

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX120420\
 Data File : VX019733.D
 Acq On : 04 Dec 2020 21:37
 Operator : JC/MD
 Sample : L4960-05
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BES-12SR

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_X\METHOD\82X120420W.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	1.294	29	34	46	rBV2	315798	1116815	5.45%	0.309%
2	1.392	46	50	56	rVB	586973	623217	3.04%	0.172%
3	1.776	107	113	129	rVB	6764775	9610346	46.93%	2.660%
4	1.983	141	147	162	rVB	2753750	3899207	19.04%	1.079%
5	2.252	186	191	201	rVB	288664	457463	2.23%	0.127%
6	2.374	204	211	223	rVB	526373	1053360	5.14%	0.292%
7	2.825	278	285	288	rVV	1754610	3779842	18.46%	1.046%
8	2.867	288	292	327	rVB	4510751	11210099	54.74%	3.102%
9	3.154	327	339	357	rVB	3559729	8145891	39.78%	2.254%
10	3.373	365	375	386	rBV	253440	584279	2.85%	0.162%
11	3.501	386	396	409	rBV	1408781	3171269	15.49%	0.878%
12	3.788	435	443	453	rVB	745959	1792936	8.76%	0.496%
13	3.898	453	461	470	rBV	457050	1172525	5.73%	0.325%
14	4.166	490	505	514	rBV2	504628	1516585	7.41%	0.420%
15	4.282	514	524	530	rVV	492026	1466872	7.16%	0.406%
16	4.373	530	539	551	rVV	3424202	9844814	48.08%	2.725%
17	4.501	551	560	575	rVB4	199953	655171	3.20%	0.181%
18	5.269	659	686	702	rVB5	426939	1944558	9.50%	0.538%
19	5.464	702	718	723	rBV3	269136	897762	4.38%	0.248%
20	5.556	723	733	748	rVV5	1331053	6705812	32.75%	1.856%
21	5.696	748	756	777	rVB	1338641	4785099	23.37%	1.324%
22	5.897	777	789	802	rBV	1265672	3801784	18.57%	1.052%
23	6.031	802	811	819	rBV	205278	512939	2.50%	0.142%
24	6.233	828	844	851	rBV2	657650	2308829	11.27%	0.639%
25	6.324	851	859	868	rVV2	1549241	4696761	22.94%	1.300%
26	6.428	868	876	887	rVB	651035	1833736	8.95%	0.508%
27	6.678	904	917	932	rVB2	418495	1070576	5.23%	0.296%
28	6.830	932	942	946	rBV	570762	1448428	7.07%	0.401%
29	6.915	946	956	967	rVB5	1049723	3803405	18.57%	1.053%
30	7.123	983	990	999	rVB	262324	559060	2.73%	0.155%
31	7.232	999	1008	1016	rBV	660641	1450685	7.08%	0.401%
32	7.440	1029	1042	1053	rBV3	2159504	5735704	28.01%	1.587%
33	7.549	1053	1060	1065	rBV2	238651	477085	2.33%	0.132%
34	7.616	1065	1071	1076	rBV	259634	497247	2.43%	0.138%

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Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
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35	7.714	1081	1087	1098	rVB	522148	1033241	5.05%	0.286%
36	7.934	1112	1123	1130	rBV4	254657	855007	4.18%	0.237%
37	8.037	1130	1140	1151	rVB	856421	1996985	9.75%	0.553%
38	8.159	1151	1160	1162	rBV	1212945	2336262	11.41%	0.647%
39	8.190	1162	1165	1170	rVV	1230396	2169840	10.60%	0.601%
40	8.238	1170	1173	1182	rVV4	521450	962130	4.70%	0.266%
41	8.324	1182	1187	1189	rVV	289610	459453	2.24%	0.127%
42	8.360	1189	1193	1199	rVB3	293195	530568	2.59%	0.147%
43	8.482	1208	1213	1217	rBV3	427615	687815	3.36%	0.190%
44	8.549	1219	1224	1233	rVB2	750244	1364248	6.66%	0.378%
45	8.653	1233	1241	1244	rBV3	433142	931580	4.55%	0.258%
46	8.696	1244	1248	1254	rVB	1206324	1900258	9.28%	0.526%
47	8.964	1285	1292	1297	rBV4	295004	562489	2.75%	0.156%
48	9.147	1313	1322	1327	rBV3	263477	575512	2.81%	0.159%
49	9.202	1327	1331	1336	rBV	556482	811239	3.96%	0.225%
50	9.543	1383	1387	1393	rVV3	273826	600007	2.93%	0.166%
51	10.092	1473	1477	1482	rVB	1063667	1400217	6.84%	0.388%
52	10.238	1495	1501	1506	rVB	12622505	16972464	82.88%	4.697%
53	10.342	1511	1518	1524	rVB	6776860	8848051	43.21%	2.449%
54	11.000	1620	1626	1634	rBV	3152726	3833990	18.72%	1.061%
55	11.122	1640	1646	1649	rBV	1003704	1283328	6.27%	0.355%
56	11.341	1677	1682	1689	rVB	8468211	10507955	51.31%	2.908%
57	11.439	1689	1698	1702	rVV	6480545	10401356	50.79%	2.879%
58	11.488	1702	1706	1718	rVB	5788754	7443656	36.35%	2.060%
59	11.652	1727	1733	1741	rVB	3180031	3922730	19.16%	1.086%
60	11.793	1751	1756	1761	rVB	16707612	19725216	96.33%	5.459%
61	11.902	1768	1774	1776	rBV	526133	674687	3.29%	0.187%
62	11.927	1776	1778	1783	rVB	723022	809824	3.95%	0.224%
63	12.012	1788	1792	1795	rVV	525786	657840	3.21%	0.182%
64	12.061	1795	1800	1804	rVV2	1281917	1900612	9.28%	0.526%
65	12.128	1804	1811	1815	rVV	10069656	12239423	59.77%	3.387%
66	12.286	1831	1837	1843	rBV	15313919	20477621	100.00%	5.667%
67	12.353	1843	1848	1855	rVB3	4427331	6734096	32.89%	1.864%
68	12.445	1858	1863	1868	rBV2	1019313	1331877	6.50%	0.369%
69	12.500	1868	1872	1879	rVB	2791079	3704504	18.09%	1.025%
70	12.573	1879	1884	1886	rBV	2825427	3862030	18.86%	1.069%
71	12.591	1886	1887	1892	rVB	1554076	1346793	6.58%	0.373%

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Integration Parameters: RTEINT.P

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72	12.652	1892	1897	1900	rBV	11399603	13331436	65.10%	3.690%
73	12.683	1900	1902	1907	rVB	1545227	1758975	8.59%	0.487%
74	12.738	1907	1911	1919	rBV2	5085046	7252354	35.42%	2.007%
75	12.884	1930	1935	1940	rVB2	1675388	2168568	10.59%	0.600%
76	12.963	1940	1948	1951	rBV	6495610	8229889	40.19%	2.278%
77	13.000	1951	1954	1960	rVV	5707450	6809091	33.25%	1.884%
78	13.061	1960	1964	1970	rVV3	628757	1289126	6.30%	0.357%
79	13.176	1980	1983	1984	rVV	451991	501291	2.45%	0.139%
80	13.207	1984	1988	1993	rVV	5604416	6763228	33.03%	1.872%
81	13.298	1998	2003	2004	rVV	647348	891788	4.35%	0.247%
82	13.329	2004	2008	2014	rVV2	10713670	15523747	75.81%	4.296%
83	13.384	2014	2017	2019	rVV3	842526	1142181	5.58%	0.316%
84	13.408	2019	2021	2026	rVV	694894	776179	3.79%	0.215%
85	13.463	2026	2030	2037	rVB	1886581	2316044	11.31%	0.641%
86	13.542	2038	2043	2046	rBV2	652152	908816	4.44%	0.252%
87	13.585	2046	2050	2052	rVV	1370305	1861997	9.09%	0.515%
88	13.621	2052	2056	2062	rVV3	1787618	3390136	16.56%	0.938%
89	13.682	2062	2066	2067	rVV	655200	934672	4.56%	0.259%
90	13.701	2067	2069	2076	rVB2	1394754	2155857	10.53%	0.597%
91	13.817	2080	2088	2097	rBV	6298487	7984078	38.99%	2.210%
92	13.963	2106	2112	2116	rBV2	720834	1064025	5.20%	0.294%
93	14.012	2116	2120	2124	rVB	592623	687169	3.36%	0.190%
94	14.115	2132	2137	2143	rBV	1267755	1536890	7.51%	0.425%
95	14.256	2155	2160	2165	rBV	1033451	1603576	7.83%	0.444%
96	14.359	2173	2177	2182	rBV	695465	898493	4.39%	0.249%
97	14.414	2182	2186	2190	rVB	700564	894928	4.37%	0.248%
98	14.676	2225	2229	2239	rVB	4821989	5802426	28.34%	1.606%
99	14.816	2247	2252	2262	rBV	2882372	3849725	18.80%	1.065%
100	15.664	2382	2391	2394	rBV	330943	514250	2.51%	0.142%

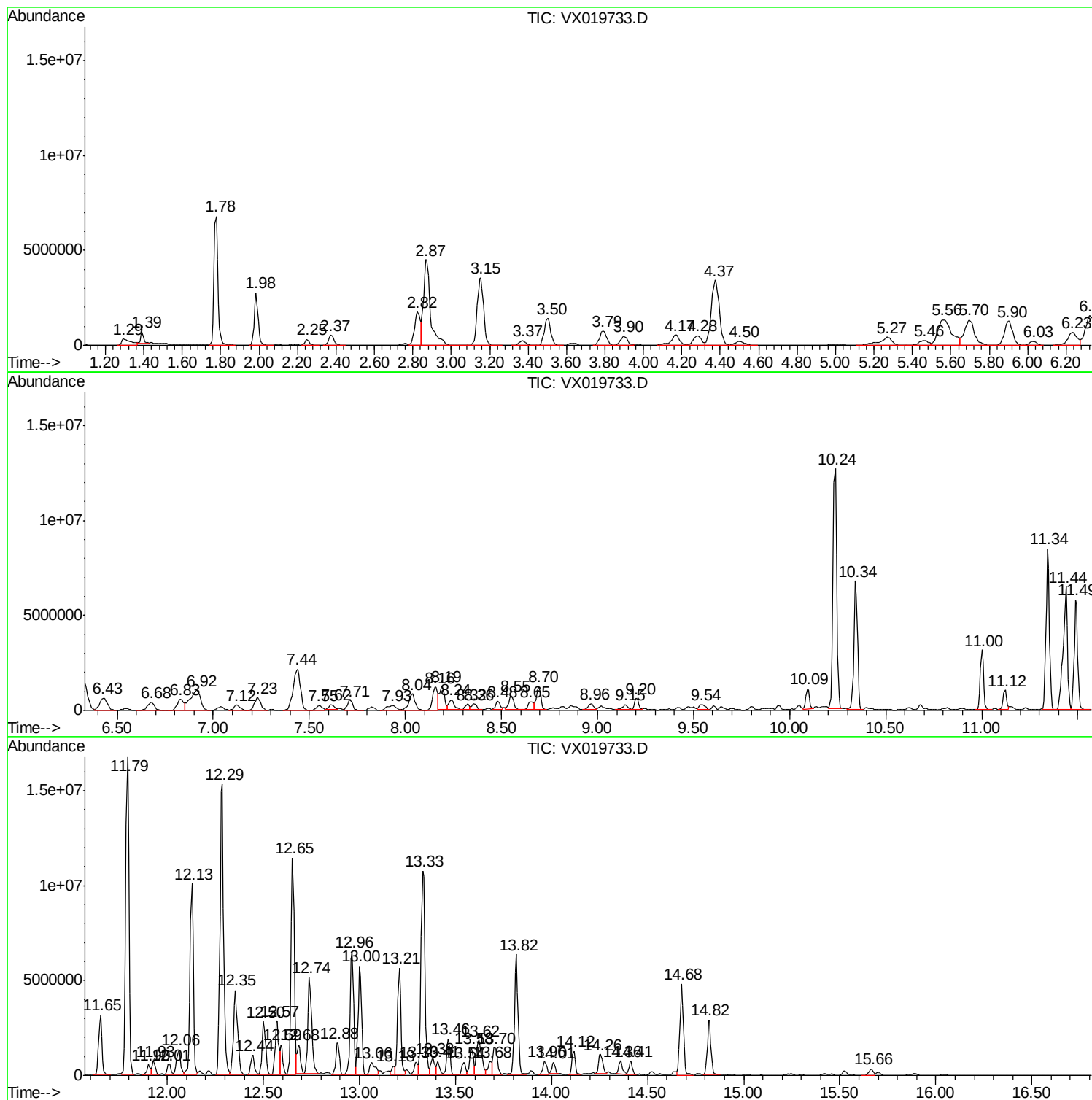
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X120420W.M
Quant Title : SW846 8260

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



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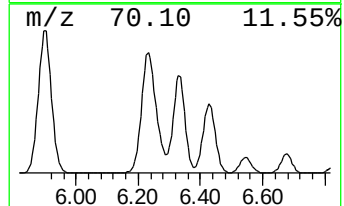
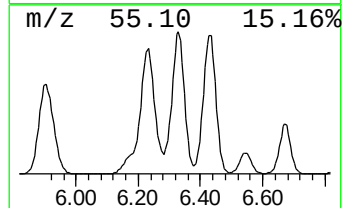
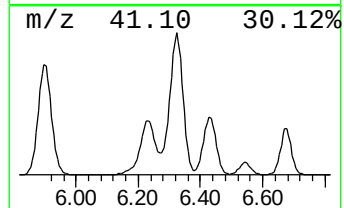
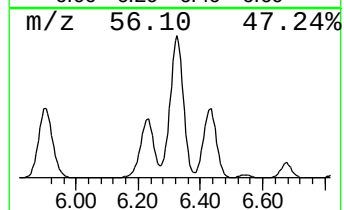
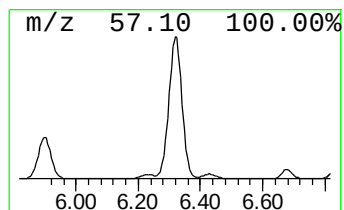
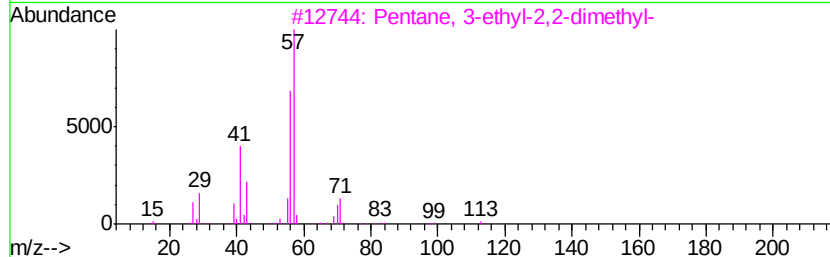
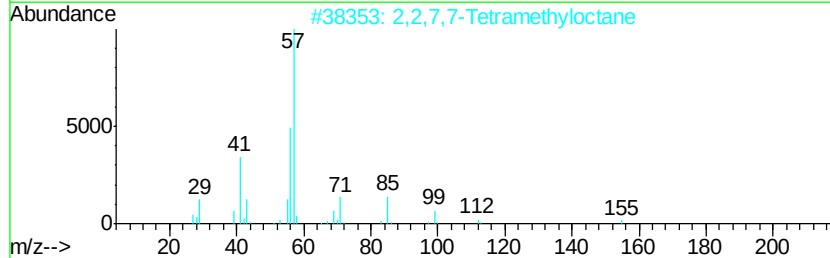
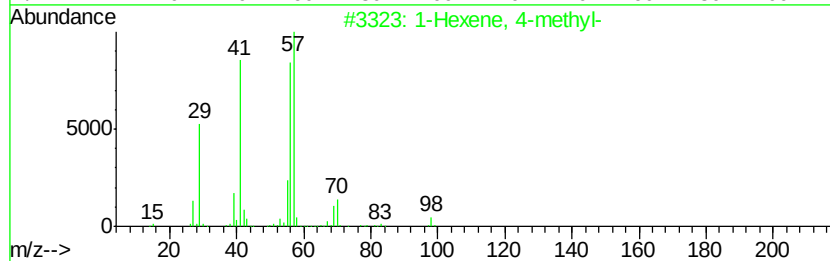
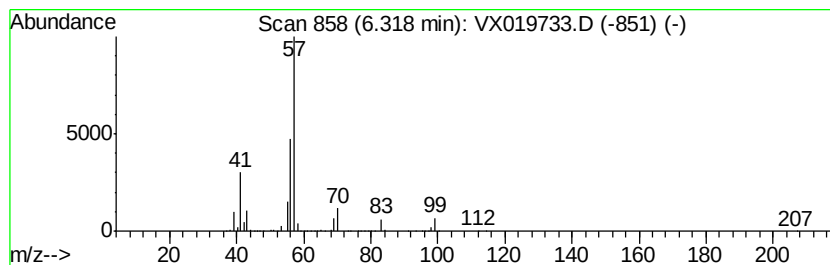
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X120420W.M
Quant Title : SW846 8260

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

Peak Number 1 1-Hexene, 4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.32	162.13 ug/l	4696760	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexene, 4-methyl-	98	C7H14	003769-23-1	59
2		2,2,7,7-Tetramethyloctane	170	C12H26	001071-31-4	50
3		Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	45
4		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	42
5		Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	40



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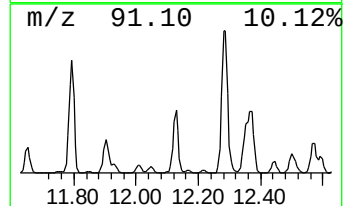
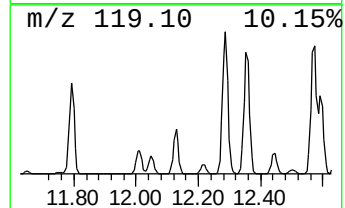
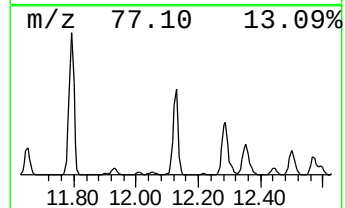
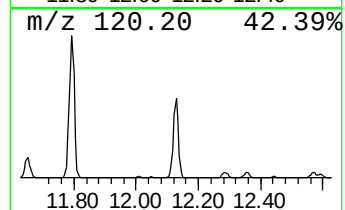
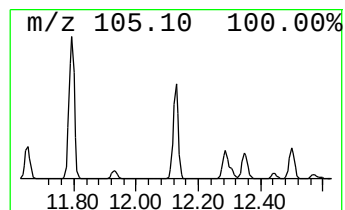
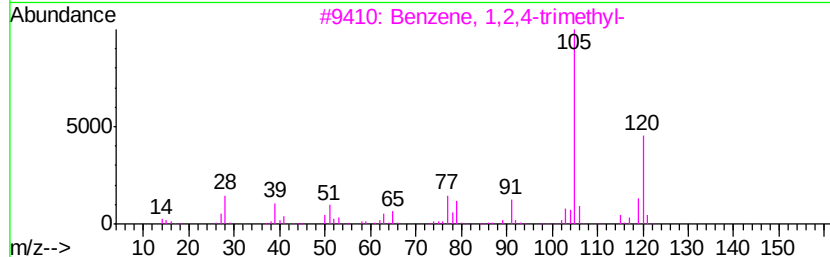
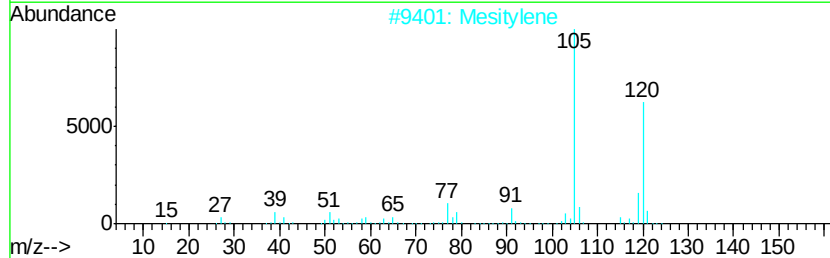
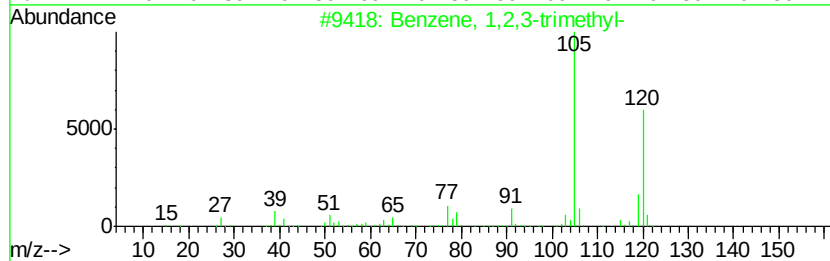
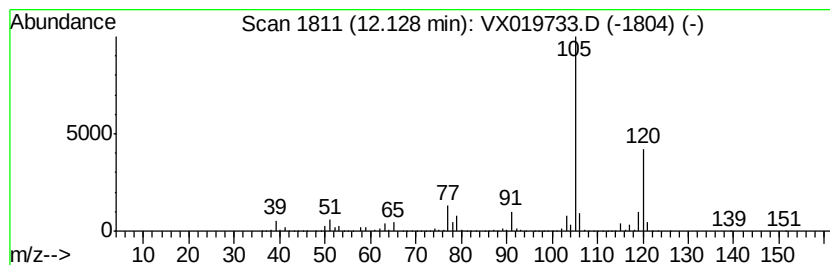
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Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1,2,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	321.99 ug/l	12239400	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Mesitylene	120	C9H12	000108-67-8	94
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



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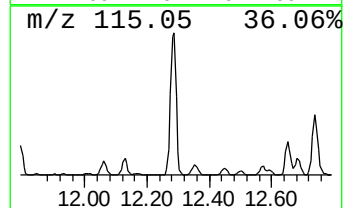
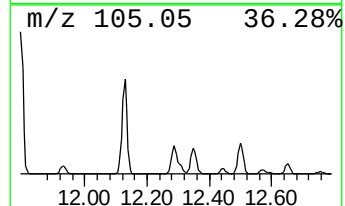
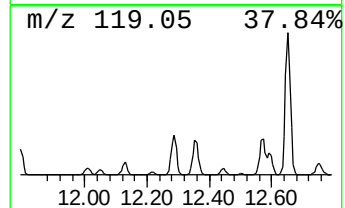
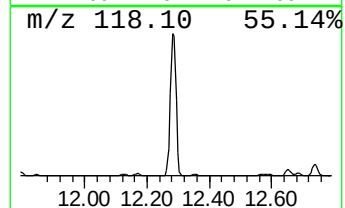
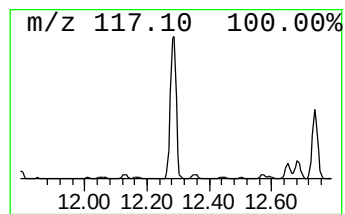
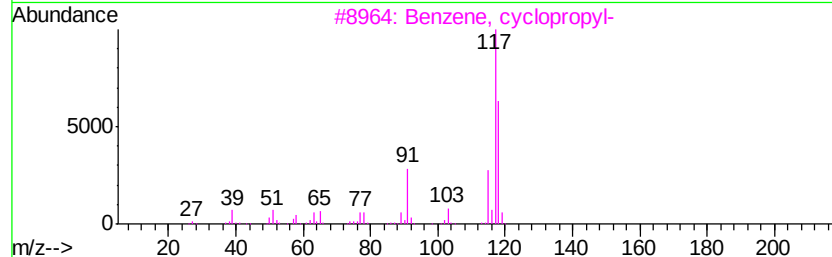
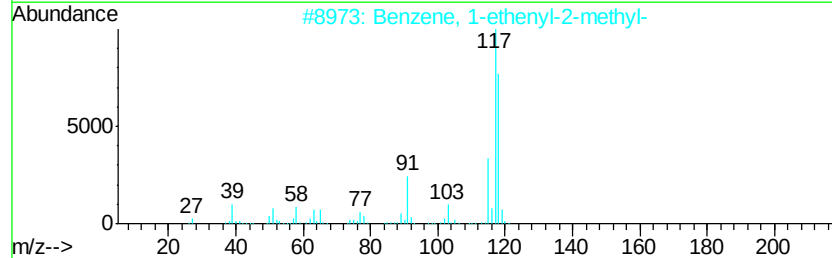
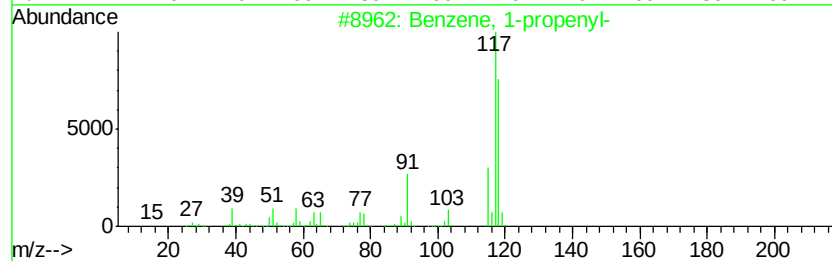
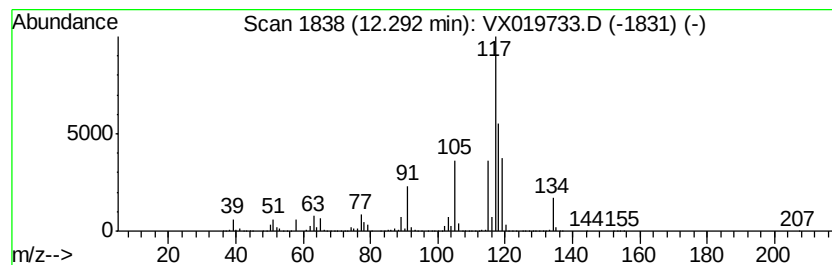
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ClientSampled :
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-propenyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	538.71 ug/l	20477600	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-propenyl-	118 C9H10	000637-50-3 92
2		Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 91
3		Benzene, cyclopropyl-	118 C9H10	000873-49-4 64
4		Indane	118 C9H10	000496-11-7 64
5		(Z)-1-Phenylpropene	118 C9H10	000766-90-5 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120420\
Data File : VX019733.D
Acq On : 04 Dec 2020 21:37
Operator : JC/MD
Sample : L4960-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 26 Sample Multiplier: 1

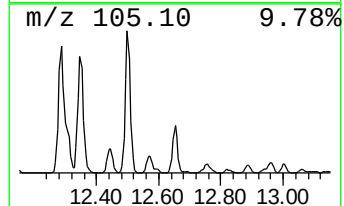
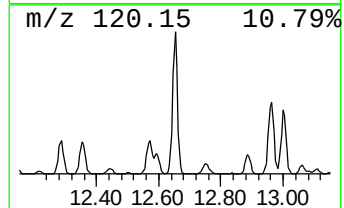
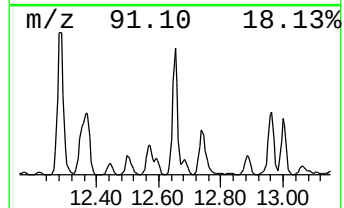
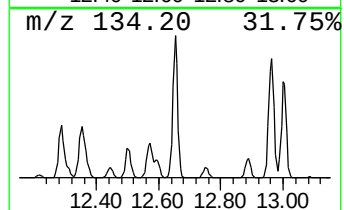
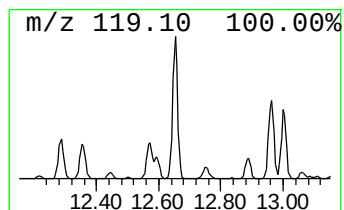
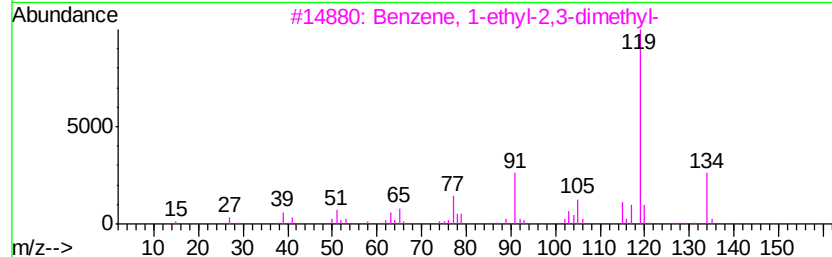
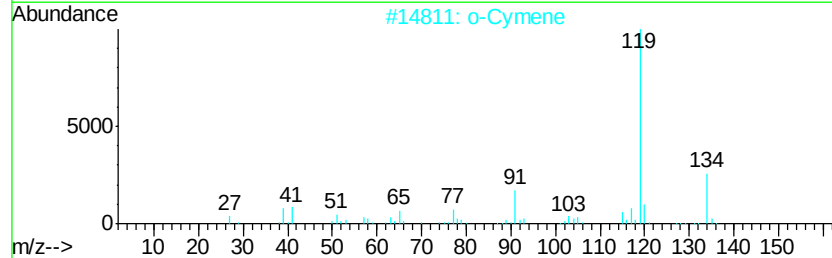
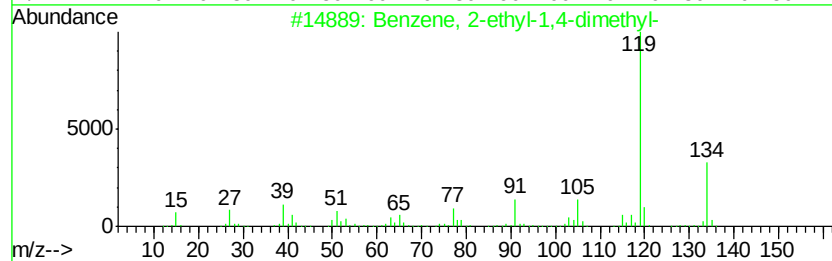
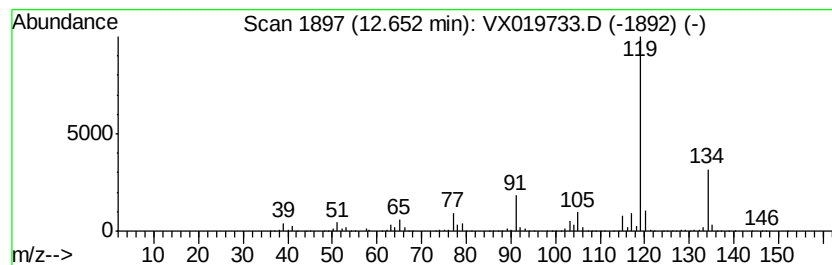
Instrument :
MSVOA_X
ClientSampleId :
BES-12SR

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X120420W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.65	350.71 ug/l	13331400	1,4-Dichlorobenzene-d4	12.06	
Hit# of	5	Tentative ID	MW MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120420\
Data File : VX019733.D
Acq On : 04 Dec 2020 21:37
Operator : JC/MD
Sample : L4960-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BES-12SR

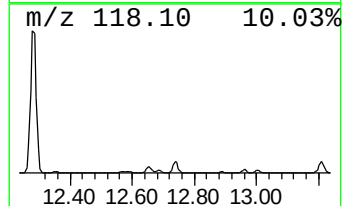
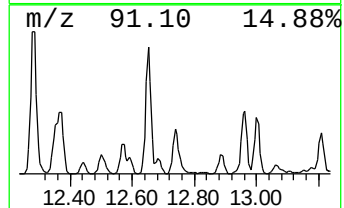
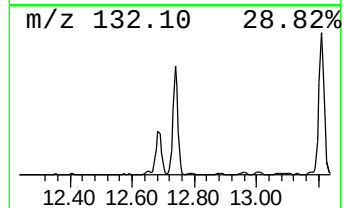
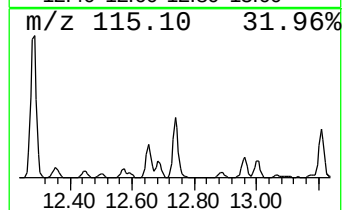
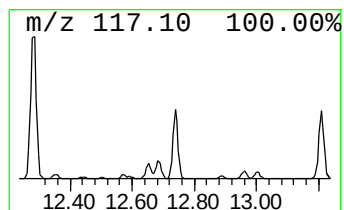
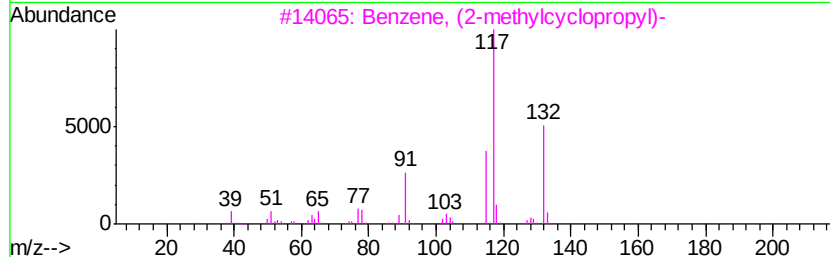
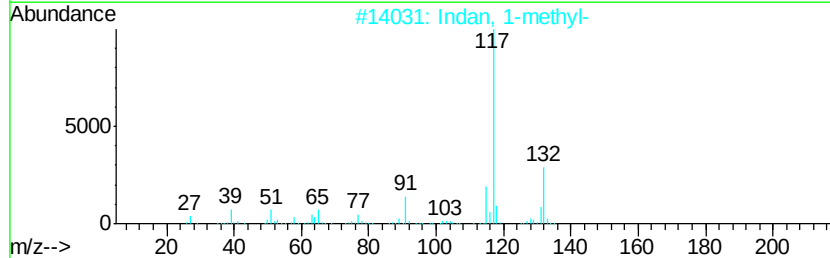
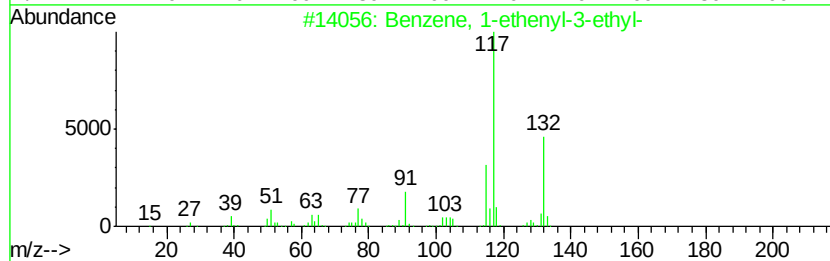
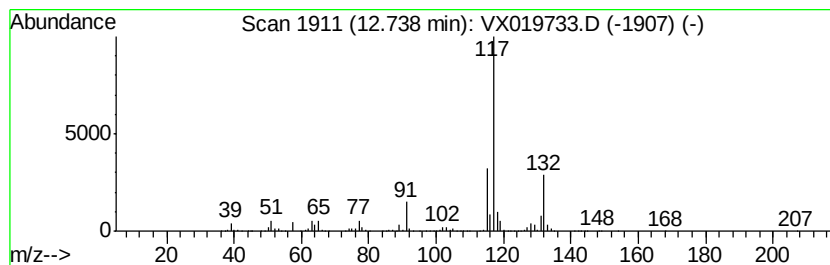
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	190.79 ug/l	7252350	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
2		Indan, 1-methyl-	132	C10H12	000767-58-8	87
3		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
4		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	83
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120420\
Data File : VX019733.D
Acq On : 04 Dec 2020 21:37
Operator : JC/MD
Sample : L4960-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BES-12SR

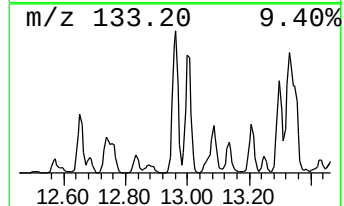
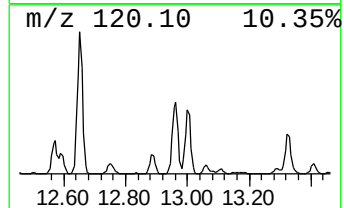
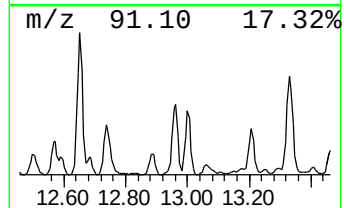
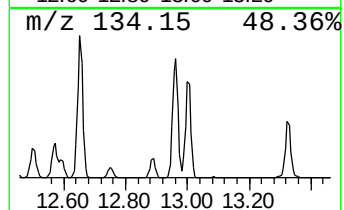
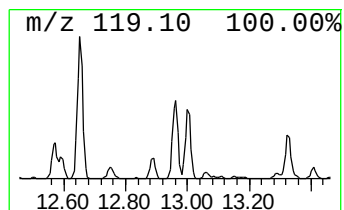
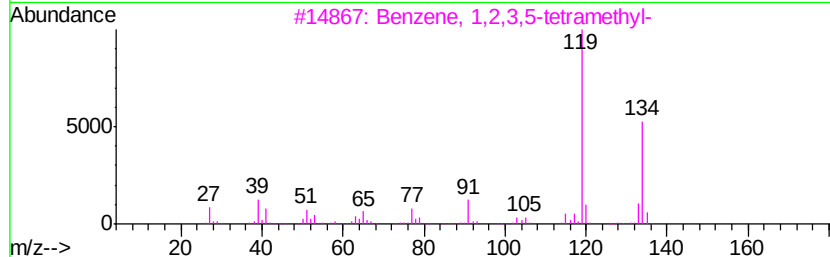
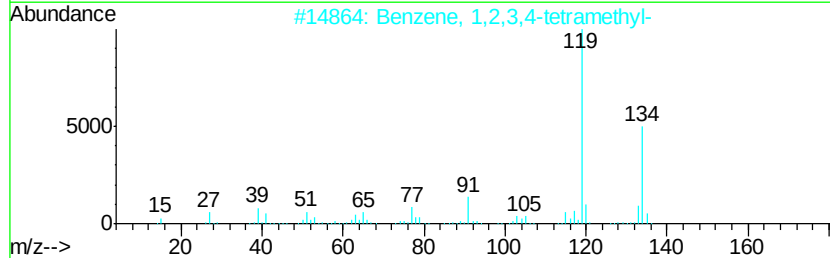
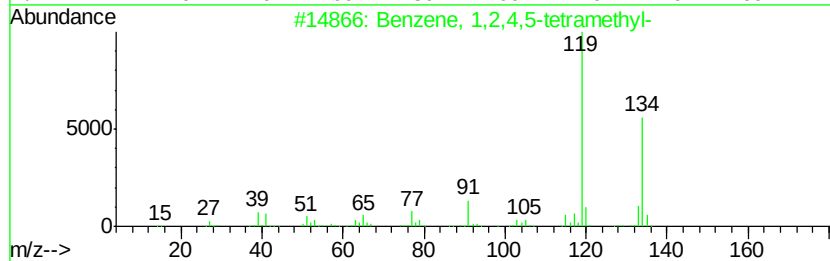
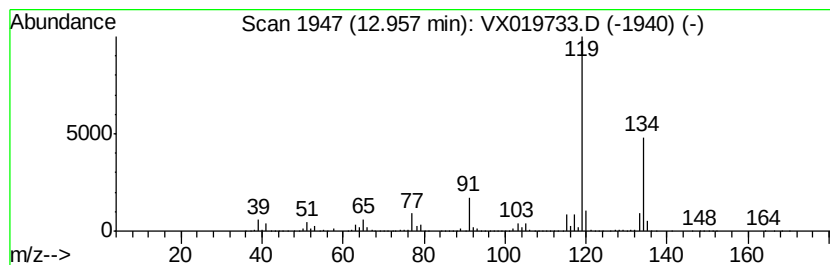
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	216.51 ug/l	8229890	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120420\
Data File : VX019733.D
Acq On : 04 Dec 2020 21:37
Operator : JC/MD
Sample : L4960-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
BES-12SR

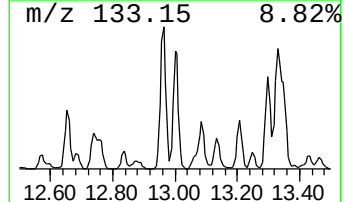
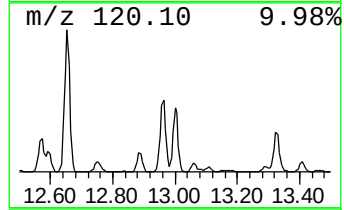
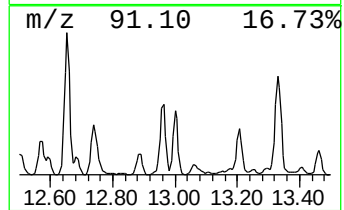
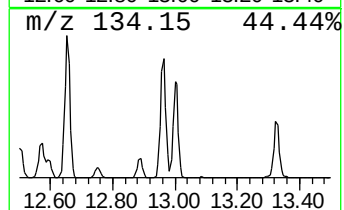
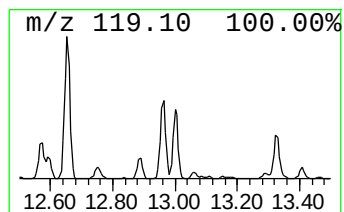
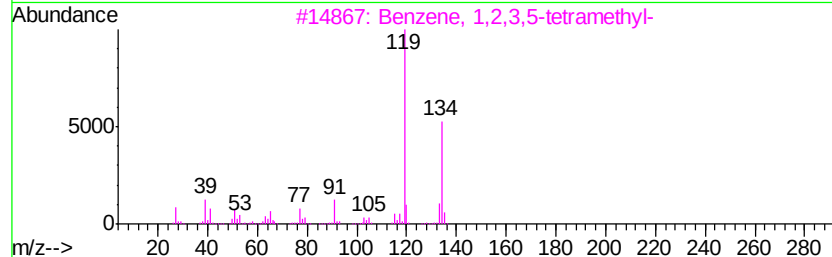
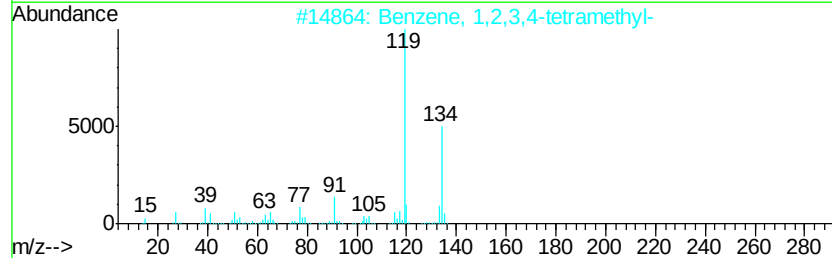
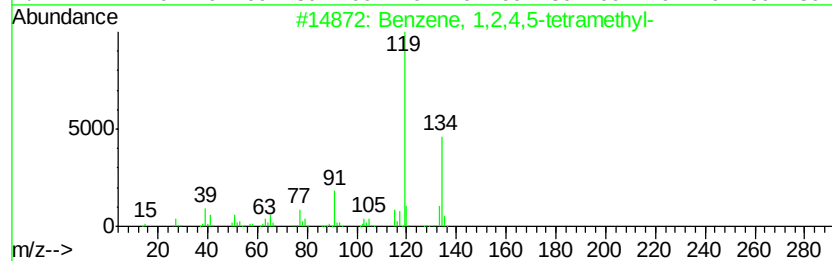
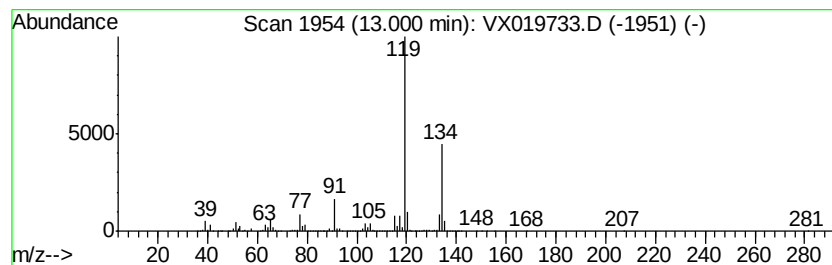
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.00	179.13 ug/l	6809090	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120420\
Data File : VX019733.D
Acq On : 04 Dec 2020 21:37
Operator : JC/MD
Sample : L4960-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BES-12SR

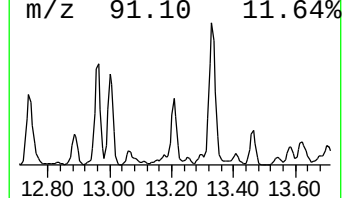
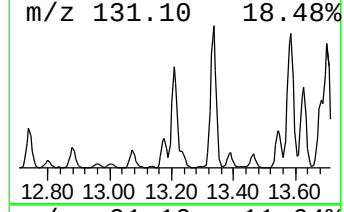
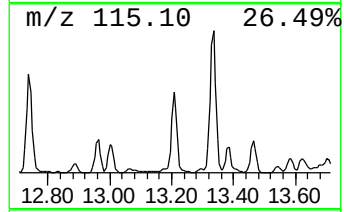
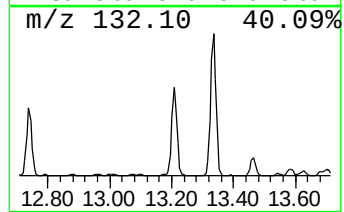
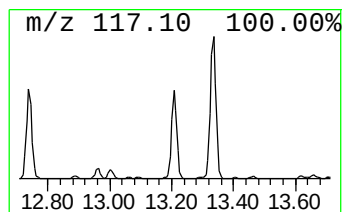
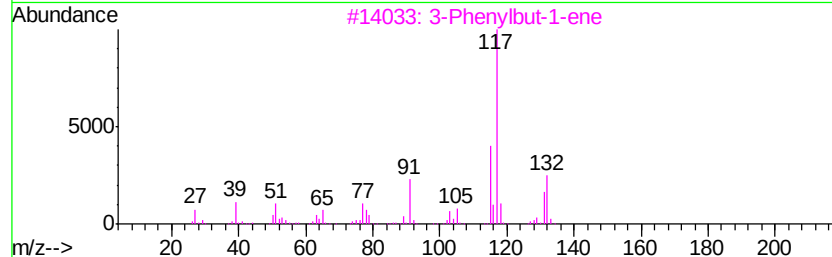
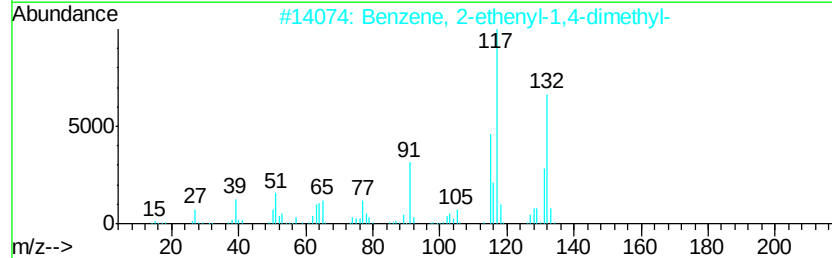
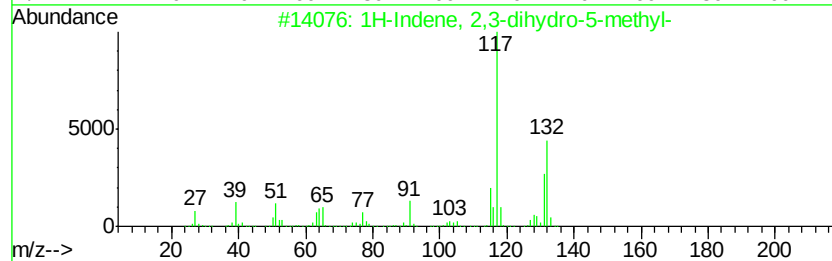
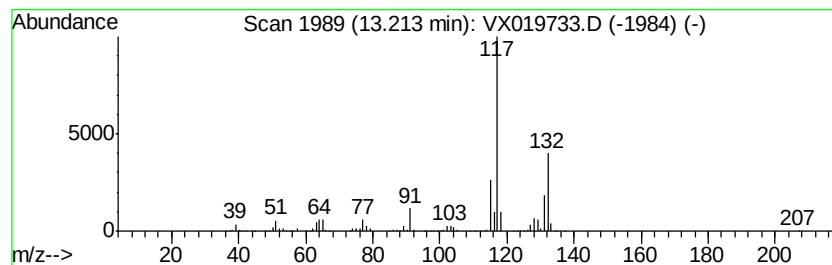
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Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.21	177.92 ug/l	6763230	1,4-Dichlorobenzene-d4	12.06

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, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4		Indan, 1-methyl-	132	C10H12	000767-58-8	90
5		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120420\
Data File : VX019733.D
Acq On : 04 Dec 2020 21:37
Operator : JC/MD
Sample : L4960-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BES-12SR

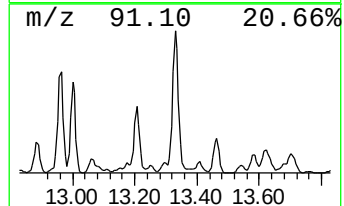
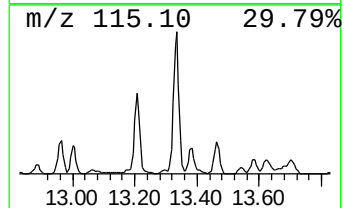
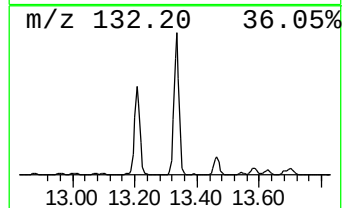
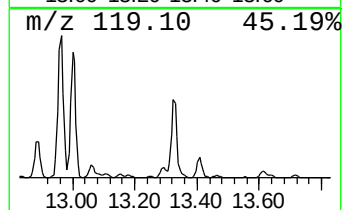
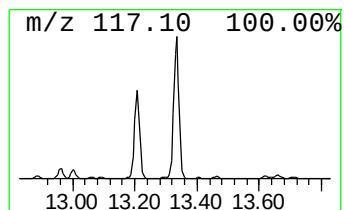
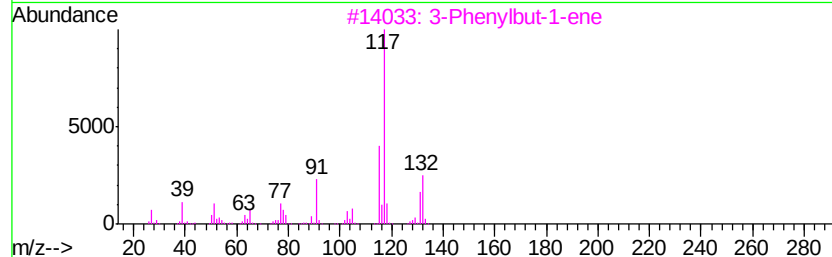
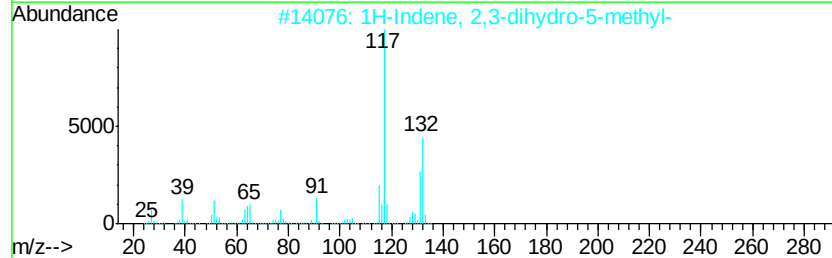
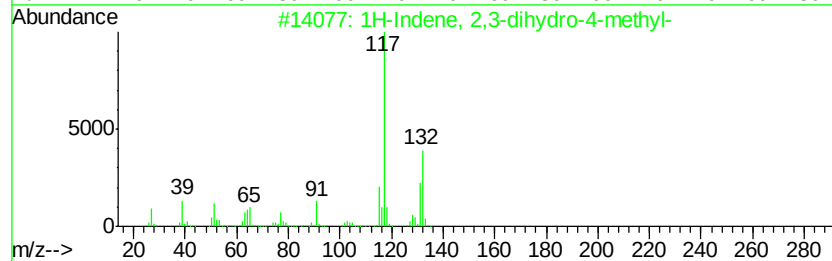
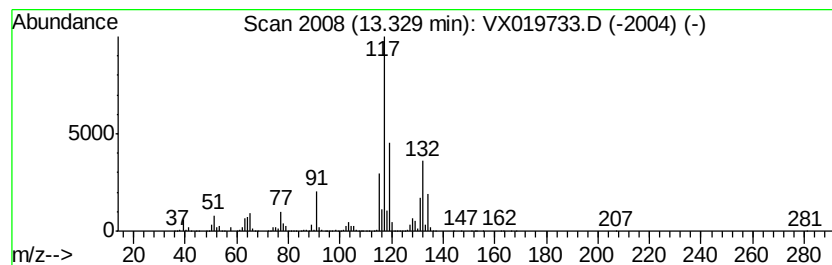
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Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.33	408.39 ug/l	15523700	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	91
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	86
4		Indan, 1-methyl-	132	C10H12	000767-58-8	70
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120420\
Data File : VX019733.D
Acq On : 04 Dec 2020 21:37
Operator : JC/MD
Sample : L4960-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
BES-12SR

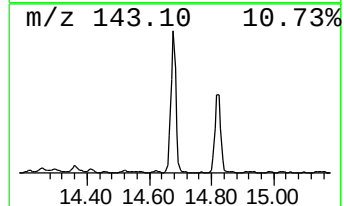
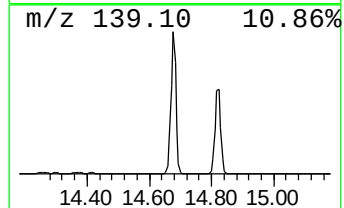
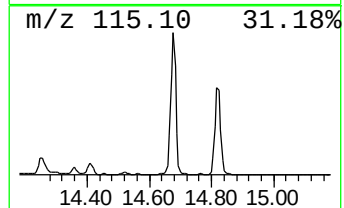
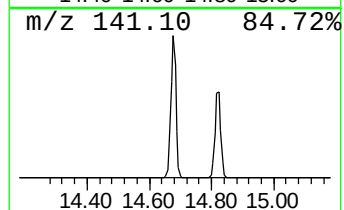
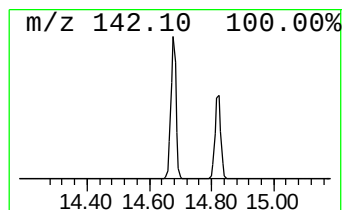
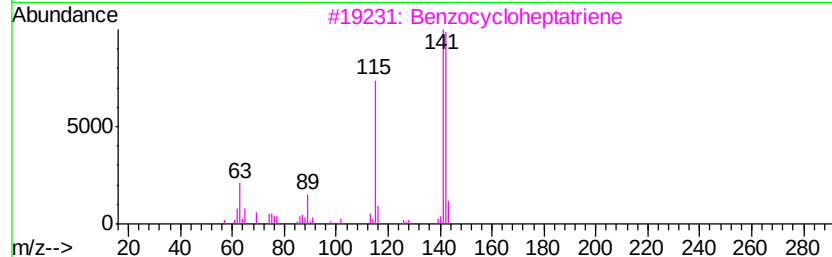
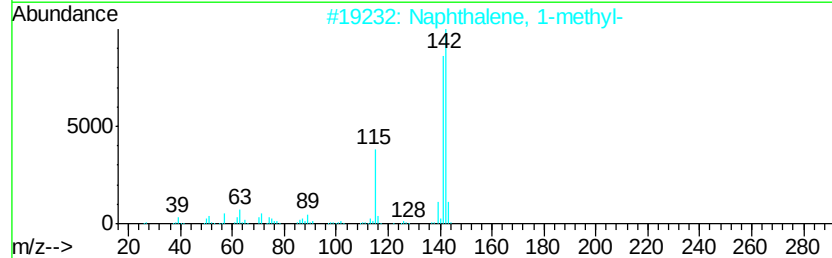
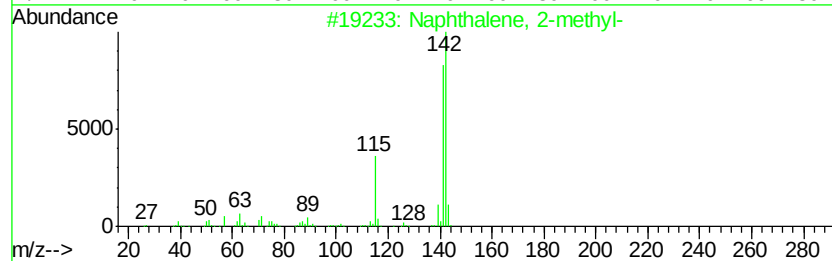
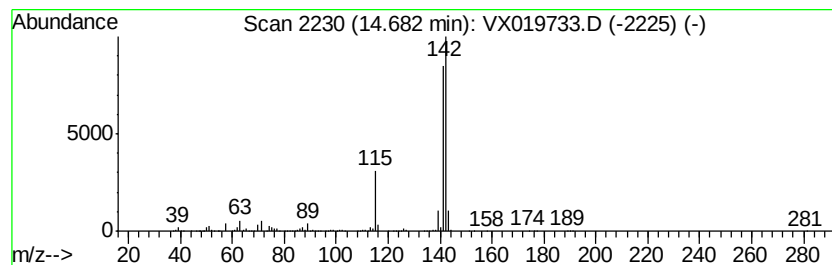
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X120420W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.68	152.65 ug/l	5802430	1,4-Dichlorobenzene-d4	12.06

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		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX120420\
Data File : VX019733.D
Acq On : 04 Dec 2020 21:37
Operator : JC/MD
Sample : L4960-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BES-12SR

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X120420W.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
1-Hexene, 4-methyl-	6.32	162.1	ug/l	4696760	2	6.83	1448430	50.0
Benzene, 1,2,3-tr...	12.13	322.0	ug/l	12239400	4	12.06	1900610	50.0
Benzene, 1-propenyl-	12.29	538.7	ug/l	20477600	4	12.06	1900610	50.0
Benzene, 2-ethyl-...	12.65	350.7	ug/l	13331400	4	12.06	1900610	50.0
Benzene, 1-etheny...	12.74	190.8	ug/l	7252350	4	12.06	1900610	50.0
Benzene, 1,2,4,5-...	12.96	216.5	ug/l	8229890	4	12.06	1900610	50.0
Benzene, 1,2,3,4-...	13.00	179.1	ug/l	6809090	4	12.06	1900610	50.0
1H-Indene, 2,3-di...	13.21	177.9	ug/l	6763230	4	12.06	1900610	50.0
1H-Indene, 2,3-di...	13.33	408.4	ug/l	15523700	4	12.06	1900610	50.0
Naphthalene, 1-me...	14.68	152.7	ug/l	5802430	4	12.06	1900610	50.0