

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
 Data File : VX000280.D
 Acq On : 12 Mar 2018 17:19
 Operator : JC/MD
 Sample : J1727-04
 Misc : 6.64g/10mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 T-4

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 : W:\HPCHEM1\MSVOA_X\METHOD\82X030718W.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	1.069	76	79	90	rBV	230635	286054	25.86%	1.370%
2	1.636	158	172	175	rBV5	22538	73190	6.62%	0.351%
3	2.300	274	281	285	rBV	33379	71128	6.43%	0.341%
4	2.386	285	295	302	rVB	80463	225180	20.36%	1.079%
5	5.519	796	809	820	rBV2	114652	349686	31.62%	1.675%
6	5.672	825	834	852	rVB	107657	345566	31.24%	1.655%
7	6.080	890	901	915	rVB2	141683	368055	33.28%	1.763%
8	6.342	934	944	954	rVV2	46078	143551	12.98%	0.688%
9	6.873	1018	1031	1049	rBV	169257	445665	40.29%	2.135%
10	7.458	1117	1127	1137	rBV3	27870	75937	6.87%	0.364%
11	8.055	1216	1225	1237	rVB2	58823	132047	11.94%	0.633%
12	8.177	1237	1245	1253	rBV3	45667	105267	9.52%	0.504%
13	8.342	1267	1272	1276	rBV2	32111	55593	5.03%	0.266%
14	8.507	1294	1299	1303	rBV	48596	81939	7.41%	0.393%
15	8.677	1319	1327	1330	rBV3	50045	103867	9.39%	0.498%
16	8.726	1330	1335	1343	rVB	608366	1006383	90.99%	4.821%
17	8.976	1370	1376	1383	rBV3	29582	57864	5.23%	0.277%
18	9.451	1447	1454	1463	rBV2	27666	50048	4.52%	0.240%
19	9.561	1467	1472	1479	rBV2	54900	113764	10.29%	0.545%
20	9.884	1520	1525	1529	rVV2	26119	59001	5.33%	0.283%
21	9.927	1529	1532	1534	rVV2	36012	51609	4.67%	0.247%
22	9.964	1534	1538	1544	rVB	125772	185591	16.78%	0.889%
23	10.067	1549	1555	1559	rBV	106508	156128	14.12%	0.748%
24	10.122	1559	1564	1574	rBV	352877	534121	48.29%	2.559%
25	10.262	1581	1587	1591	rVB	81426	116957	10.57%	0.560%
26	10.366	1591	1604	1610	rBV	195607	325084	29.39%	1.557%
27	10.421	1610	1613	1624	rVB5	49815	109945	9.94%	0.527%
28	10.646	1645	1650	1654	rBV	66643	89513	8.09%	0.429%
29	10.713	1657	1661	1662	rVV	37993	48627	4.40%	0.233%
30	10.732	1662	1664	1669	rVV	57068	64299	5.81%	0.308%
31	10.841	1674	1682	1686	rBV2	79414	131178	11.86%	0.628%
32	10.921	1691	1695	1699	rVB2	44529	56652	5.12%	0.271%
33	11.024	1706	1712	1717	rBV2	32948	59614	5.39%	0.286%
34	11.091	1717	1723	1726	rBV2	68819	98672	8.92%	0.473%

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35	11.146	1726	1732	1742	rVB3	673403	1106071	100.00%	5.299%
36	11.250	1746	1749	1753	rVB	94112	98620	8.92%	0.472%
37	11.372	1765	1769	1776	rVB	96315	140548	12.71%	0.673%
38	11.445	1776	1781	1787	rBV	154637	283171	25.60%	1.357%
39	11.518	1787	1793	1803	rVB2	179738	354209	32.02%	1.697%
40	11.677	1815	1819	1826	rBV2	95924	180093	16.28%	0.863%
41	11.744	1826	1830	1833	rBV2	44525	51999	4.70%	0.249%
42	11.774	1833	1835	1838	rBV2	52294	53685	4.85%	0.257%
43	11.817	1838	1842	1851	rVB	393424	562699	50.87%	2.696%
44	11.896	1851	1855	1858	rBV2	57774	82673	7.47%	0.396%
45	11.951	1862	1864	1868	rVB2	58915	71232	6.44%	0.341%
46	12.000	1868	1872	1875	rBV3	34262	50963	4.61%	0.244%
47	12.091	1880	1887	1891	rBV2	487358	675413	61.06%	3.236%
48	12.158	1891	1898	1902	rBV3	136182	244481	22.10%	1.171%
49	12.231	1908	1910	1915	rVB2	84431	97587	8.82%	0.468%
50	12.311	1915	1923	1930	rBV3	257880	501770	45.37%	2.404%
51	12.378	1930	1934	1941	rVB3	285745	490006	44.30%	2.347%
52	12.481	1946	1951	1955	rBV2	71635	108504	9.81%	0.520%
53	12.530	1955	1959	1965	rVB	100435	144521	13.07%	0.692%
54	12.597	1965	1970	1972	rBV	110791	153136	13.85%	0.734%
55	12.622	1972	1974	1976	rVV	63594	63040	5.70%	0.302%
56	12.677	1979	1983	1987	rVV	473952	544106	49.19%	2.607%
57	12.768	1991	1998	2006	rVB3	163589	376707	34.06%	1.805%
58	12.859	2006	2013	2015	rBV5	31726	66940	6.05%	0.321%
59	12.987	2025	2034	2038	rBV	365186	559922	50.62%	2.682%
60	13.030	2038	2041	2046	rVB	224142	270494	24.46%	1.296%
61	13.091	2046	2051	2056	rBV3	106252	240912	21.78%	1.154%
62	13.176	2061	2065	2068	rBV2	123897	150630	13.62%	0.722%
63	13.231	2071	2074	2078	rVB2	149106	189597	17.14%	0.908%
64	13.323	2083	2089	2091	rVV4	92456	148122	13.39%	0.710%
65	13.359	2091	2095	2100	rVV4	552630	772212	69.82%	3.699%
66	13.439	2104	2108	2113	rBV2	162803	222466	20.11%	1.066%
67	13.487	2113	2116	2124	rVB7	90836	204125	18.45%	0.978%
68	13.573	2124	2130	2132	rBV2	87496	136566	12.35%	0.654%
69	13.609	2132	2136	2139	rBV	122803	154229	13.94%	0.739%
70	13.646	2139	2142	2149	rVV3	245649	371621	33.60%	1.780%
71	13.737	2150	2157	2162	rVB3	160938	270046	24.41%	1.294%

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72	13.841	2171	2174	2180	rVB	72811	116607	10.54%	0.559%
73	13.920	2180	2187	2192	rVB5	79945	210151	19.00%	1.007%
74	13.993	2192	2199	2203	rBV4	137236	261888	23.68%	1.255%
75	14.036	2203	2206	2210	rBV	63431	88231	7.98%	0.423%
76	14.140	2219	2223	2230	rBV	189565	267596	24.19%	1.282%
77	14.225	2235	2237	2242	rVV2	52368	71061	6.42%	0.340%
78	14.280	2242	2246	2256	rVV2	219556	397552	35.94%	1.905%
79	14.390	2260	2264	2268	rVV	123934	164793	14.90%	0.789%
80	14.438	2268	2272	2276	rVB	149316	182277	16.48%	0.873%
81	14.487	2276	2280	2285	rVB3	35464	61039	5.52%	0.292%
82	14.548	2285	2290	2295	rBV2	92164	162132	14.66%	0.777%
83	14.682	2308	2312	2313	rVV	153510	181479	16.41%	0.869%
84	14.707	2313	2316	2319	rVV	303855	392885	35.52%	1.882%
85	14.743	2319	2322	2325	rVB	67968	85120	7.70%	0.408%
86	14.792	2326	2330	2333	rBV4	46401	70706	6.39%	0.339%
87	14.847	2335	2339	2343	rVV	235853	310985	28.12%	1.490%
88	14.883	2343	2345	2348	rVV3	50543	67545	6.11%	0.324%
89	14.914	2348	2350	2356	rVV4	53093	80454	7.27%	0.385%
90	14.975	2356	2360	2367	rVV4	31169	75064	6.79%	0.360%
91	15.030	2367	2369	2377	rVB6	26750	58210	5.26%	0.279%
92	15.261	2402	2407	2410	rBV	91000	147588	13.34%	0.707%
93	15.457	2434	2439	2442	rBV2	65125	95002	8.59%	0.455%
94	15.560	2451	2456	2461	rVV	147921	223573	20.21%	1.071%
95	15.700	2474	2479	2482	rBV	165434	252111	22.79%	1.208%
96	15.737	2482	2485	2494	rVB	106814	172297	15.58%	0.825%
97	15.932	2509	2517	2527	rVB3	39575	107543	9.72%	0.515%
98	16.456	2599	2603	2611	rVB3	32130	67088	6.07%	0.321%
99	16.706	2640	2644	2649	rVB	46238	80843	7.31%	0.387%
100	16.767	2649	2654	2657	rBV3	31613	51122	4.62%	0.245%

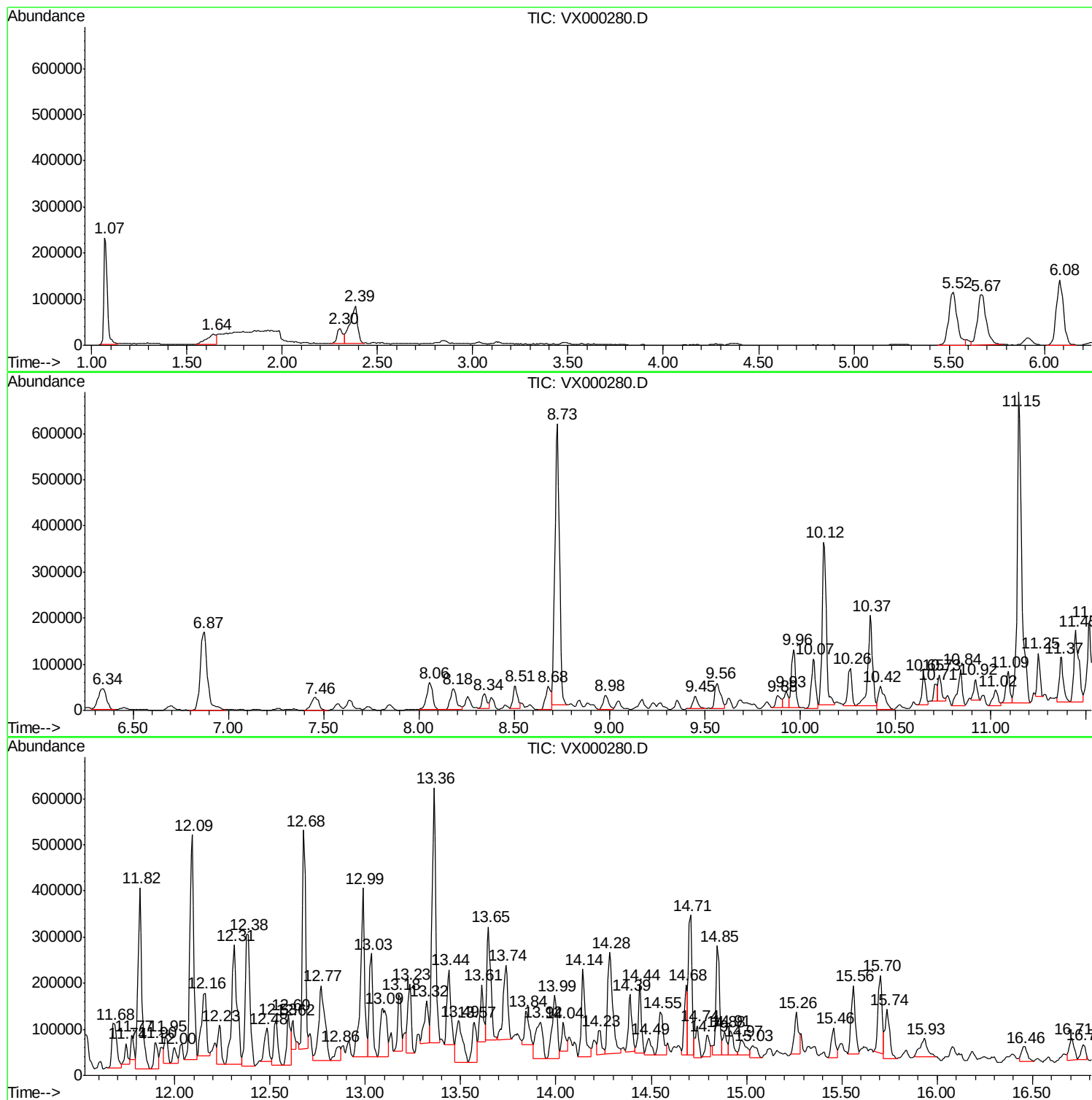
Sum of corrected areas: 20873833

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Instrument :
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ClientSampled :
T-4

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X030718W.M
Quant Title : SW846 8260

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



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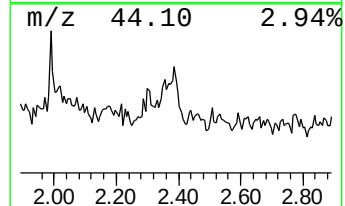
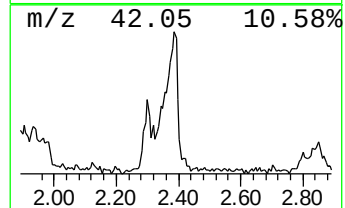
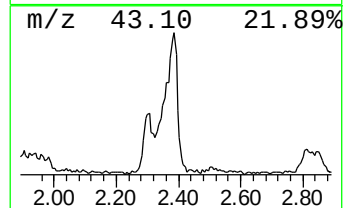
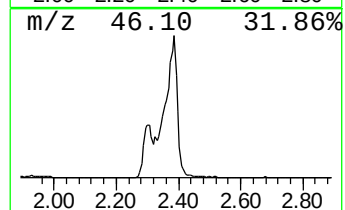
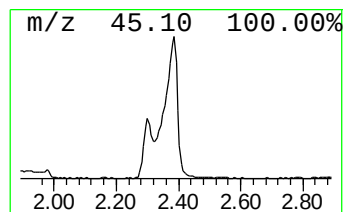
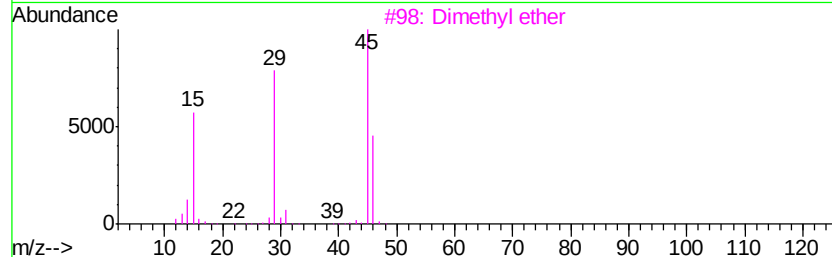
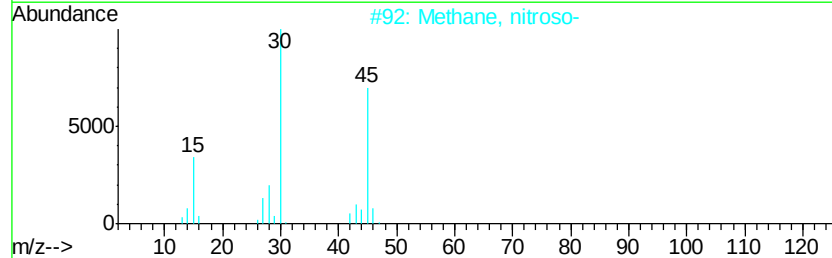
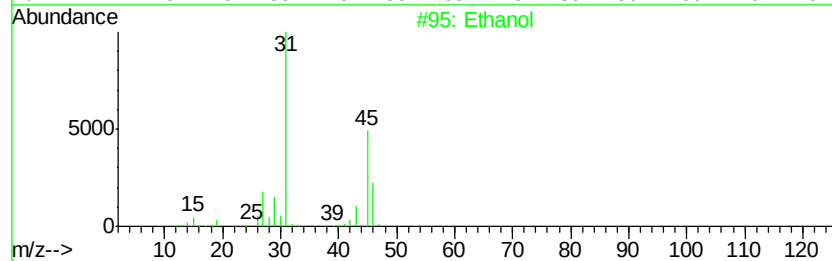
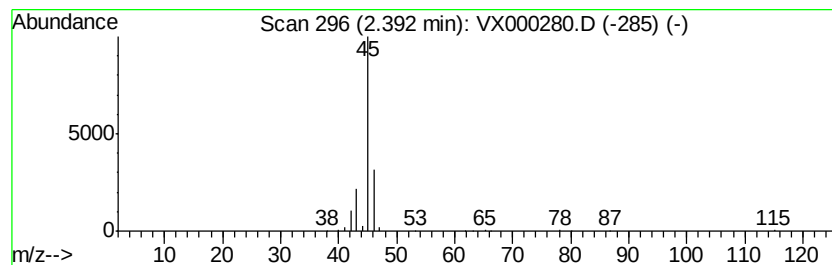
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X030718W.M
Quant Title : SW846 8260

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

Peak Number 2 unknown2.39 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.39	32.58 ug/l	225180	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol	46	C2H6O	000064-17-5	43
2		Methane, nitroso-	45	CH3NO	000865-40-7	4
3		Dimethyl ether	46	C2H6O	000115-10-6	4
4		Formic acid	46	CH2O2	000064-18-6	4
5		Propanedioic acid, dihydroxy-	136	C3H4O6	000560-27-0	4



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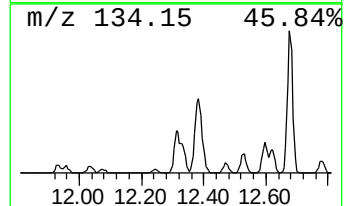
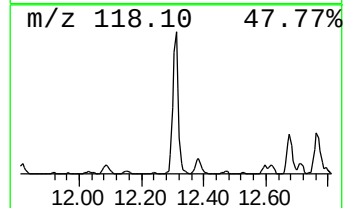
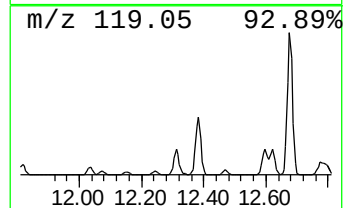
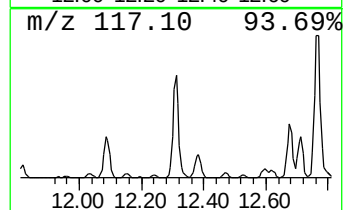
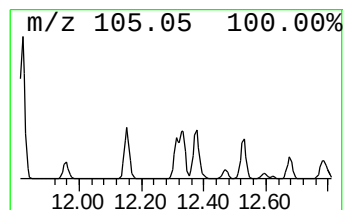
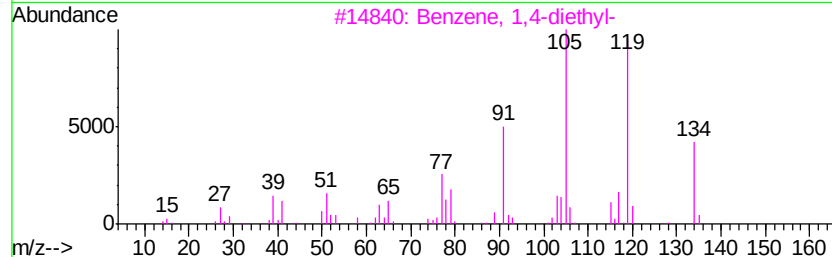
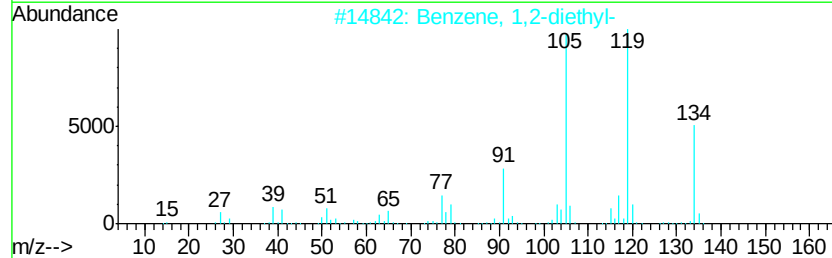
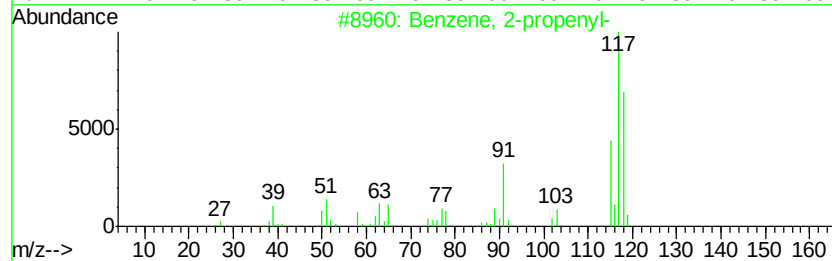
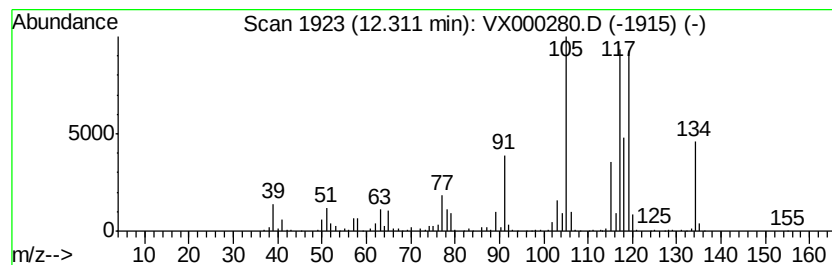
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Benzene, 2-propenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.31	37.15 ug/l	501770	1,4-Dichlorobenzene-d4	12.09		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
2		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	55
3		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	55
4		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	55
5		Benzene, cyclopropyl-	118	C9H10	000873-49-4	52



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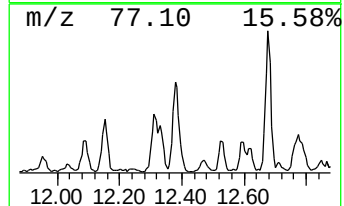
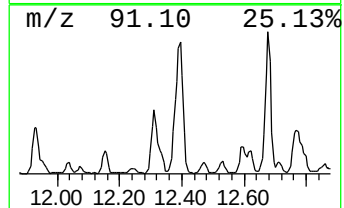
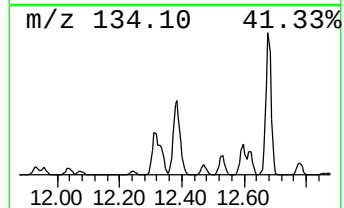
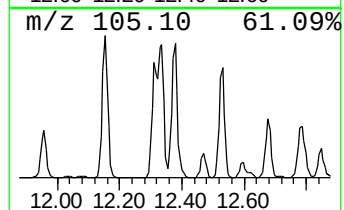
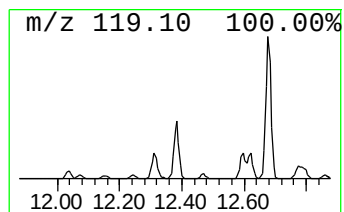
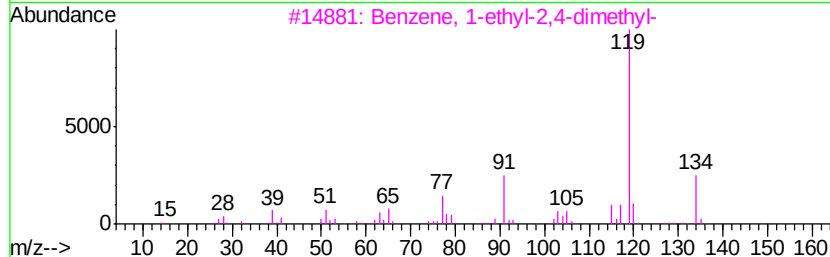
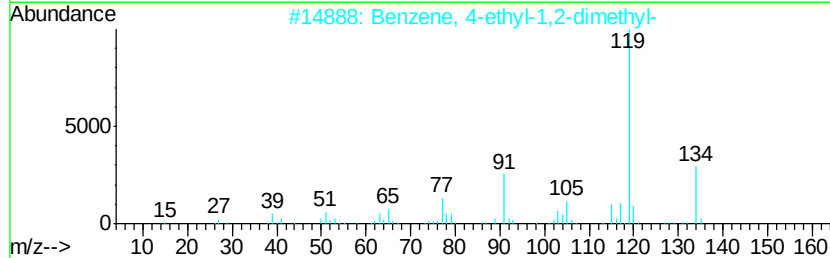
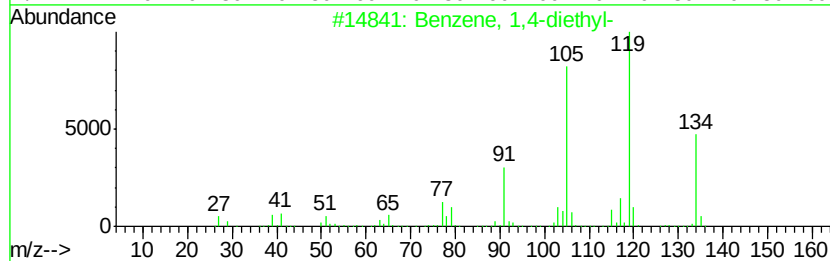
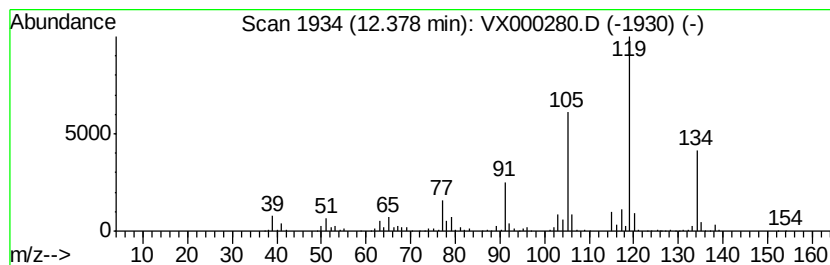
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Peak Number 4 Benzene, 1,4-diethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.38	36.27 ug/l	490006	1,4-Dichlorobenzene-d4	12.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,4-diethyl-	134 C10H14	000105-05-5 95
2		Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5 95
3		Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9 94
4		Benzene, 1,3-diethyl-	134 C10H14	000141-93-5 94
5		o-Cymene	134 C10H14	000527-84-4 93



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
Data File : VX000280.D
Acq On : 12 Mar 2018 17:19
Operator : JC/MD
Sample : J1727-04
Misc : 6.64g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
T-4

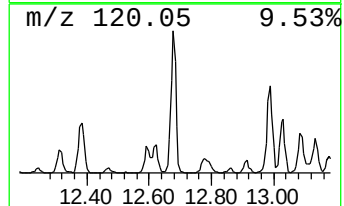
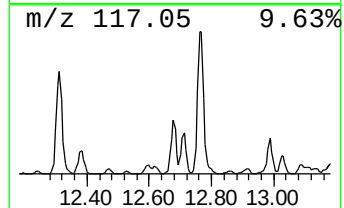
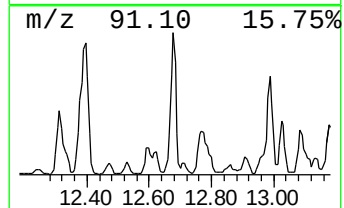
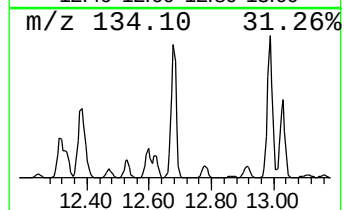
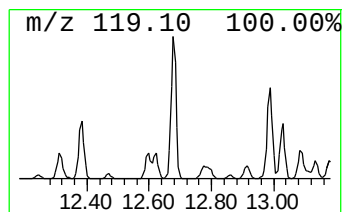
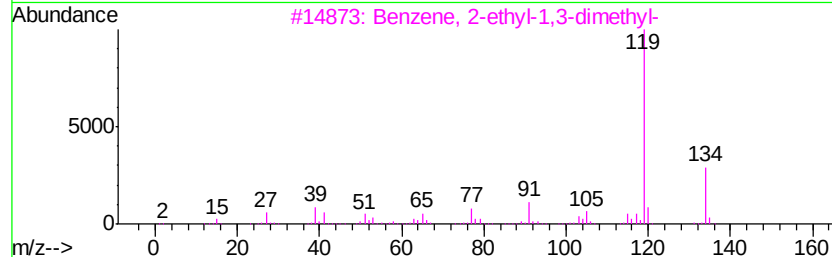
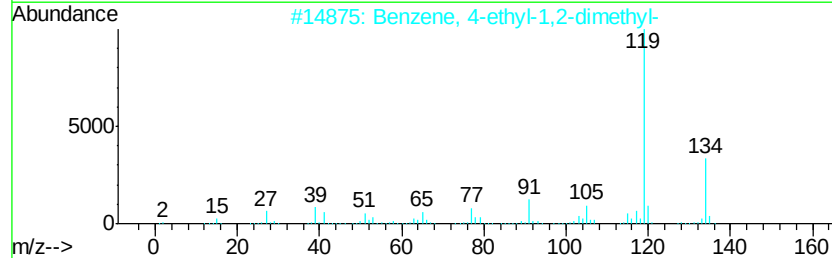
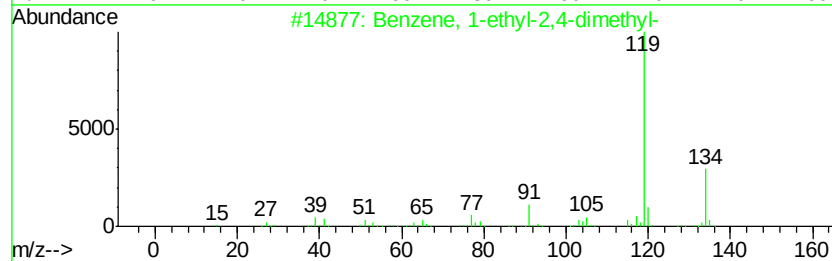
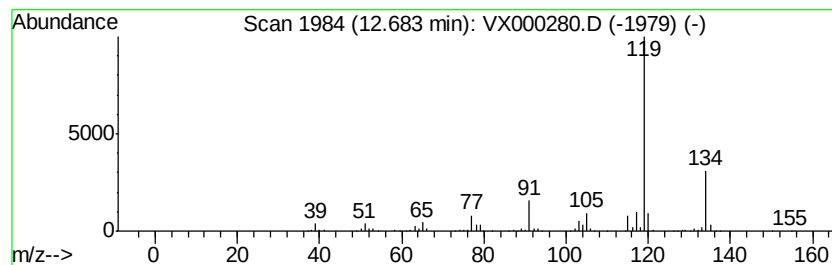
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.68	40.28 ug/l	544106	1,4-Dichlorobenzene-d4	12.09

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
Data File : VX000280.D
Acq On : 12 Mar 2018 17:19
Operator : JC/MD
Sample : J1727-04
Misc : 6.64g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
T-4

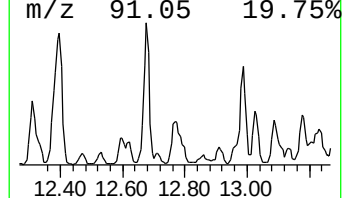
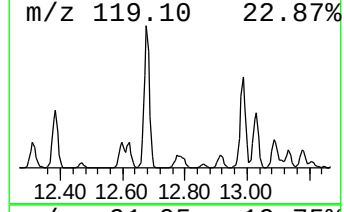
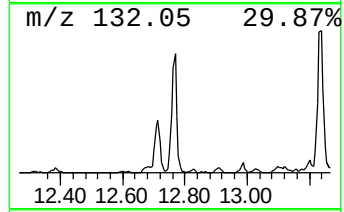
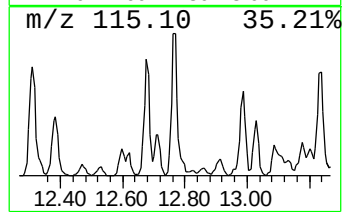
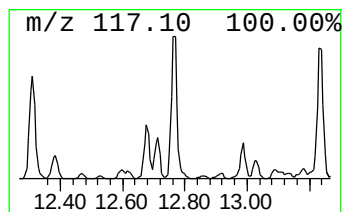
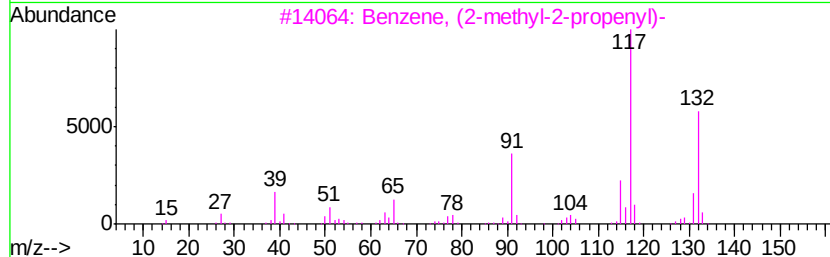
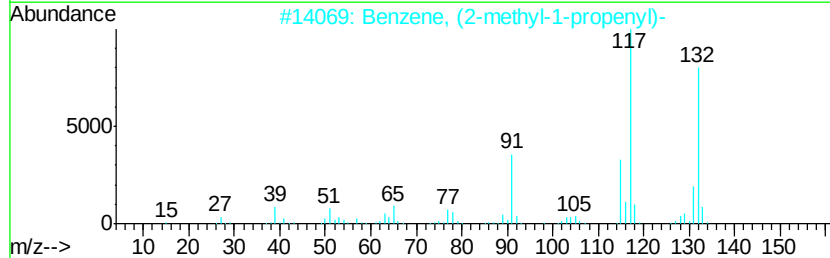
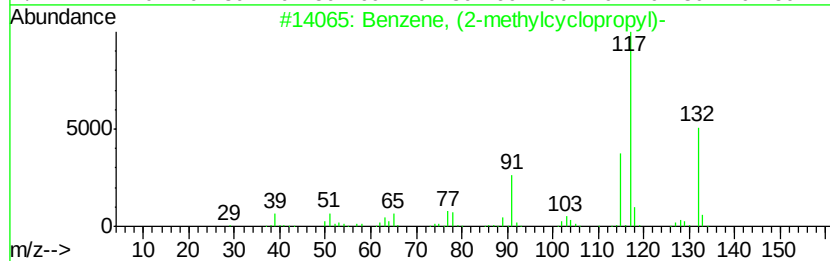
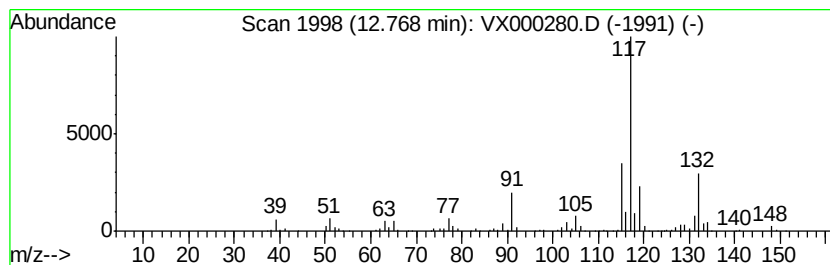
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X030718W.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, (2-methylcycloprop... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	27.89 ug/l	376707	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	93
2		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
3		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	83
4		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
5		Indan, 1-methyl-	132	C10H12	000767-58-8	76



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
Data File : VX000280.D
Acq On : 12 Mar 2018 17:19
Operator : JC/MD
Sample : J1727-04
Misc : 6.64µL/10mL/100µL/5.0mL/MSVOA_X/MEOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T-4

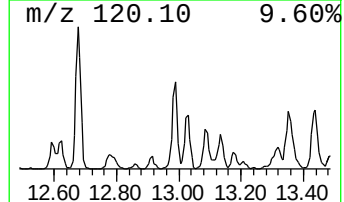
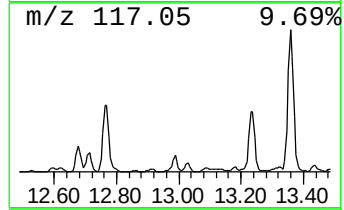
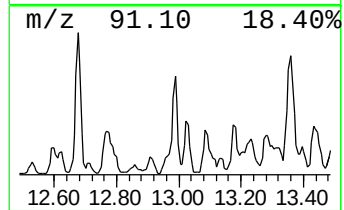
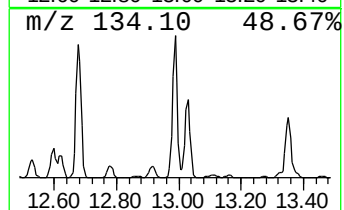
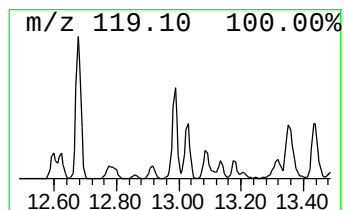
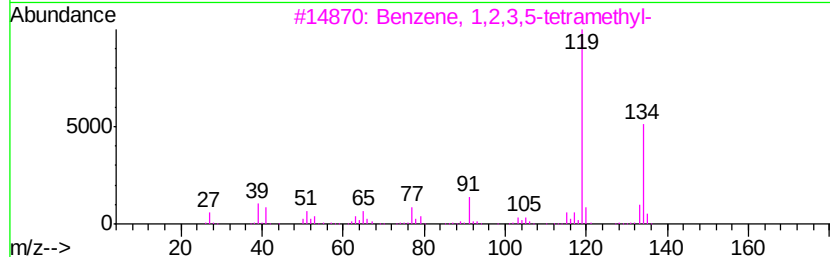
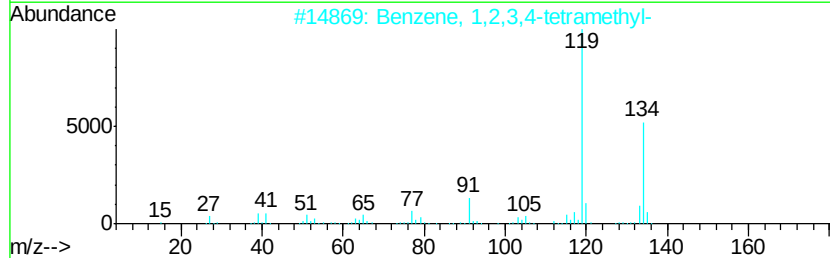
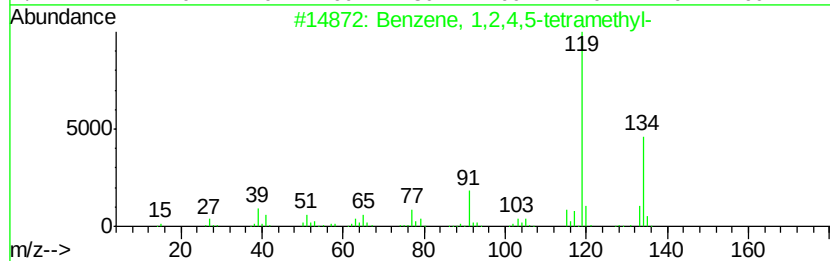
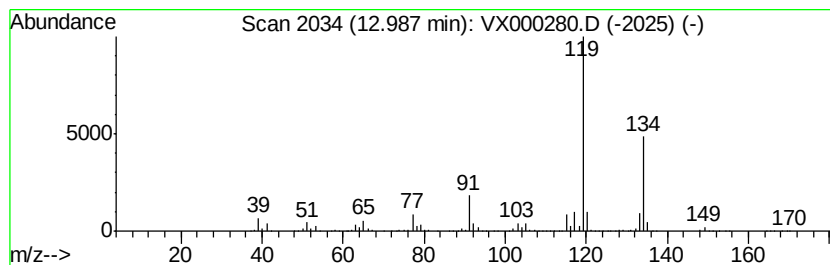
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X030718W.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.99	41.45 ug/l	559922	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	93



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
Data File : VX000280.D
Acq On : 12 Mar 2018 17:19
Operator : JC/MD
Sample : J1727-04
Misc : 6.64g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 1 Sample Multiplier: 1

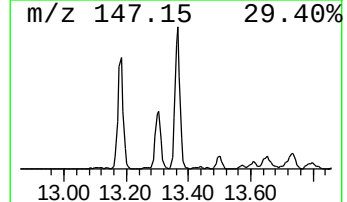
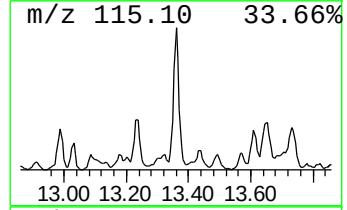
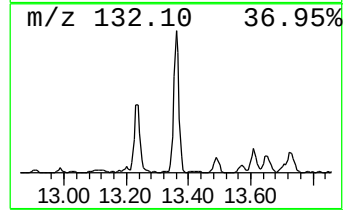
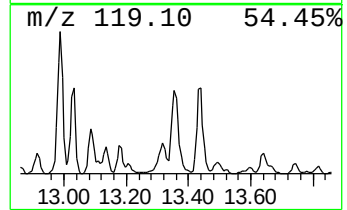
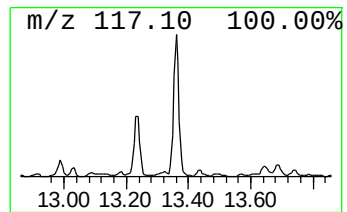
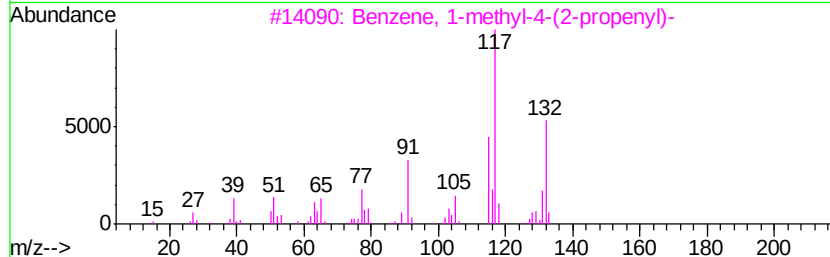
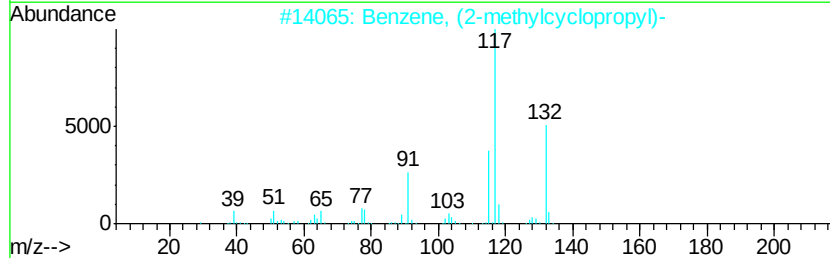
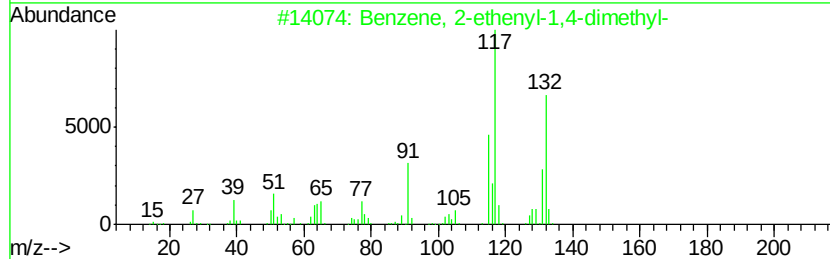
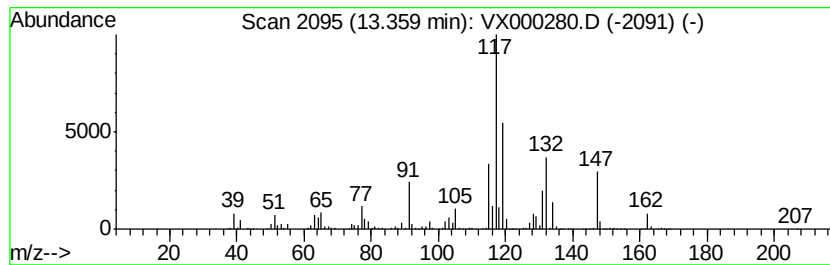
Instrument :
MSVOA_X
ClientSampled :
T-4

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X030718W.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.36	57.17 ug/l	772212	1,4-Dichlorobenzene-d4	12.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 2-ethenyl-1,4-dimethyl-	132 C10H12	002039-89-6 89
2		Benzene, (2-methylcyclopropyl)-	132 C10H12	1000327-39-0 87
3		Benzene, 1-methyl-4-(2-propenyl)-	132 C10H12	003333-13-9 83
4		2,4-Dimethylstyrene	132 C10H12	002234-20-0 83
5		Benzene, 2-butenyl-	132 C10H12	001560-06-1 83



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
Data File : VX000280.D
Acq On : 12 Mar 2018 17:19
Operator : JC/MD
Sample : J1727-04
Misc : 6.64g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T-4

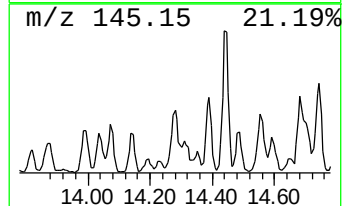
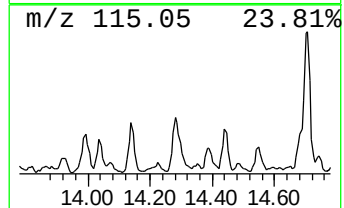
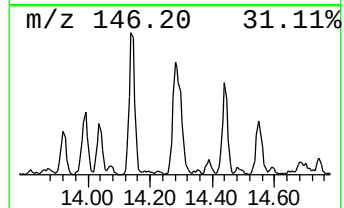
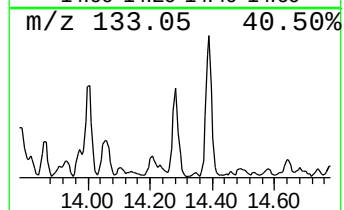
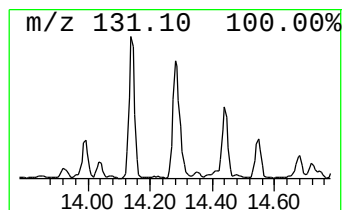
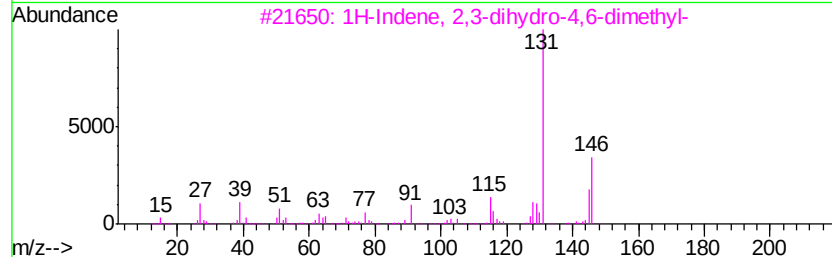
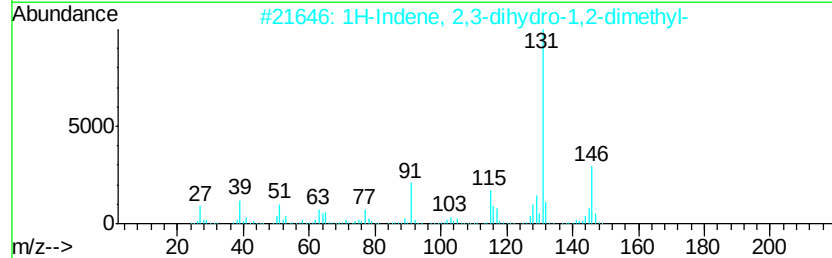
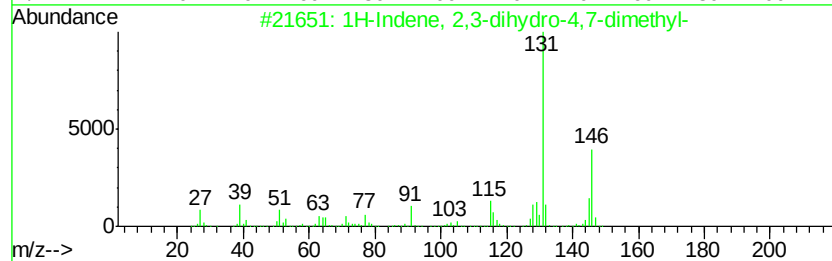
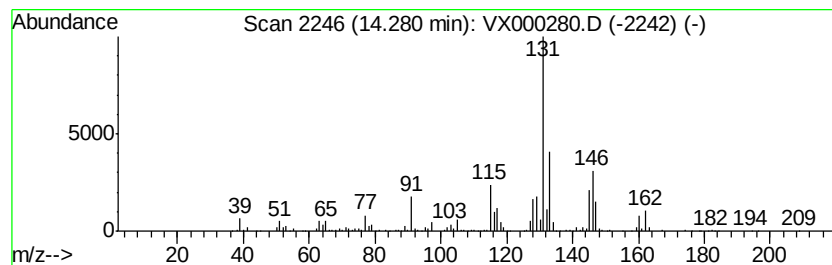
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X030718W.M
Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.28	29.43 ug/l	397552	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	64
2		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	64
3		1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	60
4		1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	60
5		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	58



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
Data File : VX000280.D
Acq On : 12 Mar 2018 17:19
Operator : JC/MD
Sample : J1727-04
Misc : 6.64g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
T-4

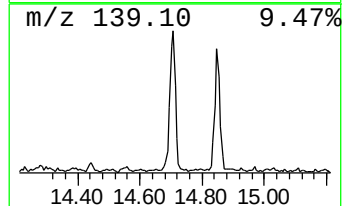
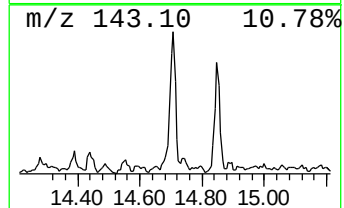
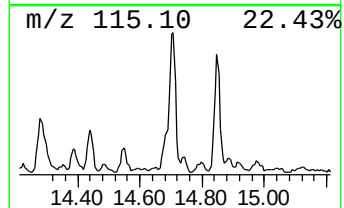
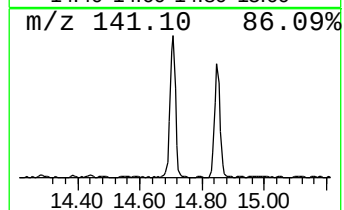
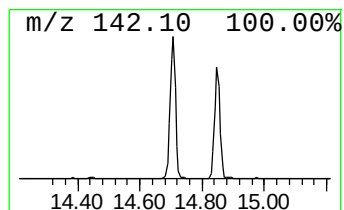
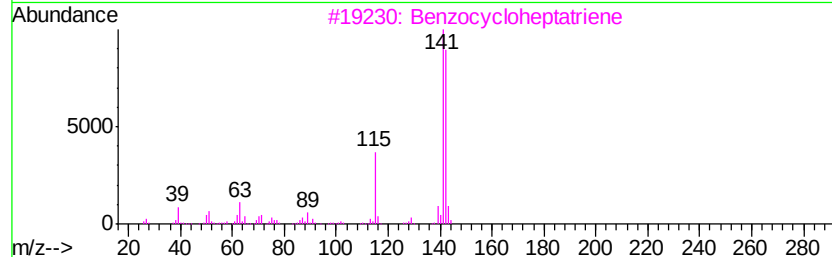
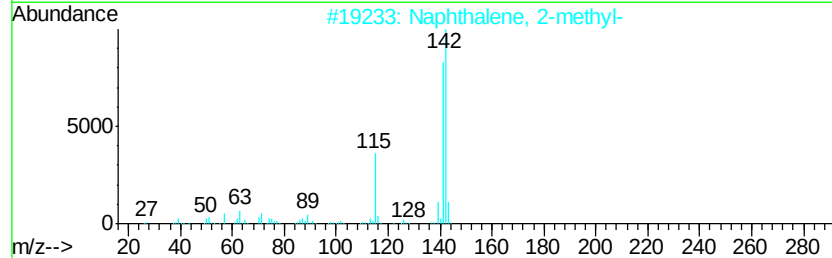
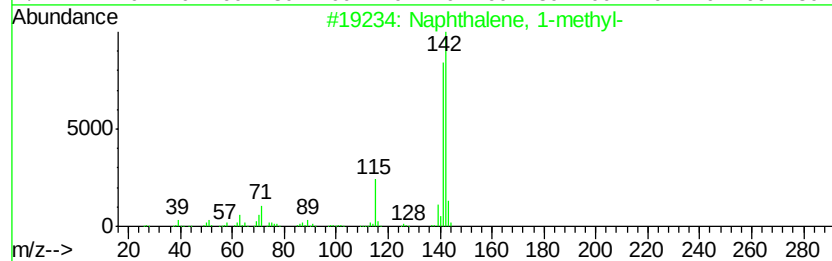
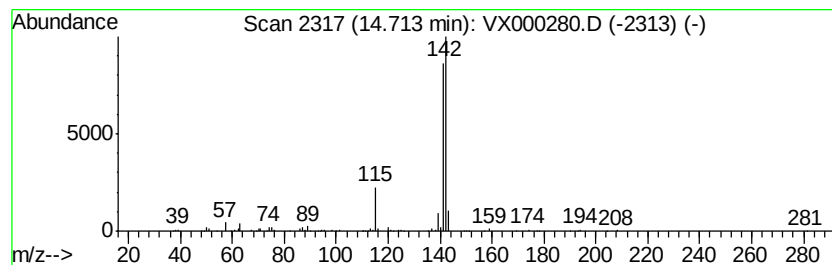
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X030718W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.71	29.08 ug/l	392885	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
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	78



Data Path : W:\HPCHEM1\MSVOA_X\DATA\VX031218\
Data File : VX000280.D
Acq On : 12 Mar 2018 17:19
Operator : JC/MD
Sample : J1727-04
Misc : 6.64g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T-4

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X030718W.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, 2-propenyl-	12.31	37.1	ug/l	501770	4	12.09	675413	50.0
Benzene, 1,4-diet...	12.38	36.3	ug/l	490006	4	12.09	675413	50.0
Benzene, 1-ethyl-...	12.68	40.3	ug/l	544106	4	12.09	675413	50.0
Benzene, (2-methy...	12.77	27.9	ug/l	376707	4	12.09	675413	50.0
Benzene, 1,2,4,5-...	12.99	41.5	ug/l	559922	4	12.09	675413	50.0
Benzene, 2-etheny...	13.36	57.2	ug/l	772212	4	12.09	675413	50.0
1H-Indene, 2,3-di...	14.28	29.4	ug/l	397552	4	12.09	675413	50.0
Naphthalene, 1-me...	14.71	29.1	ug/l	392885	4	12.09	675413	50.0