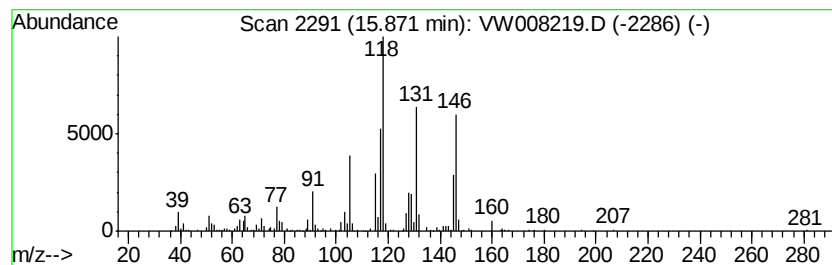
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

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

Peak Number 10 1H-1,3,2-Benzodiazaborole, ... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.87	25.15 ug/l	241616	1,4-Dichlorobenzene-d4	13.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	70	
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	58	
3	2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	49	
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	49	
5	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	43	



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW011119\
Data File : VW008219.D
Acq On : 12 Jan 2019 04:28
Operator : SY/VA
Sample : K1119-03
Misc : 5.14G/5ML/MSVOA_W/SOIL
ALS Vial : 34 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DS-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-ethyl-...	12.87	29.1	ug/l	279591	4	13.56	480280	50.0
unknown13.76	13.76	34.5	ug/l	331510	4	13.56	480280	50.0
Benzene, 1-methyl...	14.10	20.3	ug/l	194869	4	13.56	480280	50.0
3-Phenylbut-1-ene	14.21	25.7	ug/l	246827	4	13.56	480280	50.0
Benzene, 2-etheny...	14.69	18.7	ug/l	179982	4	13.56	480280	50.0
Naphthalene, 1,2,...	14.96	20.6	ug/l	198092	4	13.56	480280	50.0
Benzene, (1-methy...	15.07	20.8	ug/l	200094	4	13.56	480280	50.0
Bicyclo[4.2.0]oct...	15.19	20.2	ug/l	194287	4	13.56	480280	50.0
1H-1,3,2-Benzodia...	15.87	25.1	ug/l	241616	4	13.56	480280	50.0