

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX021919\
 Data File : VX007612.D
 Acq On : 19 Feb 2019 14:44
 Operator : JC/SP
 Sample : K1528-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-3

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\82X020119W.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	1.605	65	74	78	rBV4	21959	45370	1.85%	0.159%
2	2.843	271	277	281	rBV2	40982	89412	3.65%	0.314%
3	2.891	281	285	292	rVB2	76994	155666	6.36%	0.547%
4	3.172	323	331	342	rVB	116005	263072	10.74%	0.924%
5	3.519	380	388	396	rVB2	21157	53352	2.18%	0.187%
6	4.306	505	517	526	rBV2	40815	129072	5.27%	0.453%
7	5.238	660	670	683	rVB6	18520	72445	2.96%	0.254%
8	5.507	701	714	723	rBV	292853	829607	33.88%	2.914%
9	5.665	723	740	748	rBV	464929	1414660	57.77%	4.970%
10	5.927	772	783	795	rVB4	67208	205583	8.40%	0.722%
11	6.068	795	806	817	rBV	290917	739719	30.21%	2.599%
12	6.257	829	837	846	rBV4	49452	162083	6.62%	0.569%
13	6.354	846	853	862	rVB4	67022	189568	7.74%	0.666%
14	6.452	862	869	877	rVB5	25195	67301	2.75%	0.236%
15	6.860	927	936	945	rBV	713739	1677678	68.51%	5.893%
16	7.451	1024	1033	1044	rVB3	78070	208931	8.53%	0.734%
17	7.573	1047	1053	1058	rBV4	30628	61244	2.50%	0.215%
18	7.634	1058	1063	1068	rVV	39331	70562	2.88%	0.248%
19	7.848	1091	1098	1106	rVB2	79713	166145	6.78%	0.584%
20	8.031	1122	1128	1130	rBV	45434	84130	3.44%	0.296%
21	8.177	1145	1152	1160	rBV3	60067	153119	6.25%	0.538%
22	8.378	1176	1185	1191	rVB6	45059	97296	3.97%	0.342%
23	8.579	1212	1218	1225	rVB5	30803	69067	2.82%	0.243%
24	8.671	1225	1233	1234	rBV3	66281	102169	4.17%	0.359%
25	8.713	1234	1240	1250	rVB	1530109	2448745	100.00%	8.602%
26	8.988	1277	1285	1289	rBV2	45866	84824	3.46%	0.298%
27	9.037	1289	1293	1299	rVB2	61053	94850	3.87%	0.333%
28	9.165	1305	1314	1319	rBV2	66351	130316	5.32%	0.458%
29	9.219	1319	1323	1328	rBV	97479	144544	5.90%	0.508%
30	9.567	1374	1380	1385	rVV4	42439	83717	3.42%	0.294%
31	9.671	1392	1397	1401	rVV4	28635	61488	2.51%	0.216%
32	9.823	1415	1422	1427	rBV2	45012	83978	3.43%	0.295%
33	10.110	1464	1469	1478	rBV	1542386	2085950	85.18%	7.328%
34	10.360	1507	1510	1515	rVB5	26305	45674	1.87%	0.160%

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35	10.640	1548	1556	1564	rBV2	40714	79935	3.26%	0.281%
36	11.018	1613	1618	1626	rBV3	24214	47203	1.93%	0.166%
37	11.134	1632	1637	1643	rBV	1275075	1676466	68.46%	5.889%
38	11.359	1669	1674	1679	rBV	64335	77953	3.18%	0.274%
39	11.432	1682	1686	1688	rBV	47721	67451	2.75%	0.237%
40	11.670	1720	1725	1736	rVB2	33812	59428	2.43%	0.209%
41	11.804	1742	1747	1752	rBV	41282	53398	2.18%	0.188%
42	11.920	1759	1766	1767	rBV	46115	59141	2.42%	0.208%
43	11.945	1767	1770	1777	rVB	72351	88581	3.62%	0.311%
44	12.079	1786	1792	1800	rVV	1665444	2075527	84.76%	7.291%
45	12.298	1824	1828	1835	rBV2	224512	387225	15.81%	1.360%
46	12.365	1835	1839	1847	rVB2	171378	280721	11.46%	0.986%
47	12.457	1848	1854	1859	rBV2	30707	44758	1.83%	0.157%
48	12.518	1859	1864	1869	rVB	85993	109989	4.49%	0.386%
49	12.585	1870	1875	1878	rBV	145537	188053	7.68%	0.661%
50	12.664	1882	1888	1891	rBV	133659	161480	6.59%	0.567%
51	12.701	1891	1894	1899	rVB	81617	98009	4.00%	0.344%
52	12.755	1899	1903	1905	rBV	174350	257247	10.51%	0.904%
53	12.896	1922	1926	1932	rVB3	53728	85670	3.50%	0.301%
54	12.975	1932	1939	1943	rBV	226230	319594	13.05%	1.123%
55	13.018	1943	1946	1951	rVV2	94466	112336	4.59%	0.395%
56	13.079	1951	1956	1963	rVV2	190441	295121	12.05%	1.037%
57	13.152	1964	1968	1971	rVV2	61466	105404	4.30%	0.370%
58	13.188	1971	1974	1977	rVV	148483	208820	8.53%	0.734%
59	13.225	1977	1980	1983	rVV	177102	242378	9.90%	0.851%
60	13.255	1983	1985	1990	rVV4	59634	94749	3.87%	0.333%
61	13.310	1990	1994	1997	rVV2	153895	221746	9.06%	0.779%
62	13.347	1997	2000	2008	rVV3	346260	559643	22.85%	1.966%
63	13.426	2008	2013	2016	rVV	173505	251079	10.25%	0.882%
64	13.481	2019	2022	2030	rVB3	53958	87527	3.57%	0.307%
65	13.560	2030	2035	2038	rBV2	165948	239725	9.79%	0.842%
66	13.597	2038	2041	2044	rVV	339192	440711	18.00%	1.548%
67	13.639	2044	2048	2052	rVV3	300798	561901	22.95%	1.974%
68	13.694	2052	2057	2058	rVV	190047	328083	13.40%	1.153%
69	13.719	2058	2061	2067	rVV2	450736	650727	26.57%	2.286%
70	13.804	2072	2075	2079	rVV	65768	86651	3.54%	0.304%
71	13.853	2079	2083	2088	rVV5	57776	115325	4.71%	0.405%

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72	13.908	2088	2092	2097	rVB	256528	340142	13.89%	1.195%
73	13.981	2097	2104	2108	rBV2	296029	434931	17.76%	1.528%
74	14.030	2108	2112	2115	rVV	139088	204489	8.35%	0.718%
75	14.060	2115	2117	2124	rVB3	71056	88414	3.61%	0.311%
76	14.133	2124	2129	2133	rBV	283936	356701	14.57%	1.253%
77	14.182	2133	2137	2140	rVV3	24846	45006	1.84%	0.158%
78	14.219	2140	2143	2146	rVV	39266	47438	1.94%	0.167%
79	14.286	2146	2154	2160	rVV4	298143	663197	27.08%	2.330%
80	14.341	2160	2163	2166	rVV	49536	66566	2.72%	0.234%
81	14.377	2166	2169	2173	rVV	133040	168262	6.87%	0.591%
82	14.432	2173	2178	2182	rVV	320174	417080	17.03%	1.465%
83	14.475	2182	2185	2190	rVB2	46535	64918	2.65%	0.228%
84	14.536	2190	2195	2200	rBV2	187548	288609	11.79%	1.014%
85	14.578	2200	2202	2208	rVB3	32448	48336	1.97%	0.170%
86	14.670	2213	2217	2224	rVV2	275787	524546	21.42%	1.843%
87	14.731	2224	2227	2231	rVB	177655	221388	9.04%	0.778%
88	14.786	2231	2236	2241	rBV2	77098	96485	3.94%	0.339%
89	14.847	2241	2246	2248	rBV2	44497	72181	2.95%	0.254%
90	14.877	2248	2251	2253	rVV3	88263	114024	4.66%	0.401%
91	14.901	2253	2255	2259	rVB2	81073	103500	4.23%	0.364%
92	14.962	2259	2265	2271	rBV4	72711	154081	6.29%	0.541%
93	15.109	2282	2289	2296	rBV5	23987	48198	1.97%	0.169%
94	15.249	2306	2312	2319	rBV2	192775	350042	14.29%	1.230%
95	15.322	2319	2324	2330	rVB4	37128	71292	2.91%	0.250%
96	15.444	2339	2344	2346	rBV2	33925	52266	2.13%	0.184%
97	15.474	2346	2349	2353	rVB	47251	63999	2.61%	0.225%
98	15.596	2364	2369	2374	rVB5	33952	63208	2.58%	0.222%
99	15.682	2379	2383	2387	rVV5	32949	57196	2.34%	0.201%
100	15.724	2387	2390	2397	rVB5	25248	49221	2.01%	0.173%

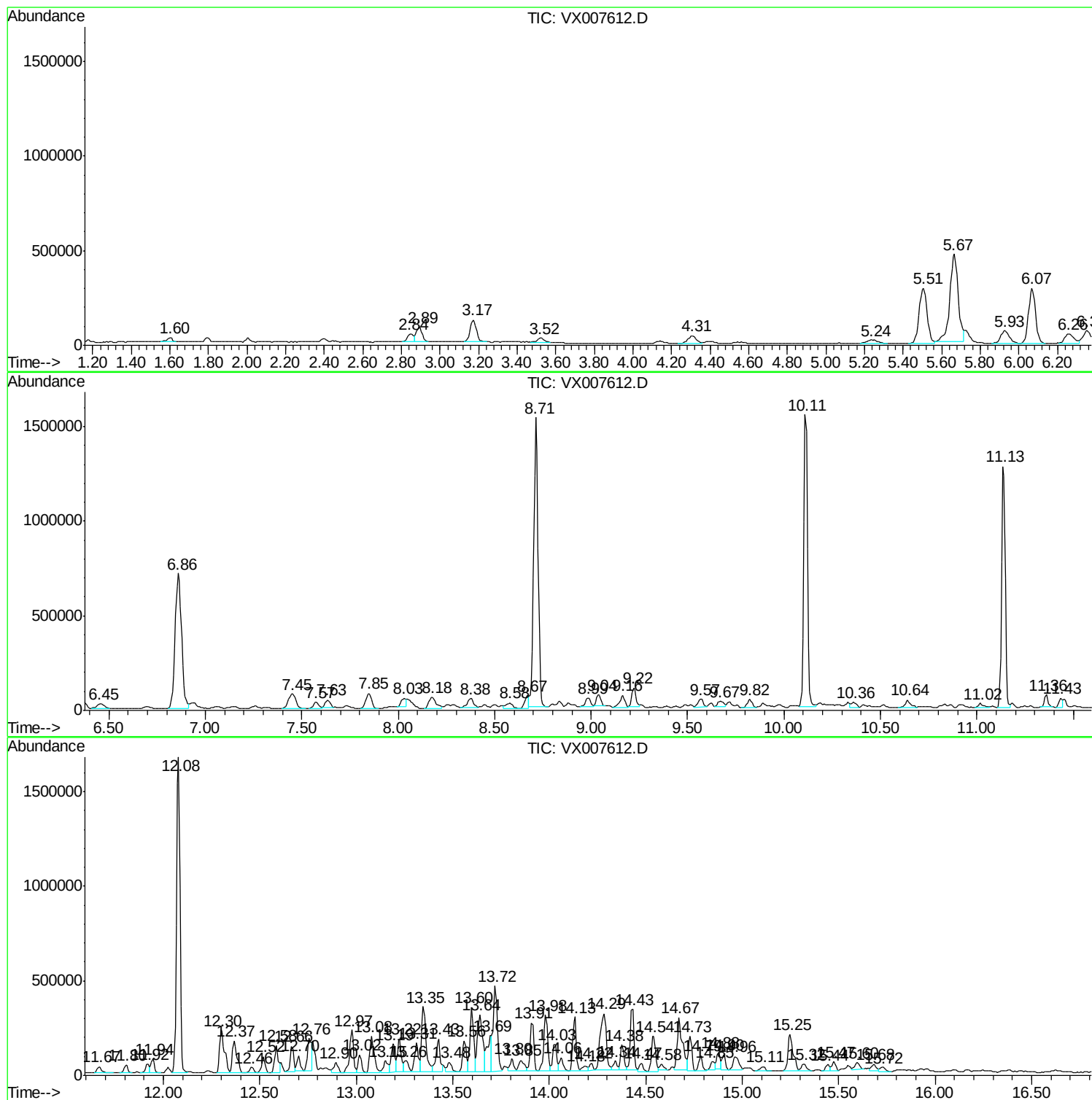
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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



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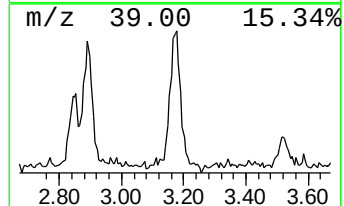
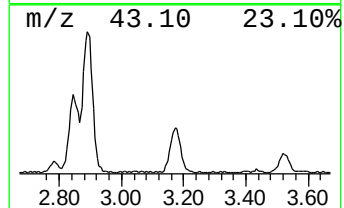
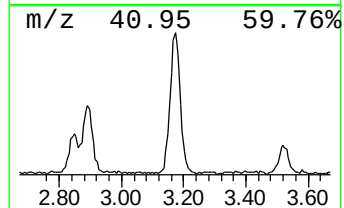
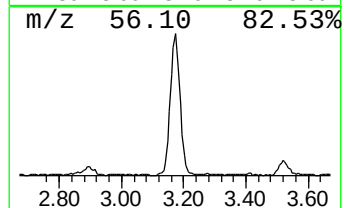
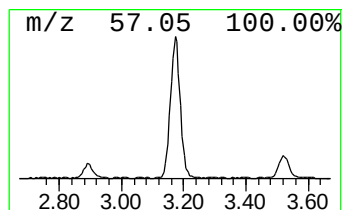
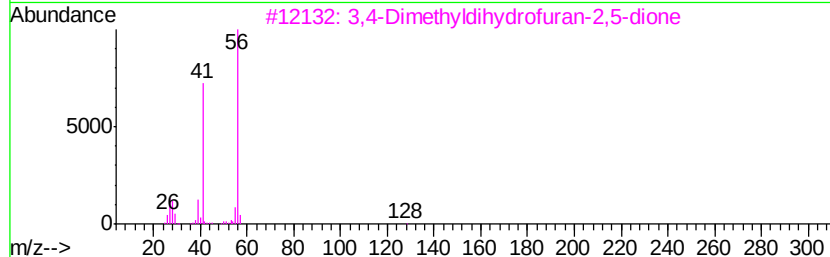
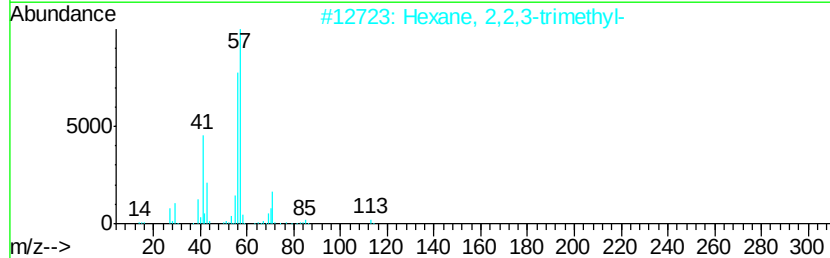
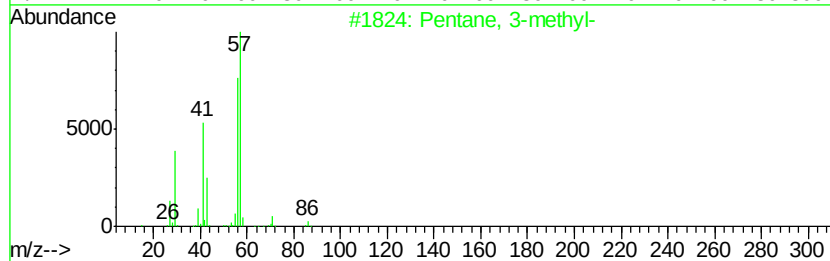
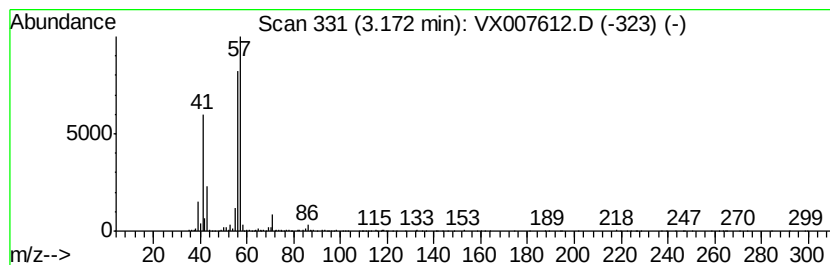
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.17	9.30 ug/l	263072	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 3-methyl-	86	C6H14	000096-14-0	86
2		Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	56
3		3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	53
4		Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	50
5		Butane, 1-chloro-	92	C4H9Cl	000109-69-3	42



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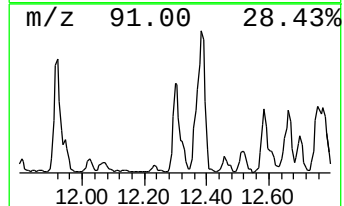
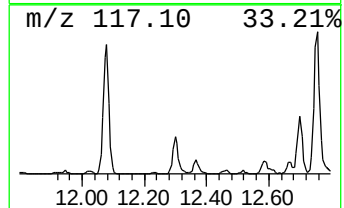
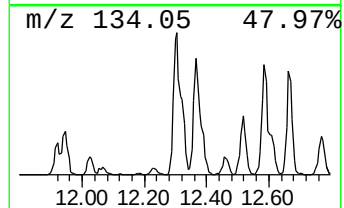
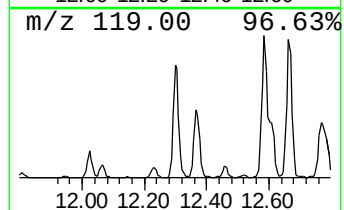
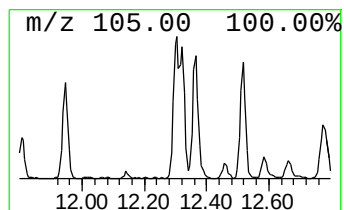
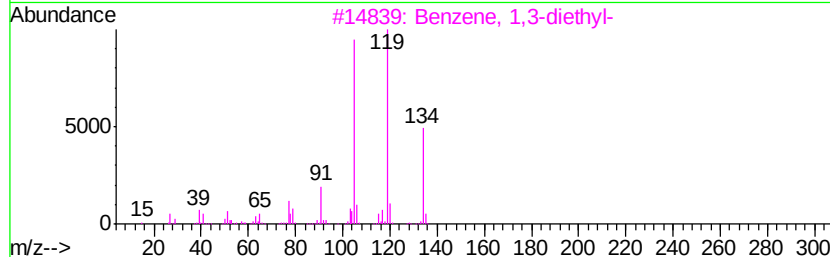
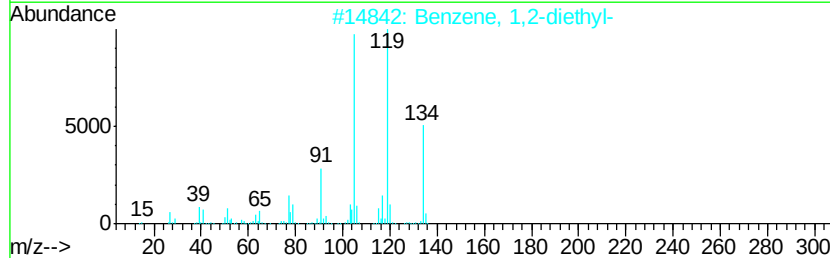
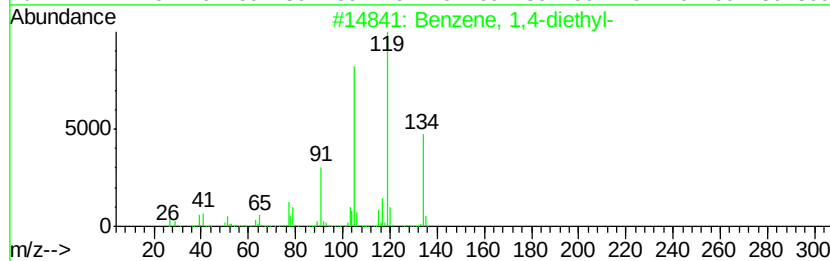
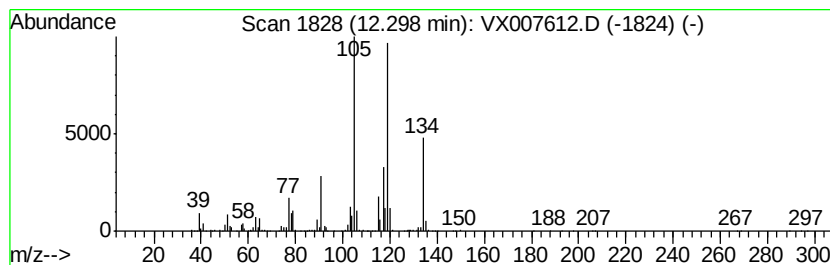
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Benzene, 1,4-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	9.33 ug/l	387225	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
2		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
3		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	87
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	83
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	72



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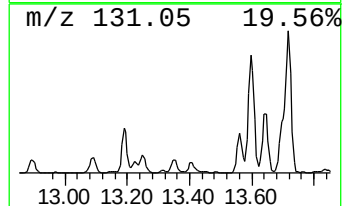
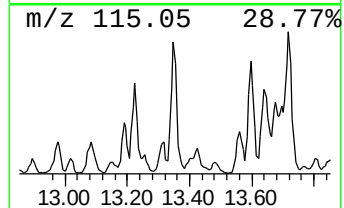
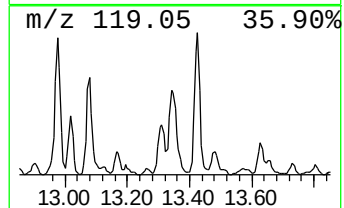
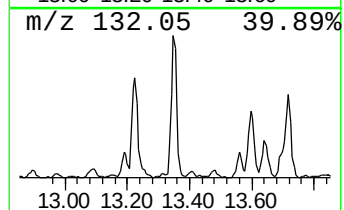
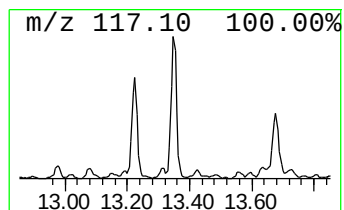
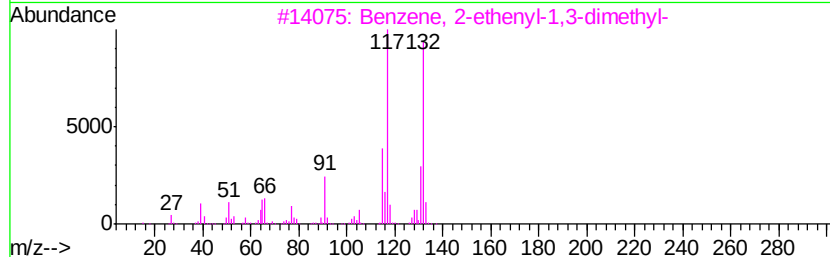
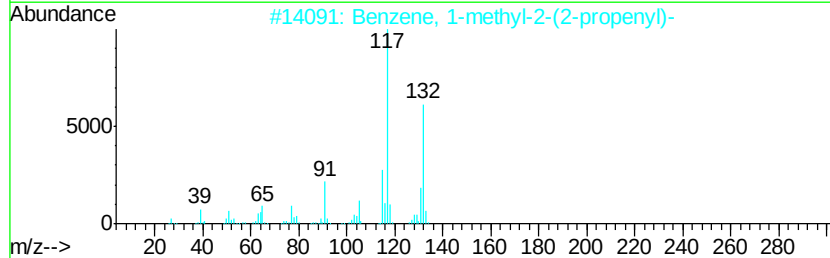
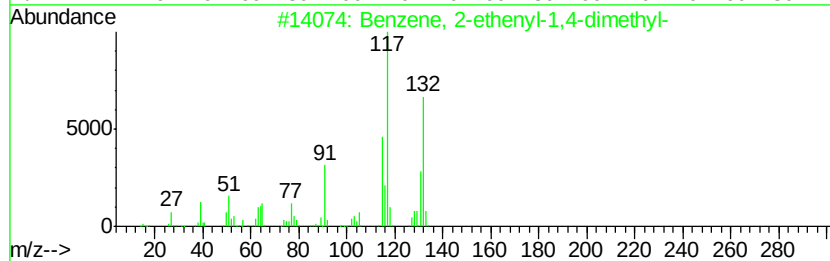
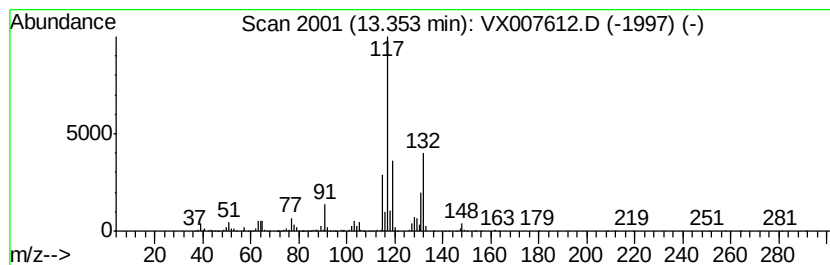
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Peak Number 3 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	13.48 ug/l	559643	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
2		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	76
3		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	76
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
5		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021919\
Data File : VX007612.D
Acq On : 19 Feb 2019 14:44
Operator : JC/SP
Sample : K1528-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

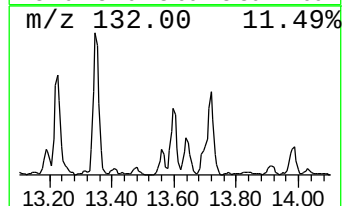
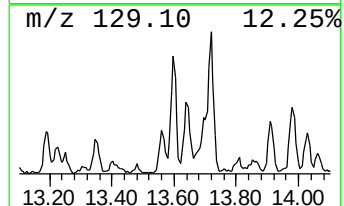
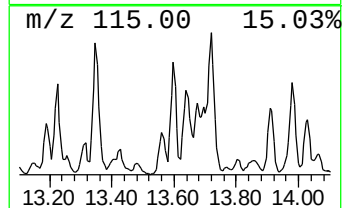
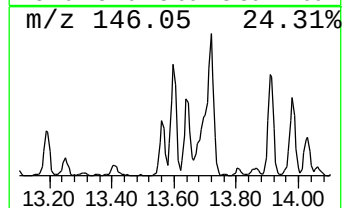
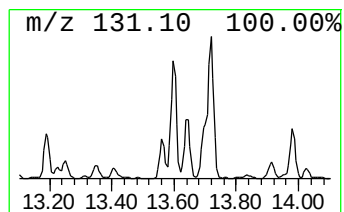
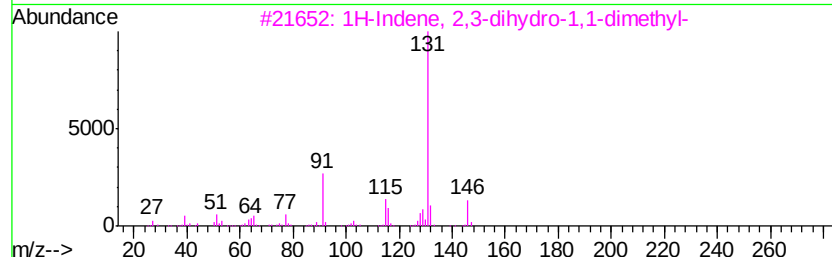
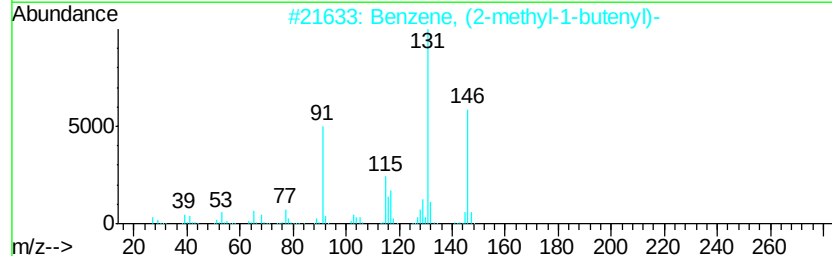
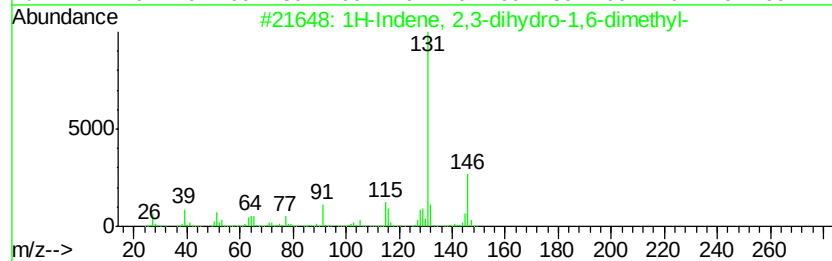
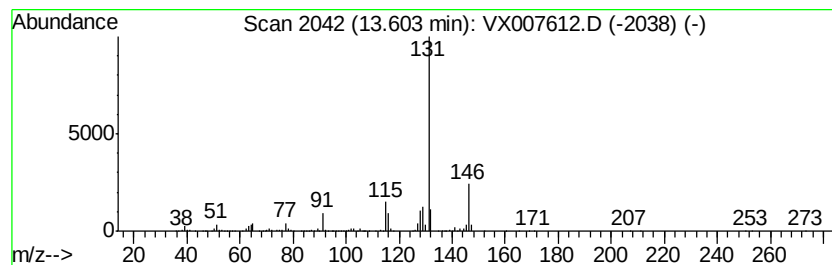
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Quant Title : SW846 8260

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

Peak Number 4 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.60	10.62 ug/l	440711	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14		017059-48-2	96
2	Benzene, (2-methyl-1-butenyl)-	146	C11H14		056253-64-6	94
3	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14		004912-92-9	91
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14		004175-53-5	91
5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14		017057-82-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021919\
Data File : VX007612.D
Acq On : 19 Feb 2019 14:44
Operator : JC/SP
Sample : K1528-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-3

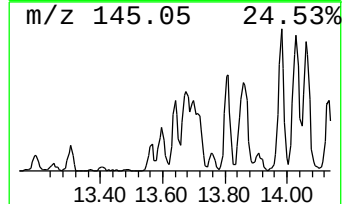
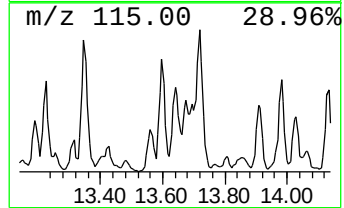
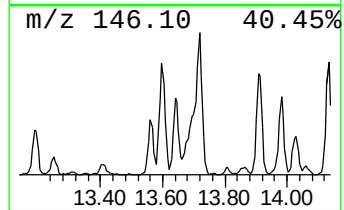
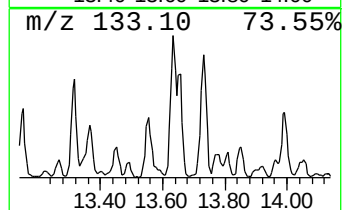
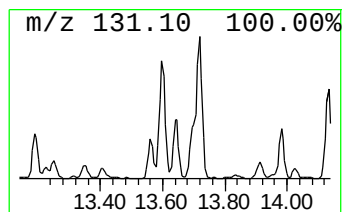
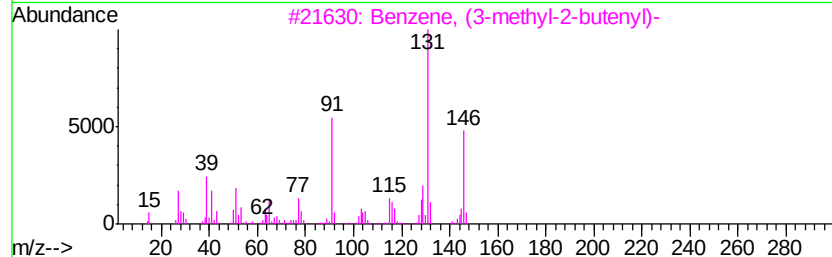
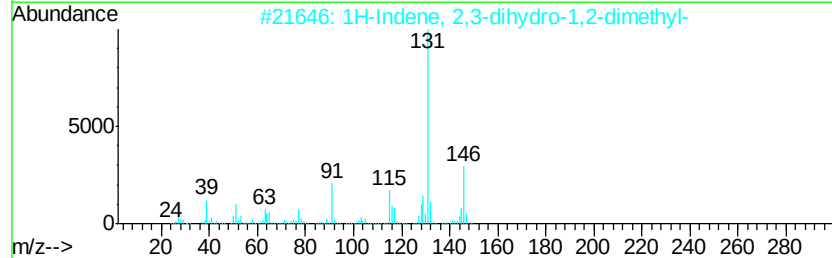
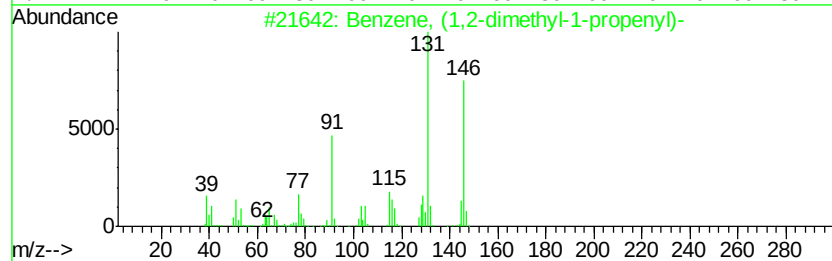
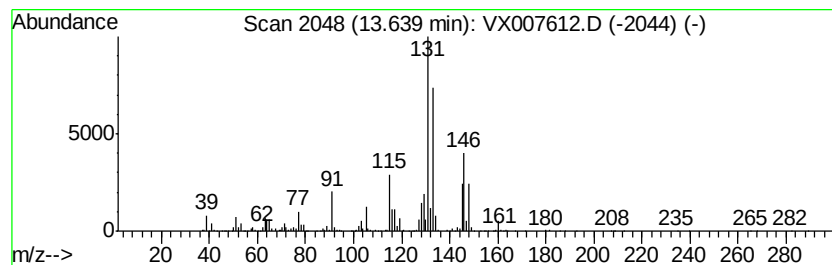
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.64	13.54 ug/l	561901	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	90
2		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	78
3		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	60
4		1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	60
5		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021919\
Data File : VX007612.D
Acq On : 19 Feb 2019 14:44
Operator : JC/SP
Sample : K1528-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

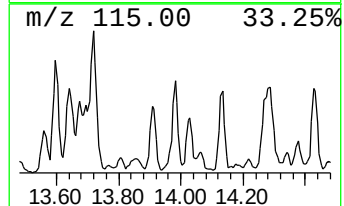
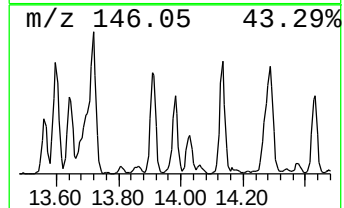
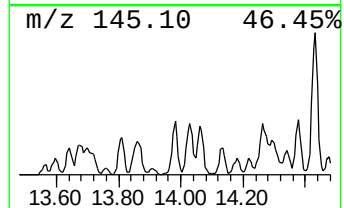
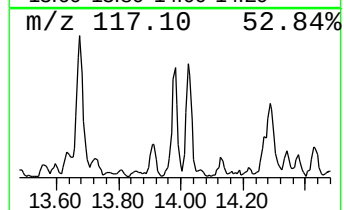
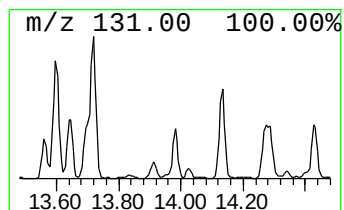
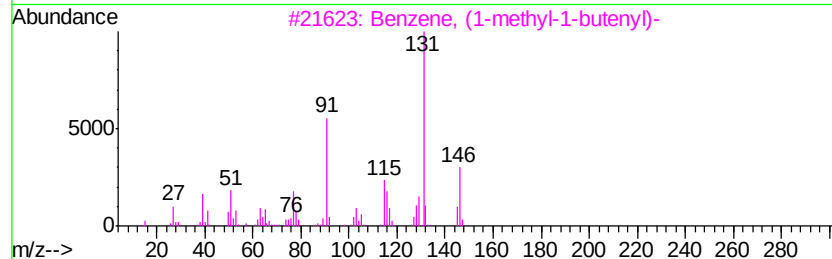
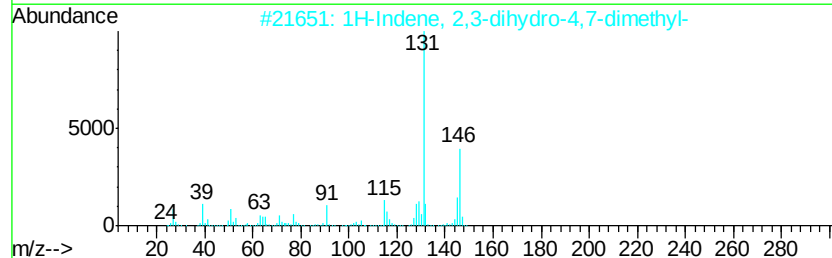
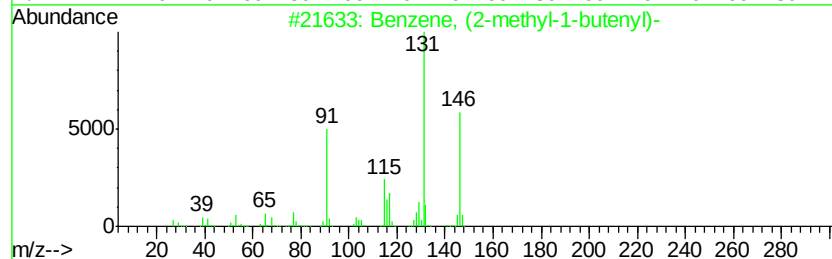
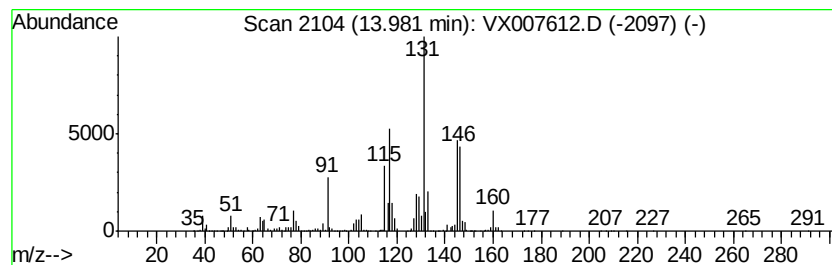
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.98	10.48 ug/l	434931	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	60
2		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	60
3		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	60
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	160	C12H16	078920-29-3	60
5		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021919\
Data File : VX007612.D
Acq On : 19 Feb 2019 14:44
Operator : JC/SP
Sample : K1528-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

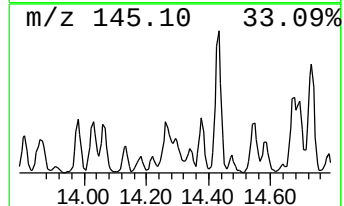
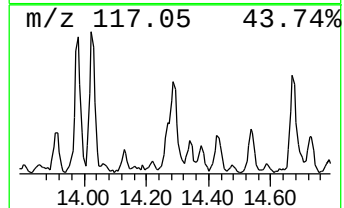
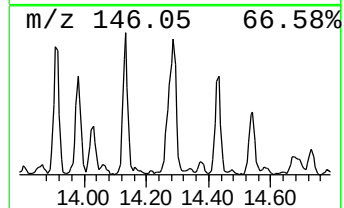
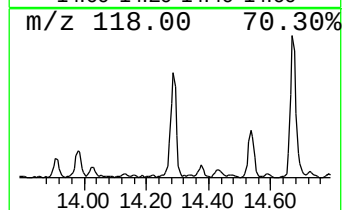
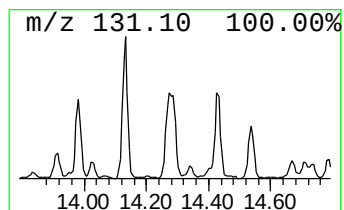
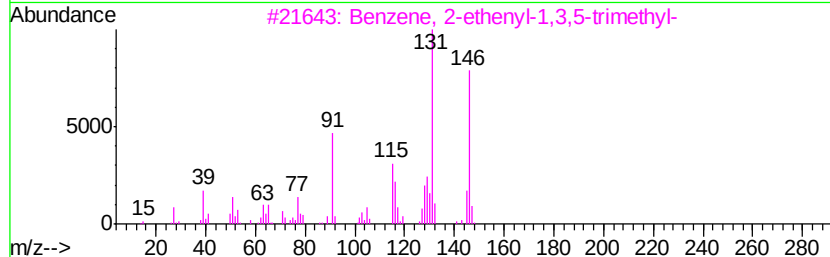
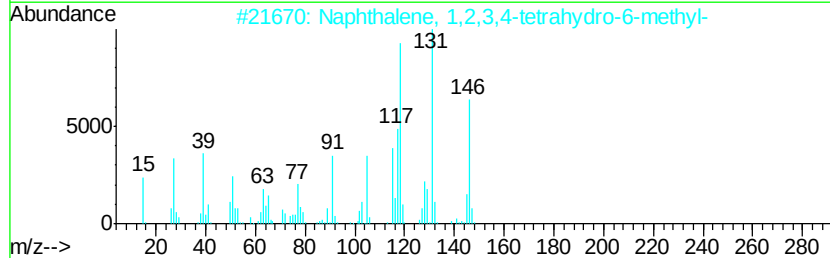
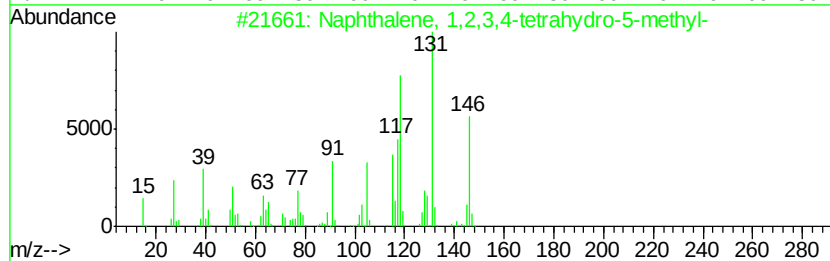
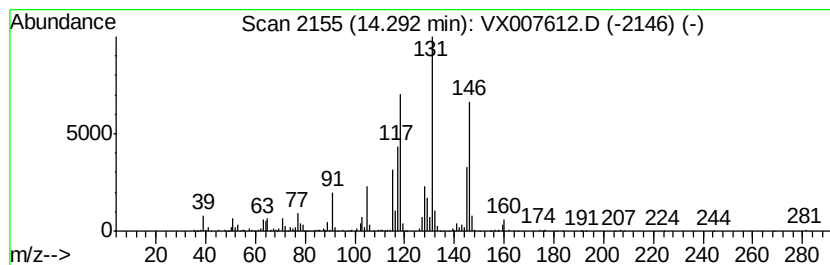
Instrument :
MSVOA_X
ClientSampleId :
MW-3

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	15.98 ug/l	663197	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 95
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 94
3		Benzene, 2-ethenyl-1,3,5-trimethyl-	146 C11H14	000769-25-5 86
4		2,3,4,5,6,7-Hexahydro-1H-cyclope...	146 C11H14	1000189-31-0 76
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021919\
Data File : VX007612.D
Acq On : 19 Feb 2019 14:44
Operator : JC/SP
Sample : K1528-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

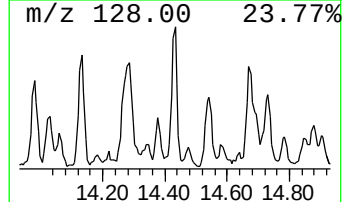
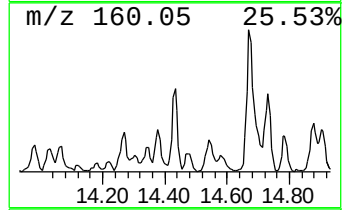
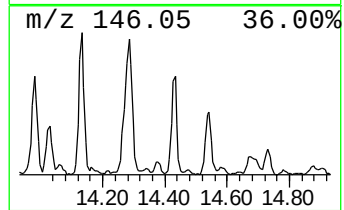
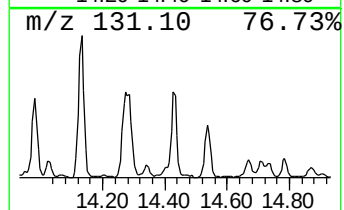
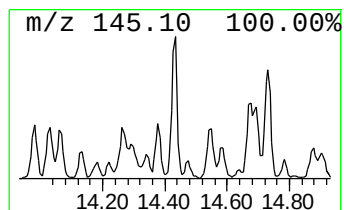
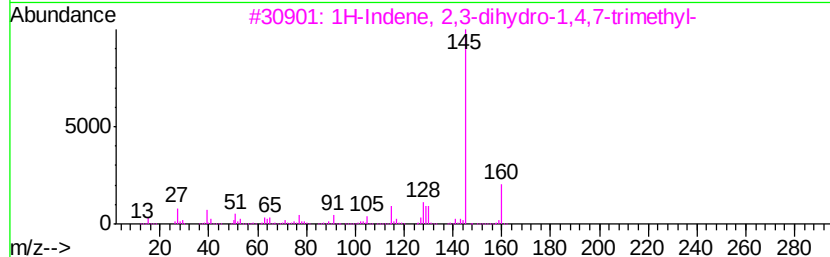
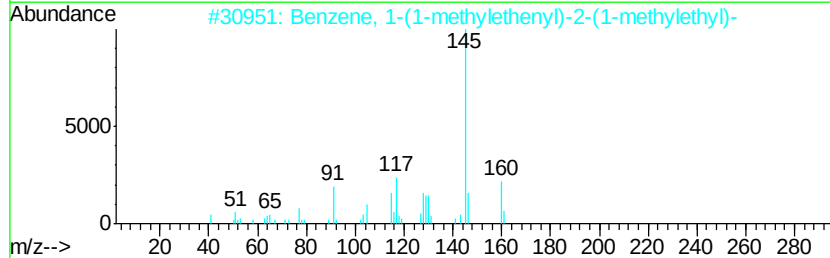
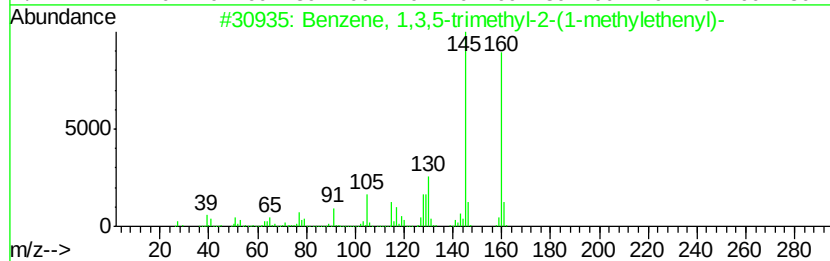
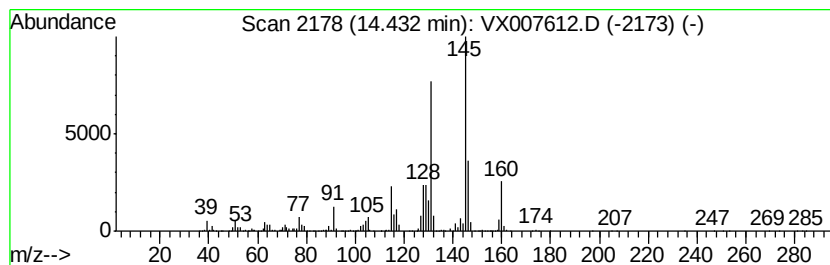
Instrument :
MSVOA_X
ClientSampleId :
MW-3

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,3,5-trimethyl-2-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.43	10.05 ug/l	417080	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,3,5-trimethyl-2-(1-me...	160 C12H16	014679-13-1 55
2		Benzene, 1-(1-methylethenyl)-2-(...	160 C12H16	005557-93-7 53
3		1H-Indene, 2,3-dihydro-1,4,7-tri...	160 C12H16	054340-87-3 53
4		Benzene, 4-(2-butenyl)-1,2-dimet...	160 C12H16	054340-86-2 46
5		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 45



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX021919\
Data File : VX007612.D
Acq On : 19 Feb 2019 14:44
Operator : JC/SP
Sample : K1528-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

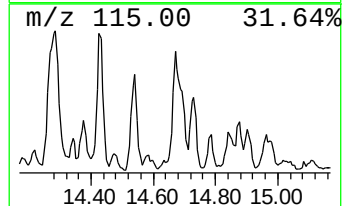
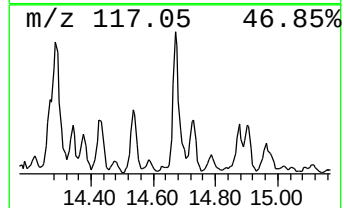
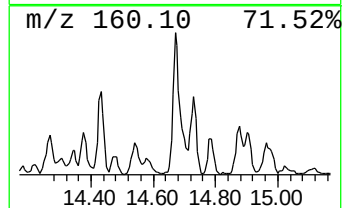
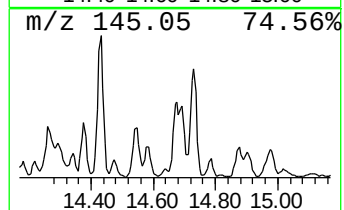
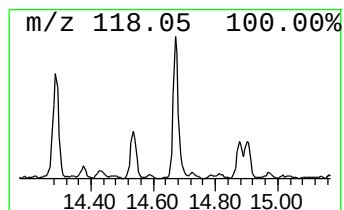
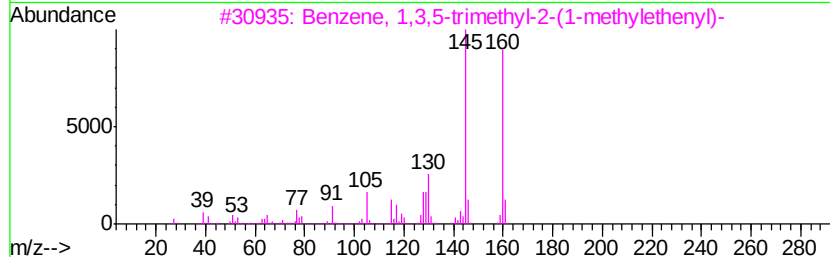
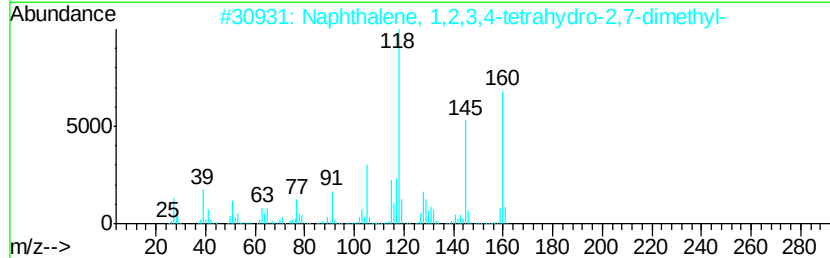
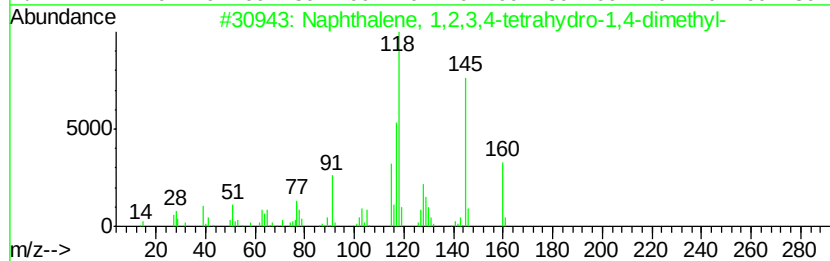
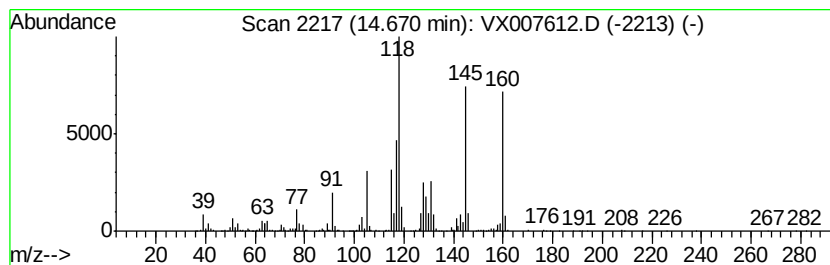
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetra... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.67	12.64 ug/l	524546	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	95
2		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	90
3		Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	70
4		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	68
5		1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	014944-23-1	46



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX021919\
Data File : VX007612.D
Acq On : 19 Feb 2019 14:44
Operator : JC/SP
Sample : K1528-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.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
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Benzene, 1,4-diet...	12.30	9.3	ug/l	387225	4	12.08	2075530	50.0
Benzene, 2-etheny...	13.35	13.5	ug/l	559643	4	12.08	2075530	50.0
1H-Indene, 2,3-di...	13.60	10.6	ug/l	440711	4	12.08	2075530	50.0
Benzene, (1,2-dim...	13.64	13.5	ug/l	561901	4	12.08	2075530	50.0
Benzene, (2-methy...	13.98	10.5	ug/l	434931	4	12.08	2075530	50.0
Naphthalene, 1,2,...	14.29	16.0	ug/l	663197	4	12.08	2075530	50.0
Benzene, 1,3,5-tr...	14.43	10.1	ug/l	417080	4	12.08	2075530	50.0
Naphthalene, 1,2,...	14.67	12.6	ug/l	524546	4	12.08	2075530	50.0