

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
 Data File : VX000287.D
 Acq On : 12 Mar 2018 20:19
 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	75	79	91	rBV	299268	358148	32.18%	1.445%
2	1.636	156	172	175	rBV4	22400	74511	6.69%	0.301%
3	2.343	276	288	295	rVB	212081	350998	31.54%	1.416%
4	5.519	797	809	819	rBV2	101126	309602	27.82%	1.249%
5	5.672	825	834	852	rVB2	177402	540106	48.53%	2.179%
6	5.909	863	873	884	rVB3	18395	54499	4.90%	0.220%
7	6.080	891	901	919	rBV	113917	310340	27.88%	1.252%
8	6.336	934	943	953	rVV2	55760	163018	14.65%	0.658%
9	6.873	1019	1031	1049	rBV	279078	714360	64.18%	2.882%
10	7.458	1117	1127	1135	rVB3	34390	89766	8.07%	0.362%
11	7.635	1150	1156	1167	rVV2	25342	61132	5.49%	0.247%
12	8.055	1212	1225	1236	rVB	68660	148019	13.30%	0.597%
13	8.177	1236	1245	1253	rBV3	53954	121139	10.88%	0.489%
14	8.342	1267	1272	1275	rBV	36731	56184	5.05%	0.227%
15	8.507	1294	1299	1302	rBV2	52982	87419	7.85%	0.353%
16	8.677	1320	1327	1330	rBV4	56483	120932	10.87%	0.488%
17	8.726	1330	1335	1343	rVB	511711	864034	77.63%	3.486%
18	8.976	1370	1376	1383	rBV3	34737	64287	5.78%	0.259%
19	9.451	1443	1454	1463	rBV2	33780	64725	5.82%	0.261%
20	9.561	1467	1472	1479	rBV2	63545	132357	11.89%	0.534%
21	9.884	1520	1525	1529	rVV2	28930	62965	5.66%	0.254%
22	9.927	1529	1532	1534	rVV2	37895	53196	4.78%	0.215%
23	9.964	1534	1538	1544	rVB	147575	213644	19.20%	0.862%
24	10.067	1548	1555	1560	rBV	126311	186954	16.80%	0.754%
25	10.128	1560	1565	1575	rBV	535429	843729	75.81%	3.404%
26	10.262	1581	1587	1592	rVB	91751	132738	11.93%	0.535%
27	10.366	1592	1604	1610	rBV	214172	372291	33.45%	1.502%
28	10.421	1610	1613	1624	rVB5	59327	126144	11.33%	0.509%
29	10.646	1646	1650	1654	rBV	72011	93928	8.44%	0.379%
30	10.732	1657	1664	1669	rBV2	63233	136839	12.29%	0.552%
31	10.841	1674	1682	1686	rBV	88746	151882	13.65%	0.613%
32	10.921	1691	1695	1699	rVB2	52706	66609	5.98%	0.269%
33	11.030	1706	1713	1716	rBV2	37567	69989	6.29%	0.282%
34	11.091	1716	1723	1726	rBV	81539	115352	10.36%	0.465%

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35	11.152	1726	1733	1742	rVB3	599372	1068590	96.01%	4.311%
36	11.250	1746	1749	1753	rVB	111681	118761	10.67%	0.479%
37	11.372	1765	1769	1776	rVB	122238	165727	14.89%	0.669%
38	11.445	1776	1781	1787	rBV	183265	332284	29.85%	1.340%
39	11.518	1787	1793	1803	rVB2	212523	409462	36.79%	1.652%
40	11.683	1815	1820	1826	rBV2	114980	207731	18.66%	0.838%
41	11.744	1826	1830	1833	rBV	45333	54089	4.86%	0.218%
42	11.774	1833	1835	1838	rBV2	59310	62534	5.62%	0.252%
43	11.817	1838	1842	1851	rVB	468709	677414	60.86%	2.733%
44	11.902	1851	1856	1858	rBV	66839	95670	8.60%	0.386%
45	11.957	1862	1865	1868	rVB2	67130	81064	7.28%	0.327%
46	12.091	1881	1887	1892	rBV	862290	1112999	100.00%	4.490%
47	12.158	1892	1898	1902	rBV3	169261	289883	26.05%	1.169%
48	12.238	1908	1911	1915	rVB2	92642	111224	9.99%	0.449%
49	12.311	1915	1923	1930	rBV3	316845	596303	53.58%	2.406%
50	12.384	1930	1935	1941	rVB3	352059	590861	53.09%	2.384%
51	12.481	1946	1951	1955	rBV2	80189	130028	11.68%	0.525%
52	12.530	1955	1959	1965	rVB	121613	175525	15.77%	0.708%
53	12.597	1965	1970	1972	rBV	140218	184593	16.59%	0.745%
54	12.677	1979	1983	1987	rBV	543848	638346	57.35%	2.575%
55	12.768	1992	1998	2006	rVB3	203955	447711	40.23%	1.806%
56	12.859	2006	2013	2015	rBV4	38033	77359	6.95%	0.312%
57	12.914	2019	2022	2025	rVB2	47030	57925	5.20%	0.234%
58	12.987	2025	2034	2038	rBV	445357	694761	62.42%	2.803%
59	13.030	2038	2041	2047	rVB	275096	333778	29.99%	1.347%
60	13.091	2047	2051	2056	rBV3	130041	287332	25.82%	1.159%
61	13.183	2061	2066	2068	rBV2	146608	190135	17.08%	0.767%
62	13.231	2068	2074	2078	rVB3	178791	298674	26.84%	1.205%
63	13.274	2078	2081	2083	rBV3	56393	66393	5.97%	0.268%
64	13.323	2083	2089	2091	rVV4	112108	179557	16.13%	0.724%
65	13.359	2091	2095	2101	rVV3	653193	937891	84.27%	3.784%
66	13.439	2104	2108	2113	rVB2	201729	280099	25.17%	1.130%
67	13.493	2113	2117	2124	rVB6	108381	245723	22.08%	0.991%
68	13.573	2124	2130	2133	rBV3	108404	184511	16.58%	0.744%
69	13.609	2133	2136	2139	rVV	136447	179427	16.12%	0.724%
70	13.646	2139	2142	2148	rVV3	297794	437570	39.31%	1.765%
71	13.737	2150	2157	2163	rVB2	183065	333997	30.01%	1.347%

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72	13.841	2171	2174	2180	rVB	94800	155123	13.94%	0.626%
73	13.908	2180	2185	2192	rVB4	119000	288955	25.96%	1.166%
74	13.993	2192	2199	2203	rBV3	168895	314197	28.23%	1.268%
75	14.036	2203	2206	2209	rBV	76850	95345	8.57%	0.385%
76	14.140	2219	2223	2227	rBV	223363	288757	25.94%	1.165%
77	14.201	2231	2233	2235	rVV3	45378	55146	4.95%	0.222%
78	14.225	2235	2237	2242	rVV2	64965	96196	8.64%	0.388%
79	14.280	2242	2246	2256	rVV3	261795	457996	41.15%	1.848%
80	14.390	2260	2264	2269	rVV	143917	188685	16.95%	0.761%
81	14.438	2269	2272	2277	rVB	169070	192931	17.33%	0.778%
82	14.487	2277	2280	2285	rVB4	46561	73289	6.58%	0.296%
83	14.548	2285	2290	2295	rBV2	110795	204811	18.40%	0.826%
84	14.682	2308	2312	2313	rVV2	201609	235574	21.17%	0.950%
85	14.707	2313	2316	2319	rVV	348529	439339	39.47%	1.772%
86	14.743	2319	2322	2326	rVB2	76804	99026	8.90%	0.399%
87	14.792	2326	2330	2333	rBV4	55974	87336	7.85%	0.352%
88	14.847	2335	2339	2343	rVV	252137	328905	29.55%	1.327%
89	14.890	2343	2346	2348	rVV3	48315	68497	6.15%	0.276%
90	14.914	2348	2350	2357	rVB5	56861	76698	6.89%	0.309%
91	15.261	2402	2407	2409	rBV	105405	163787	14.72%	0.661%
92	15.286	2409	2411	2414	rVB2	85393	88498	7.95%	0.357%
93	15.457	2434	2439	2442	rBV2	75285	114612	10.30%	0.462%
94	15.560	2451	2456	2462	rBV	159244	247813	22.27%	1.000%
95	15.700	2474	2479	2482	rVV	196593	307162	27.60%	1.239%
96	15.737	2482	2485	2493	rVB2	117033	194946	17.52%	0.786%
97	15.932	2514	2517	2522	rVB3	35641	54522	4.90%	0.220%
98	16.182	2553	2558	2563	rBV3	33468	66530	5.98%	0.268%
99	16.456	2599	2603	2611	rVB	37115	73697	6.62%	0.297%
100	16.706	2640	2644	2649	rVB2	48245	82338	7.40%	0.332%

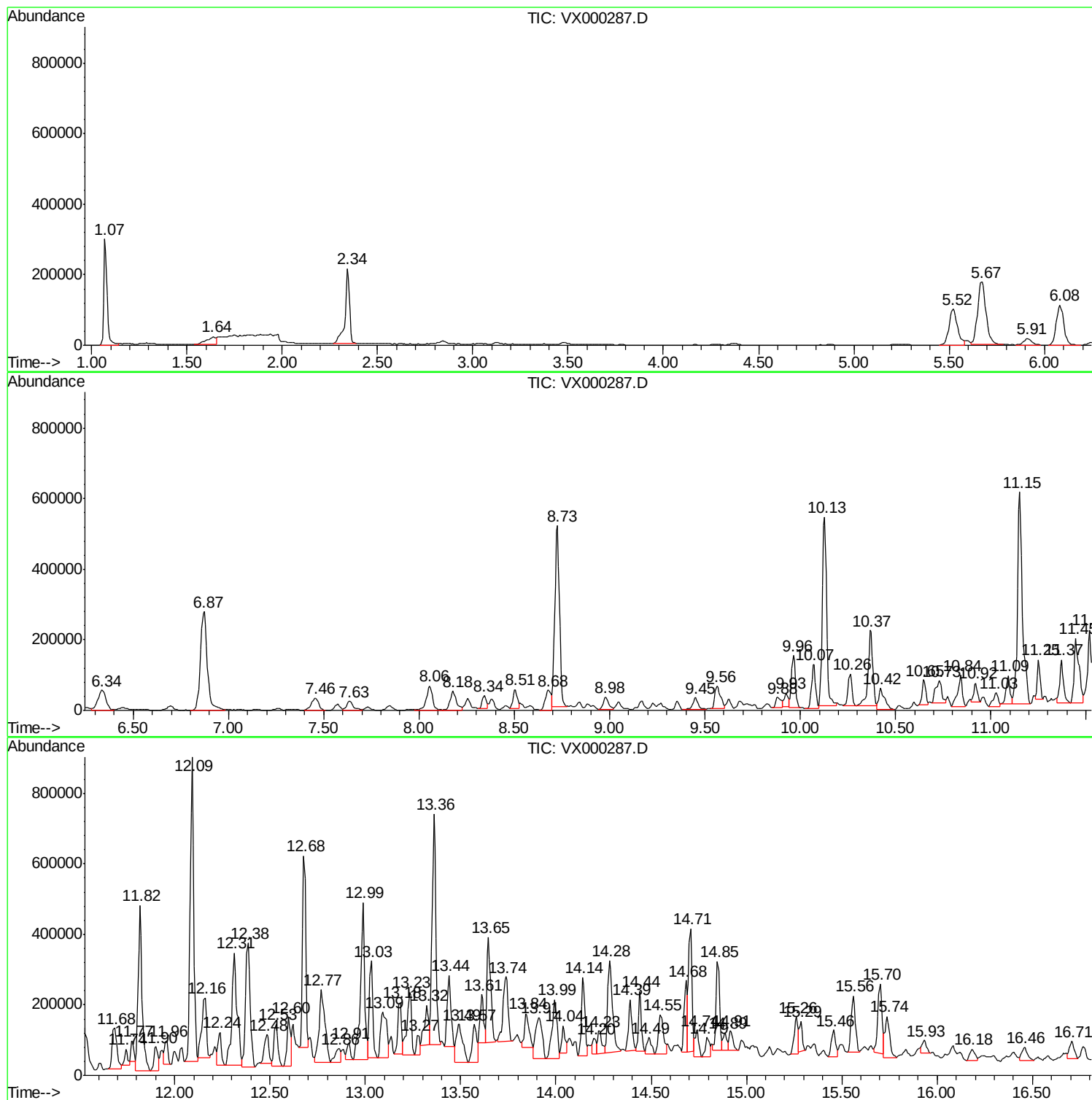
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ALS Vial : 1 Sample Multiplier: 1

Instrument :
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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



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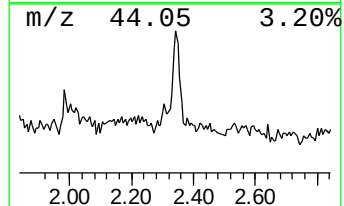
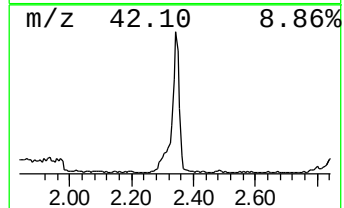
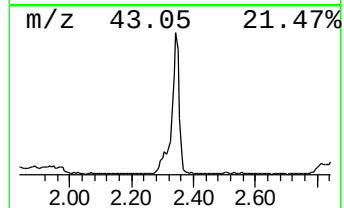
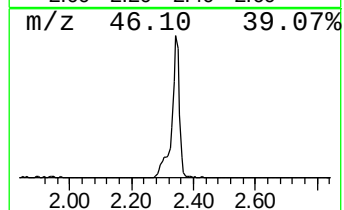
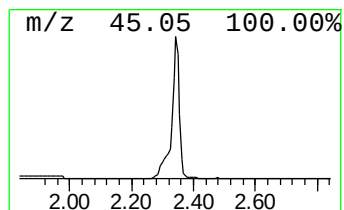
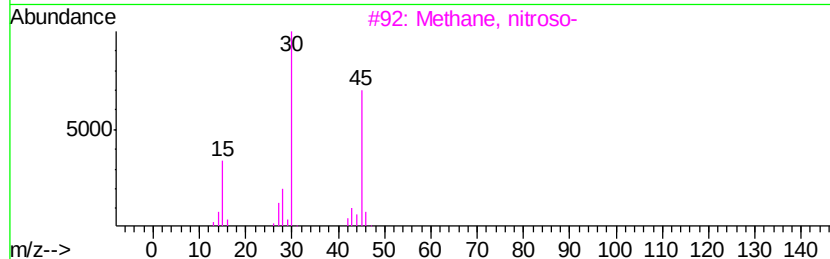
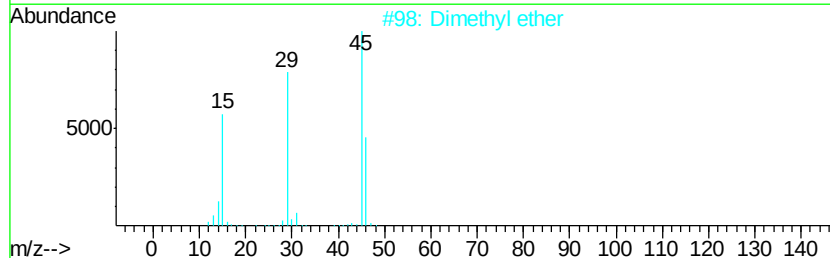
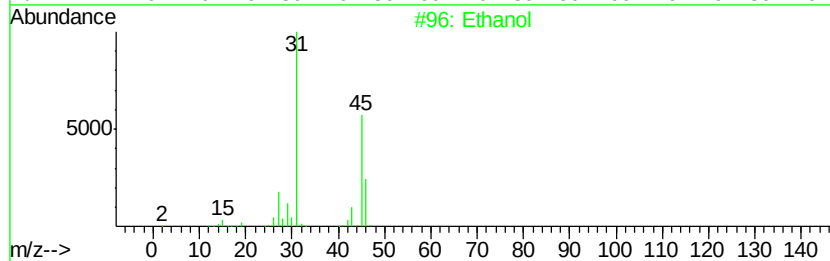
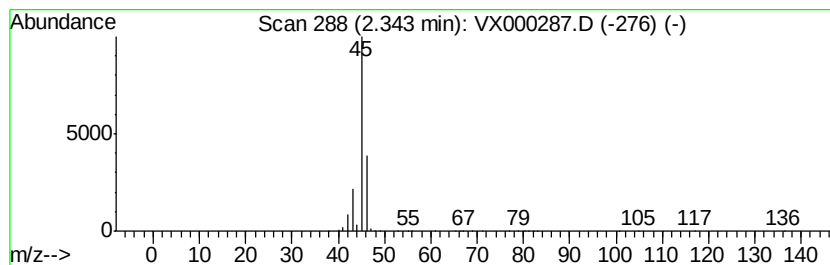
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X030718W.M
Quant Title : SW846 8260

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

Peak Number 2 Ethanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.34	32.49 ug/l	350998	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol	46	C2H6O	000064-17-5	64
2		Dimethyl ether	46	C2H6O	000115-10-6	9
3		Methane, nitroso-	45	CH3NO	000865-40-7	4
4		Formic acid	46	CH2O2	000064-18-6	4
5		Oxalic acid	90	C2H2O4	000144-62-7	4



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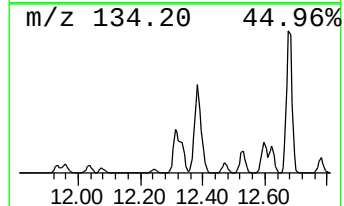
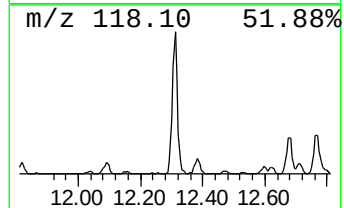
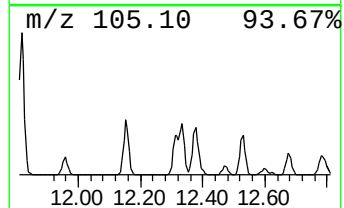
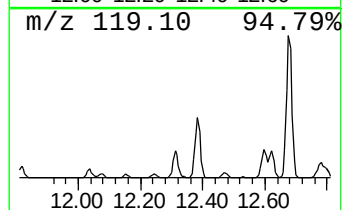
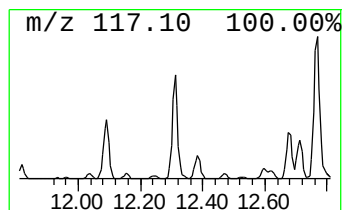
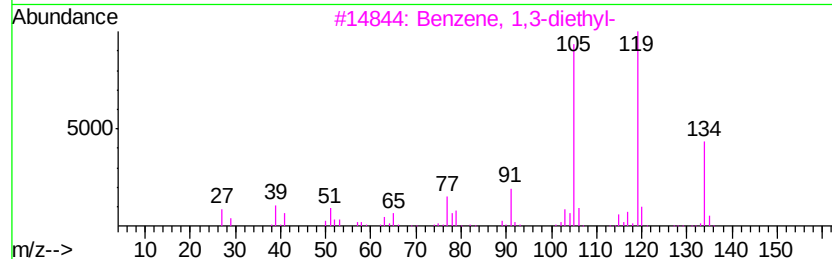
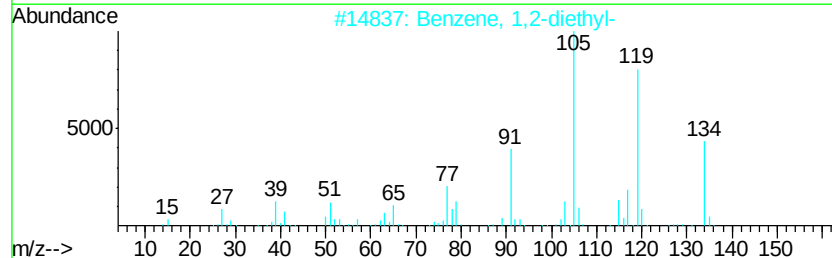
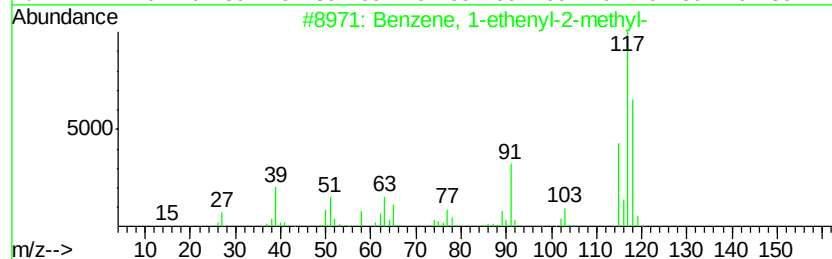
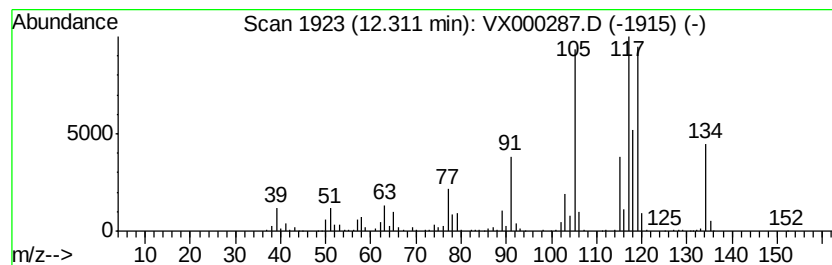
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-ethenyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.31	26.79 ug/l	596303	1,4-Dichlorobenzene-d4	12.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 80
2		Benzene, 1,2-diethyl-	134 C10H14	000135-01-3 60
3		Benzene, 1,3-diethyl-	134 C10H14	000141-93-5 55
4		Benzene, 1,4-diethyl-	134 C10H14	000105-05-5 55
5		Benzene, 2-propenyl-	118 C9H10	000300-57-2 42



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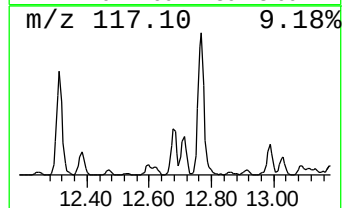
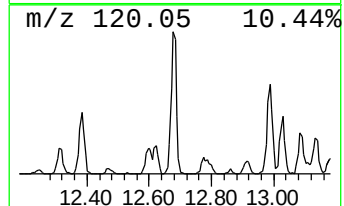
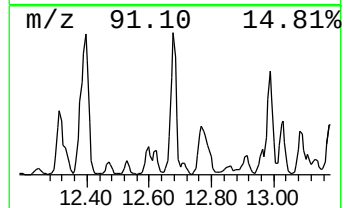
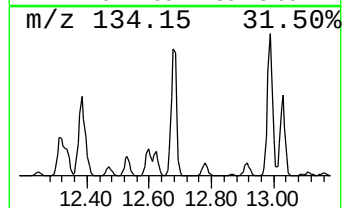
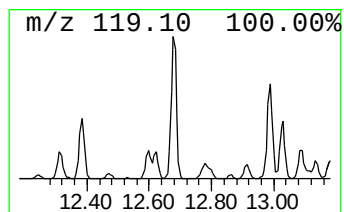
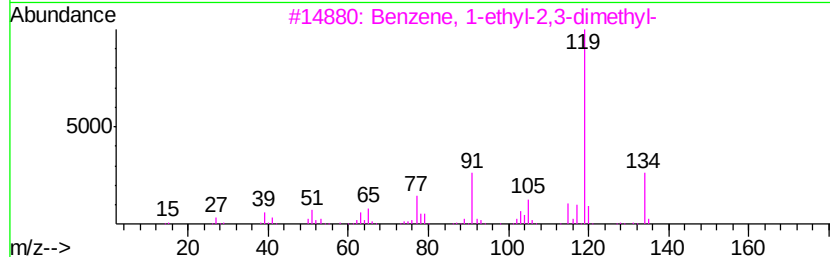
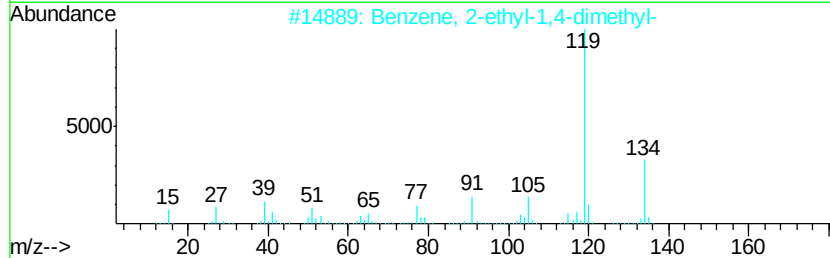
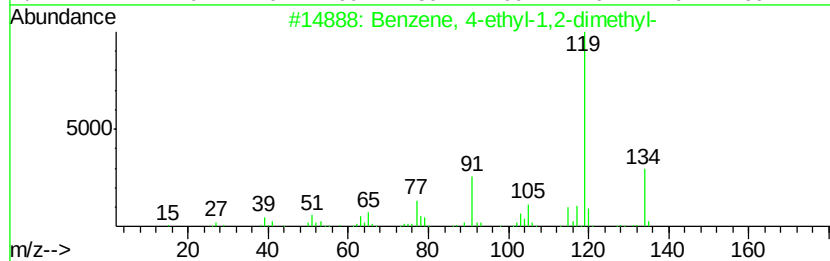
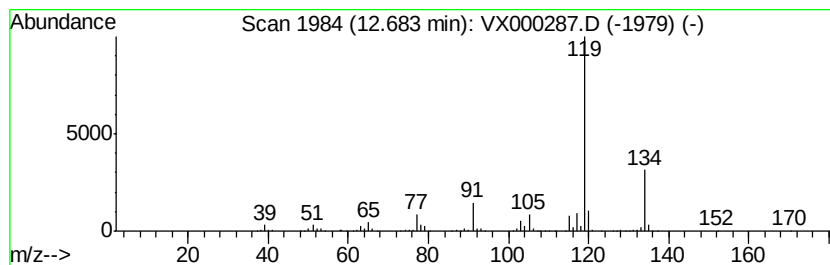
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Peak Number 4 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.68	28.68 ug/l	638346	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
Data File : VX000287.D
Acq On : 12 Mar 2018 20:19
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

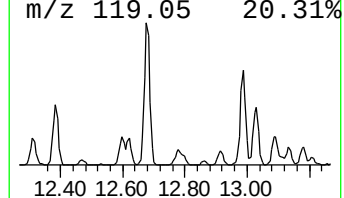
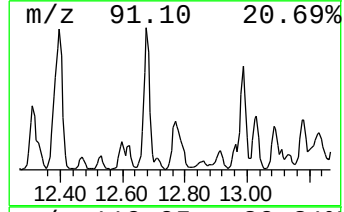
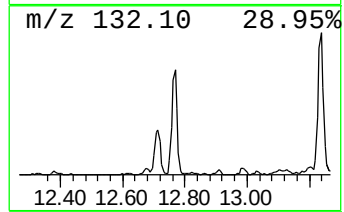
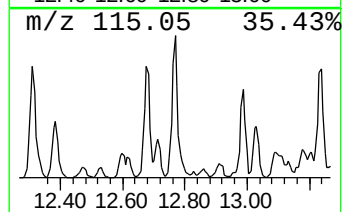
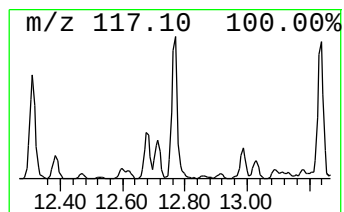
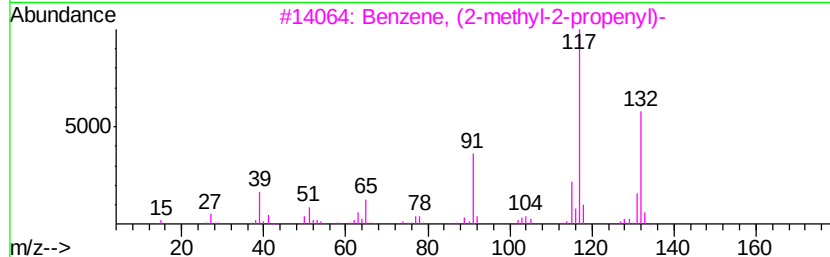
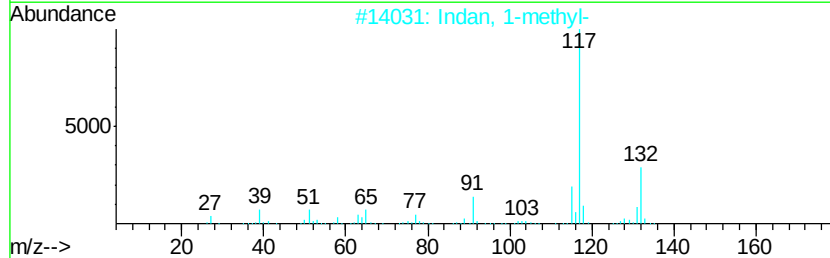
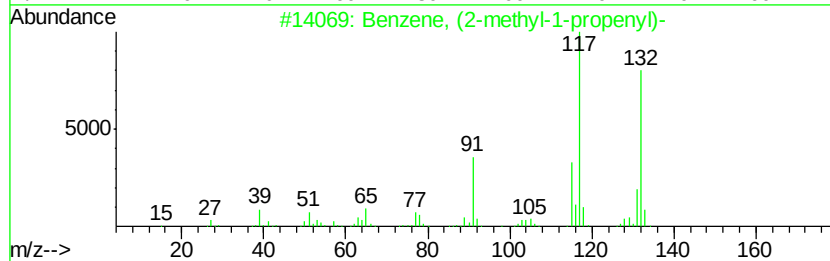
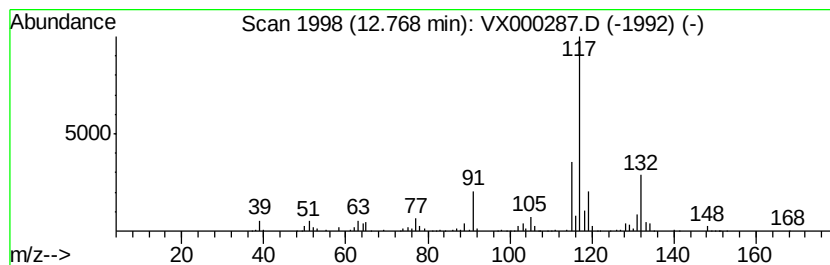
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, (2-methyl-1-propen... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	20.11 ug/l	447711	1,4-Dichlorobenzene-d4	12.09

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	81
2		Indan, 1-methyl-	132	C10H12	000767-58-8	81
3		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	80
4		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	76
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	74



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
Data File : VX000287.D
Acq On : 12 Mar 2018 20:19
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

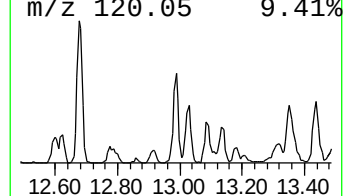
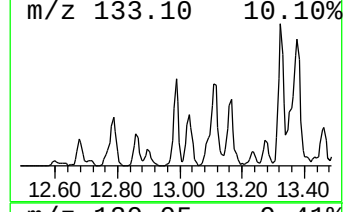
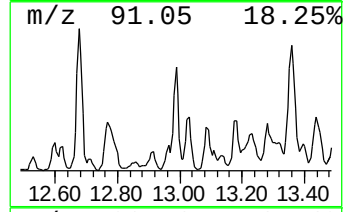
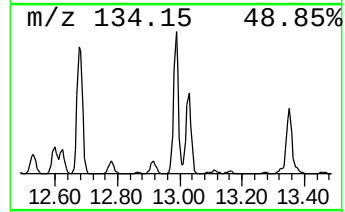
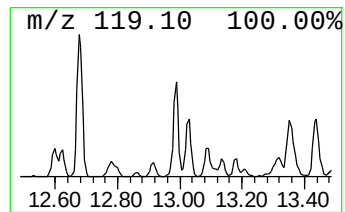
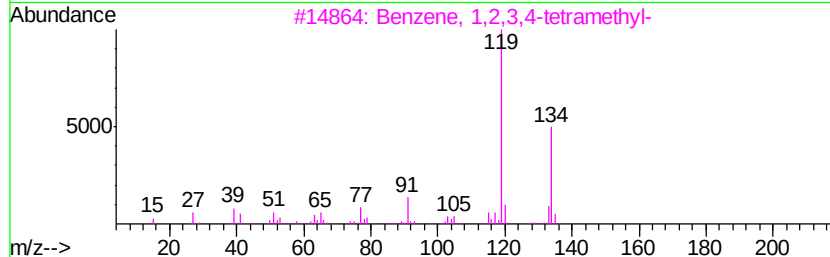
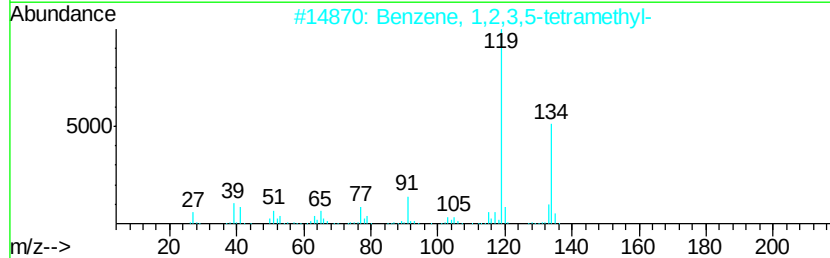
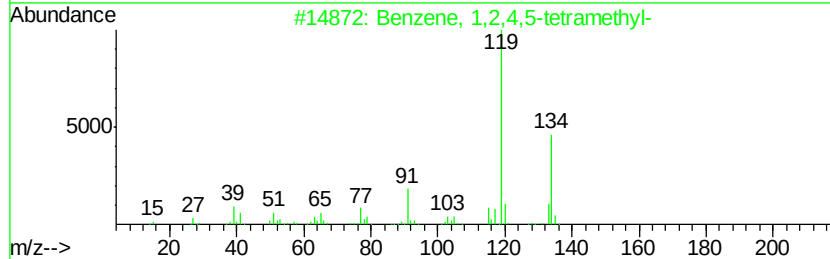
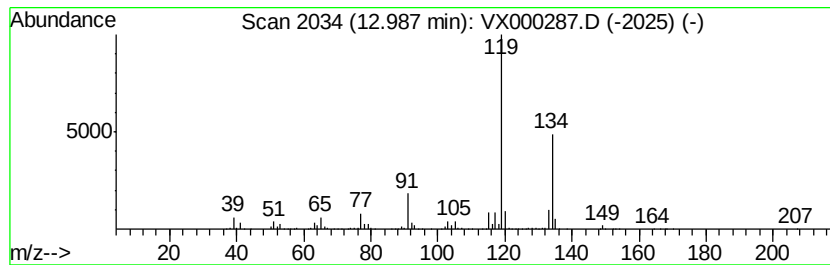
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.99	31.21 ug/l	694761	1,4-Dichlorobenzene-d4	12.09

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	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 : VX000287.D
Acq On : 12 Mar 2018 20:19
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

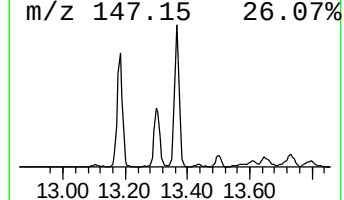
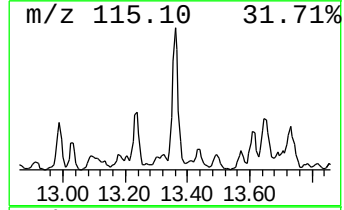
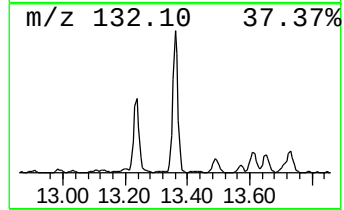
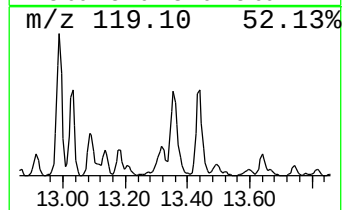
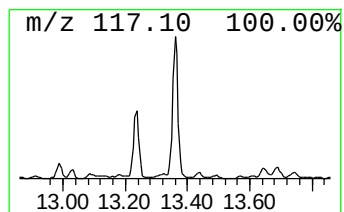
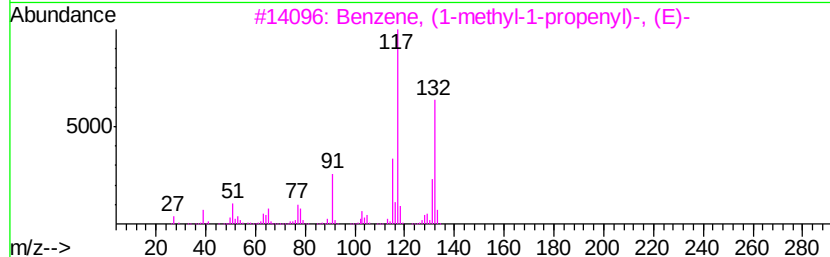
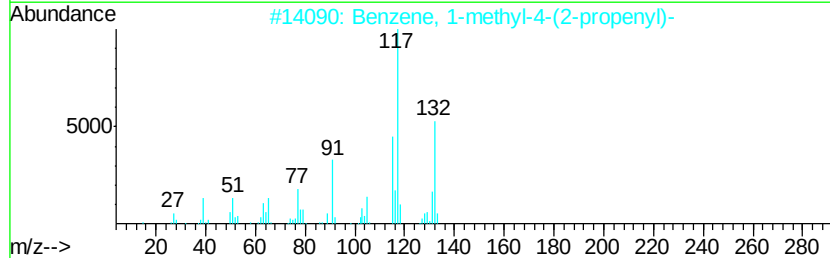
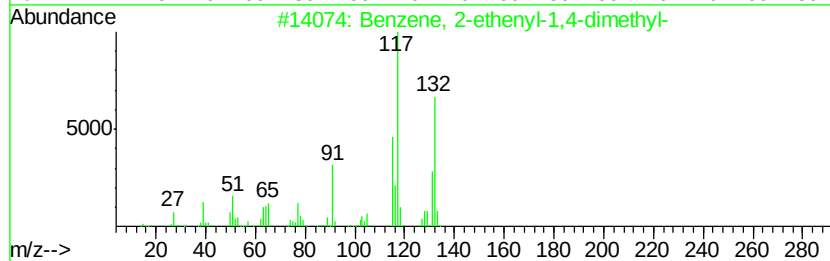
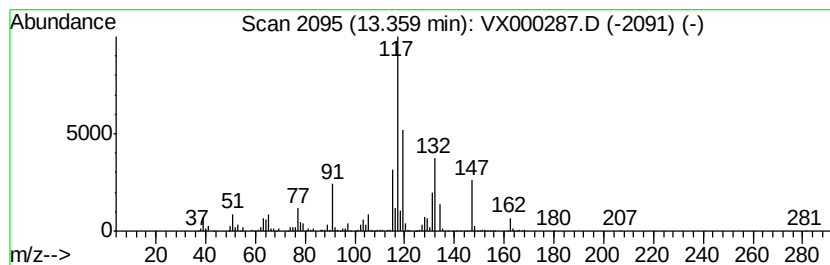
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.36	42.13 ug/l	937891	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
2		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	83
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	64
5		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	60



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
Data File : VX000287.D
Acq On : 12 Mar 2018 20:19
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

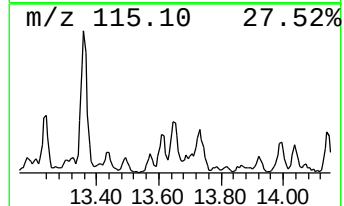
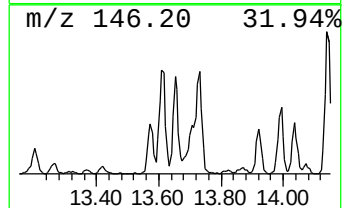
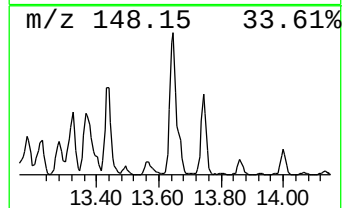
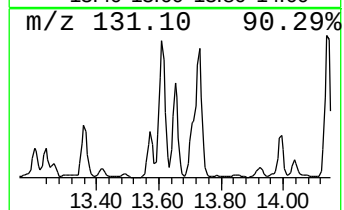
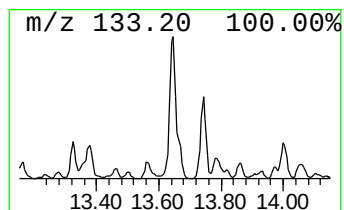
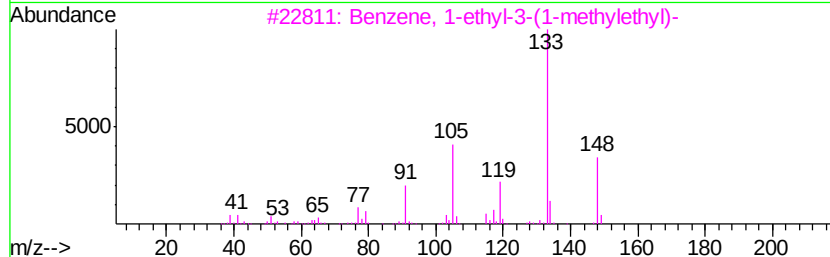
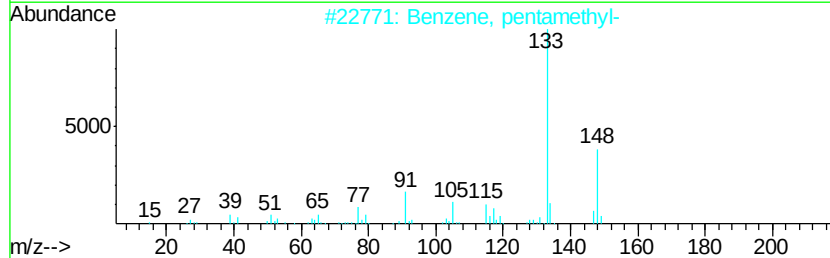
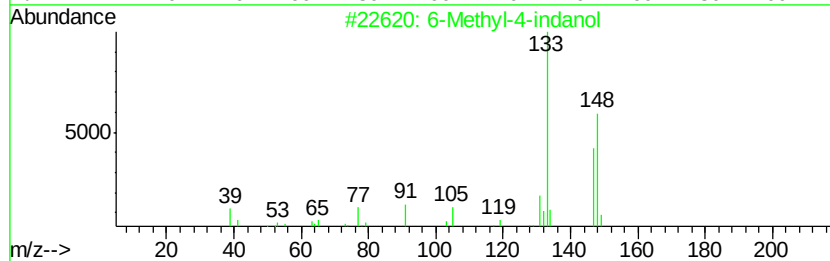
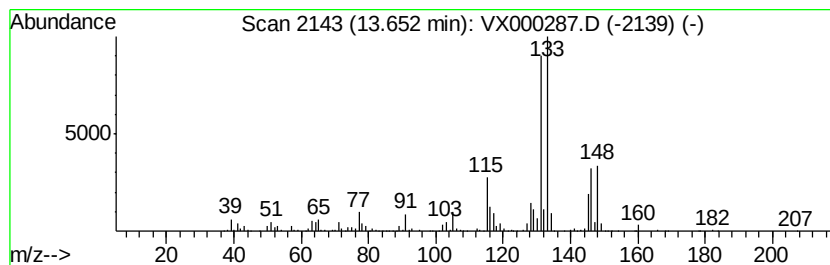
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Quant Title : SW846 8260

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

Peak Number 8 6-Methyl-4-indanol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.65	19.66 ug/l	437570	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	6-Methyl-4-indanol	148	C10H12O	020294-32-0	78
2		Benzene, pentamethyl-	148	C11H16	000700-12-9	76
3		Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	68
4		3,4-Dimethylcumene	148	C11H16	1000370-34-1	64
5		Pyrazine, 2,5-dimethyl-3-(1-prop...	148	C9H12N2	055138-73-3	59



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
Data File : VX000287.D
Acq On : 12 Mar 2018 20:19
Operator : JC/MD
Sample : J1727-04
Misc : 6.64g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 1 Sample Multiplier: 1

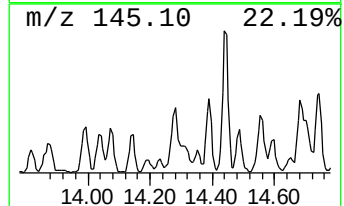
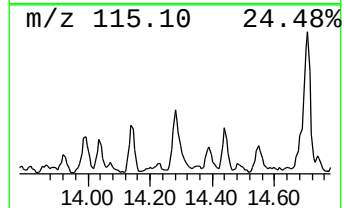
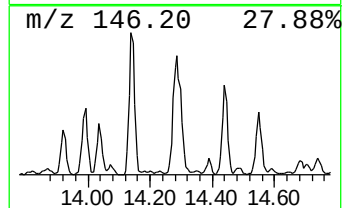
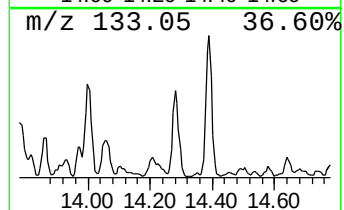
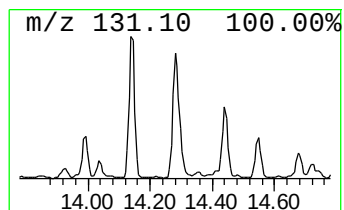
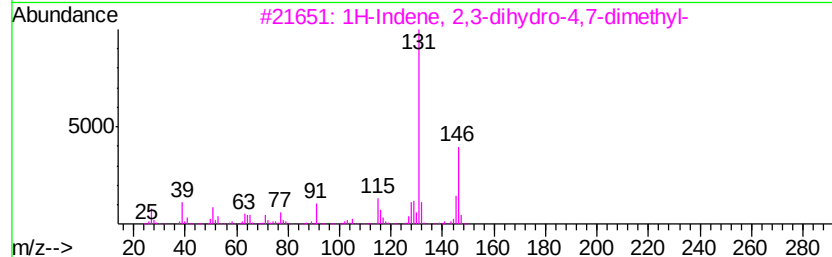
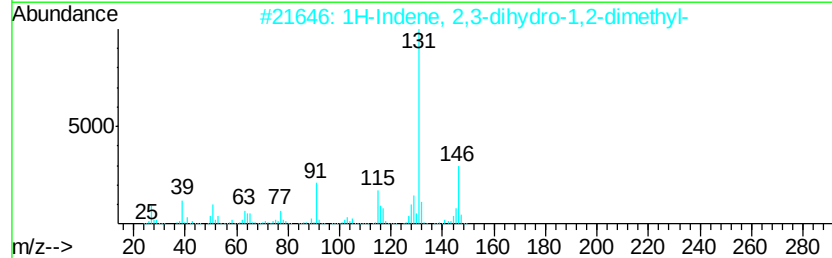
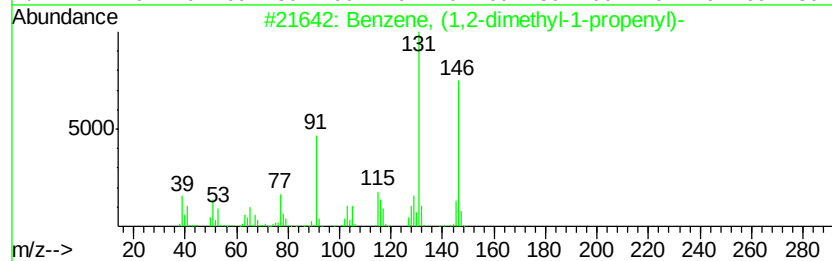
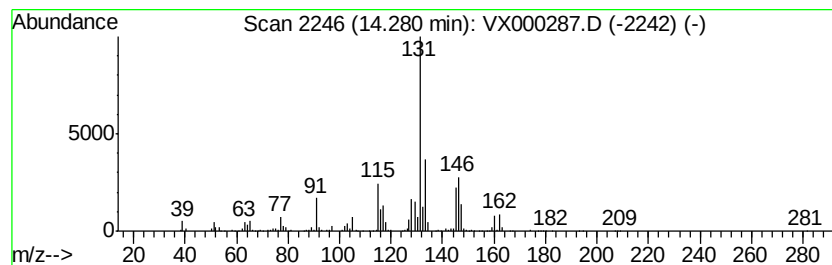
Instrument :
MSVOA_X
ClientSampled :
T-4

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 9 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.28	20.57 ug/l	457996	1,4-Dichlorobenzene-d4	12.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3 93
2		1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8 81
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 76
4		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 76
5		1H-Indene, 2,3-dihydro-4,6-dimet...	146 C11H14	001685-82-1 70



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
Data File : VX000287.D
Acq On : 12 Mar 2018 20:19
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

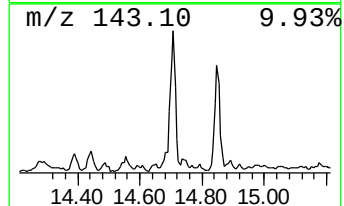
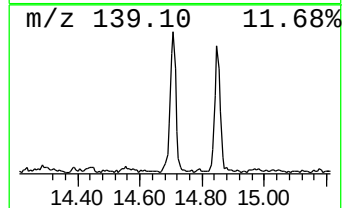
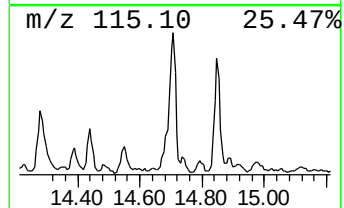
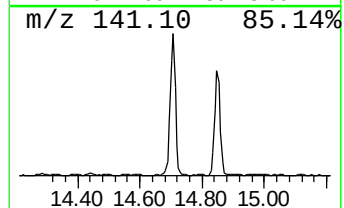
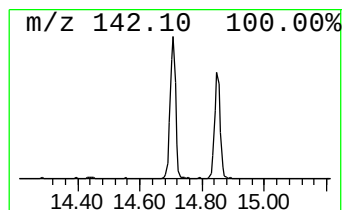
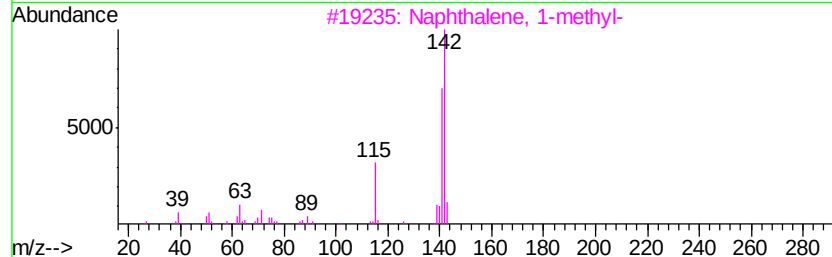
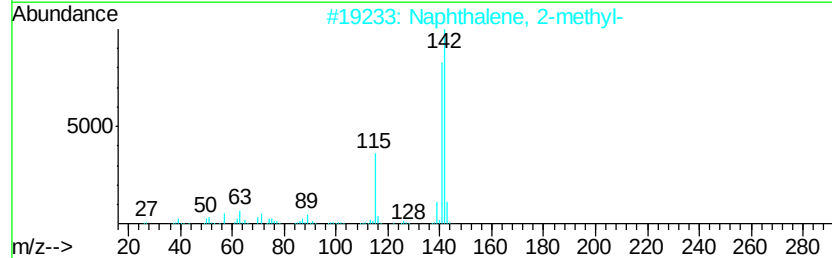
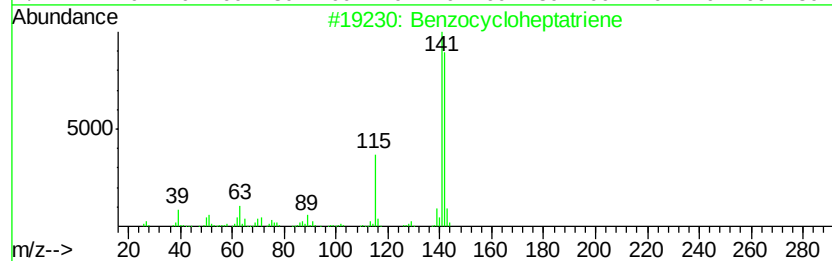
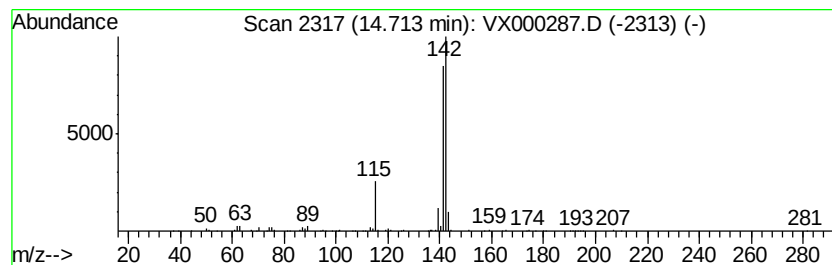
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Quant Title : SW846 8260

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

Peak Number 10 Benzocycloheptatriene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.71	19.74 ug/l	439339	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzocycloheptatriene	142	C11H10	000264-09-5	91
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	91
3		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	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 : VX000287.D
Acq On : 12 Mar 2018 20:19
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
Ethanol	2.34	32.5	ug/l	350998	1	5.67	540106	50.0
Benzene, 1-etheny...	12.31	26.8	ug/l	596303	4	12.09	1113000	50.0
Benzene, 4-ethyl-...	12.68	28.7	ug/l	638346	4	12.09	1113000	50.0
Benzene, (2-methy...	12.77	20.1	ug/l	447711	4	12.09	1113000	50.0
Benzene, 1,2,4,5-...	12.99	31.2	ug/l	694761	4	12.09	1113000	50.0
Benzene, 2-etheny...	13.36	42.1	ug/l	937891	4	12.09	1113000	50.0
6-Methyl-4-indanol	13.65	19.7	ug/l	437570	4	12.09	1113000	50.0
Benzene, (1,2-dim...	14.28	20.6	ug/l	457996	4	12.09	1113000	50.0
Benzocycloheptatr...	14.71	19.7	ug/l	439339	4	12.09	1113000	50.0