

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
 Data File : VX000279.D
 Acq On : 12 Mar 2018 16:54
 Operator : JC/MD
 Sample : J1727-02
 Misc : 5.27g/10mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 T-2

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	1.075	67	80	93	rBV	253960	339710	30.65%	2.068%
2	1.636	156	172	174	rBV5	22700	69382	6.26%	0.422%
3	2.300	274	281	284	rBV	41247	78188	7.05%	0.476%
4	2.380	288	294	305	rVB	93256	224642	20.27%	1.368%
5	5.519	797	809	819	rBV2	63291	189225	17.07%	1.152%
6	5.672	824	834	849	rVB	188011	551551	49.76%	3.358%
7	5.915	862	874	886	rVB7	7374	25461	2.30%	0.155%
8	6.080	890	901	911	rBV2	75581	204769	18.48%	1.247%
9	6.342	936	944	954	rVB2	9317	27456	2.48%	0.167%
10	6.867	1019	1030	1051	rBV	297534	743451	67.08%	4.526%
11	7.458	1115	1127	1136	rVB4	13317	36879	3.33%	0.225%
12	8.055	1218	1225	1234	rVB4	10557	26629	2.40%	0.162%
13	8.177	1237	1245	1253	rBV8	8972	24243	2.19%	0.148%
14	8.342	1267	1272	1276	rVV3	20947	36508	3.29%	0.222%
15	8.378	1276	1278	1284	rVB2	14953	23537	2.12%	0.143%
16	8.500	1293	1298	1303	rBV	29308	48691	4.39%	0.296%
17	8.671	1320	1326	1329	rBV3	19561	37127	3.35%	0.226%
18	8.726	1329	1335	1343	rVB	335988	555149	50.09%	3.380%
19	8.976	1371	1376	1383	rBV3	11922	23102	2.08%	0.141%
20	9.165	1399	1407	1413	rBV3	12631	26347	2.38%	0.160%
21	9.451	1447	1454	1463	rBV4	19094	35252	3.18%	0.215%
22	9.561	1467	1472	1479	rBV2	33898	74472	6.72%	0.453%
23	9.628	1479	1483	1487	rVB4	16976	24913	2.25%	0.152%
24	9.683	1487	1492	1495	rBV4	13619	24235	2.19%	0.148%
25	9.884	1520	1525	1529	rBV2	16043	36882	3.33%	0.225%
26	9.964	1534	1538	1549	rVB	95084	139635	12.60%	0.850%
27	10.067	1549	1555	1559	rBV	84019	121962	11.00%	0.743%
28	10.122	1559	1564	1573	rBV	613738	898329	81.05%	5.469%
29	10.262	1581	1587	1591	rVB3	18596	33749	3.05%	0.205%
30	10.342	1591	1600	1601	rVV5	17278	38447	3.47%	0.234%
31	10.372	1601	1605	1610	rVV2	33455	62668	5.65%	0.382%
32	10.421	1610	1613	1625	rVB6	23857	57765	5.21%	0.352%
33	10.646	1645	1650	1654	rBV	49755	70987	6.40%	0.432%
34	10.732	1658	1664	1669	rBV3	46983	75206	6.79%	0.458%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	10.841	1674	1682	1686	rVB2	64789	100495	9.07%	0.612%
36	10.921	1686	1695	1699	rBV2	45482	86633	7.82%	0.527%
37	11.024	1706	1712	1716	rBV4	20334	36691	3.31%	0.223%
38	11.091	1718	1723	1726	rBV	58651	84485	7.62%	0.514%
39	11.146	1726	1732	1742	rVB2	404661	702427	63.38%	4.276%
40	11.250	1742	1749	1753	rBV	106707	145800	13.15%	0.888%
41	11.372	1764	1769	1776	rVB	60327	83680	7.55%	0.509%
42	11.445	1777	1781	1787	rBV	45153	96976	8.75%	0.590%
43	11.518	1787	1793	1795	rBV2	41000	63382	5.72%	0.386%
44	11.695	1815	1822	1826	rBV3	49718	96100	8.67%	0.585%
45	11.744	1826	1830	1833	rVV2	45676	55386	5.00%	0.337%
46	11.774	1833	1835	1838	rVV	50990	53924	4.87%	0.328%
47	11.817	1838	1842	1851	rVB2	164585	269838	24.35%	1.643%
48	11.902	1851	1856	1858	rBV	51854	75426	6.81%	0.459%
49	11.933	1858	1861	1862	rVV	31482	35514	3.20%	0.216%
50	11.951	1862	1864	1869	rVB2	57245	71093	6.41%	0.433%
51	12.091	1880	1887	1892	rBV	824490	1108324	100.00%	6.747%
52	12.134	1892	1894	1896	rBV	56977	58309	5.26%	0.355%
53	12.158	1896	1898	1902	rVB2	79828	94173	8.50%	0.573%
54	12.231	1908	1910	1916	rVB2	81566	99050	8.94%	0.603%
55	12.311	1916	1923	1930	rBV2	178310	307491	27.74%	1.872%
56	12.378	1930	1934	1941	rVB3	177781	322239	29.07%	1.962%
57	12.481	1946	1951	1955	rBV3	35909	67258	6.07%	0.409%
58	12.530	1955	1959	1965	rVB	45331	67805	6.12%	0.413%
59	12.597	1965	1970	1972	rBV	86920	115973	10.46%	0.706%
60	12.676	1979	1983	1987	rBV	404717	455961	41.14%	2.776%
61	12.768	1992	1998	2006	rVB4	111133	280110	25.27%	1.705%
62	12.914	2019	2022	2025	rVB3	19063	23596	2.13%	0.144%
63	12.987	2025	2034	2038	rBV	327205	526514	47.51%	3.205%
64	13.030	2038	2041	2046	rVB	154545	192997	17.41%	1.175%
65	13.091	2046	2051	2056	rBV2	127353	235341	21.23%	1.433%
66	13.176	2060	2065	2068	rBV2	114049	157144	14.18%	0.957%
67	13.231	2071	2074	2078	rVB2	122395	156963	14.16%	0.956%
68	13.280	2078	2082	2084	rBV3	40620	54110	4.88%	0.329%
69	13.323	2084	2089	2091	rVV4	104249	157457	14.21%	0.959%
70	13.359	2091	2095	2103	rVV4	397609	607966	54.85%	3.701%
71	13.432	2103	2107	2113	rVV	179688	265549	23.96%	1.617%

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72	13.493	2113	2117	2124	rVB4	58953	139269	12.57%	0.848%
73	13.573	2124	2130	2133	rBV2	95575	153446	13.84%	0.934%
74	13.609	2133	2136	2139	rVV	167661	223787	20.19%	1.362%
75	13.646	2139	2142	2148	rVV3	296119	446150	40.25%	2.716%
76	13.737	2150	2157	2163	rVB3	185556	372384	33.60%	2.267%
77	13.896	2180	2183	2185	rBV2	29145	33864	3.06%	0.206%
78	13.993	2192	2199	2203	rBV3	128789	239464	21.61%	1.458%
79	14.036	2203	2206	2209	rVV	65590	82592	7.45%	0.503%
80	14.066	2209	2211	2214	rVB2	24168	28981	2.61%	0.176%
81	14.140	2219	2223	2227	rVV	185074	233712	21.09%	1.423%
82	14.201	2231	2233	2234	rVV2	22936	24776	2.24%	0.151%
83	14.225	2235	2237	2241	rVV	49159	61280	5.53%	0.373%
84	14.280	2241	2246	2256	rVV	211429	354239	31.96%	2.157%
85	14.390	2260	2264	2269	rVV	126830	173318	15.64%	1.055%
86	14.438	2269	2272	2276	rVB	140065	165234	14.91%	1.006%
87	14.487	2276	2280	2285	rVB5	28642	47333	4.27%	0.288%
88	14.554	2285	2291	2295	rBV3	47795	89901	8.11%	0.547%
89	14.707	2308	2316	2320	rBV	258994	383424	34.59%	2.334%
90	14.743	2320	2322	2326	rVB2	26305	28036	2.53%	0.171%
91	14.792	2326	2330	2334	rBV5	25965	42559	3.84%	0.259%
92	14.847	2335	2339	2343	rVV	190482	240696	21.72%	1.465%
93	14.877	2343	2344	2348	rVV2	19922	22338	2.02%	0.136%
94	15.261	2402	2407	2414	rBV2	26825	50251	4.53%	0.306%
95	15.456	2434	2439	2442	rBV3	29669	47447	4.28%	0.289%
96	15.560	2450	2456	2472	rVB	64811	125301	11.31%	0.763%
97	15.700	2472	2479	2482	rBV	61229	102153	9.22%	0.622%
98	15.737	2482	2485	2494	rVB	41219	62273	5.62%	0.379%
99	15.926	2513	2516	2522	rVB3	13913	25935	2.34%	0.158%
100	16.078	2538	2541	2553	rVB3	13417	26878	2.43%	0.164%

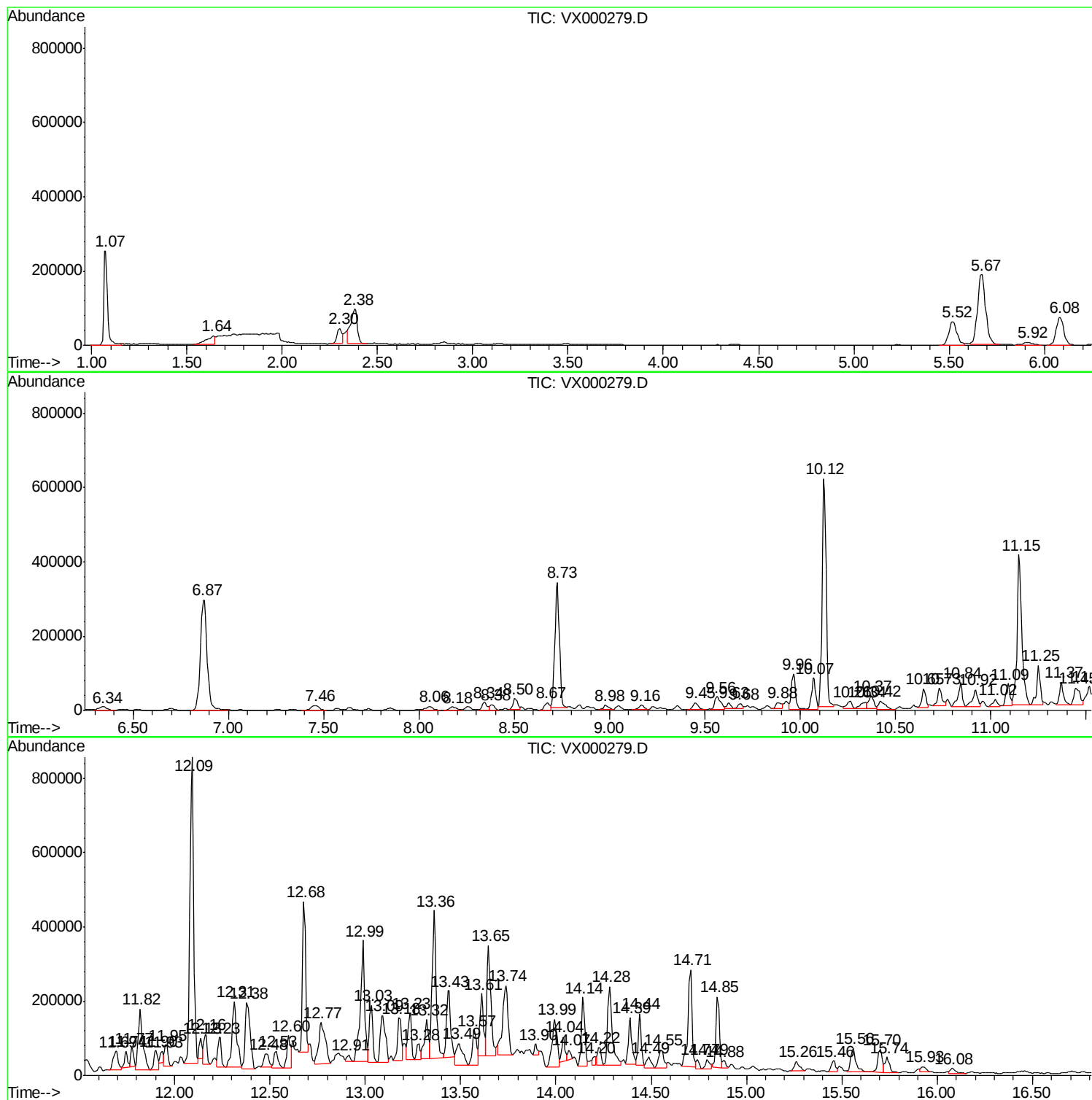
Sum of corrected areas: 16425750

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T-2

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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ClientSampled :
T-2

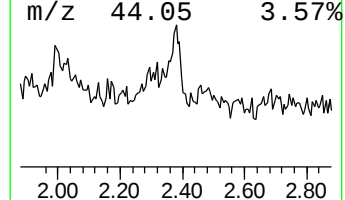
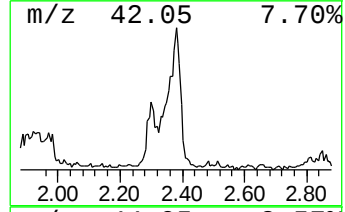
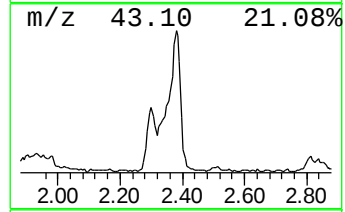
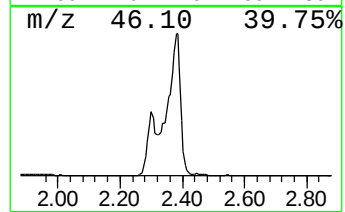
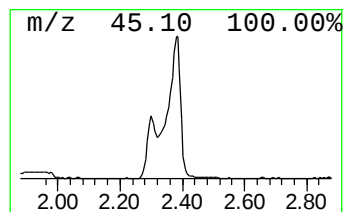
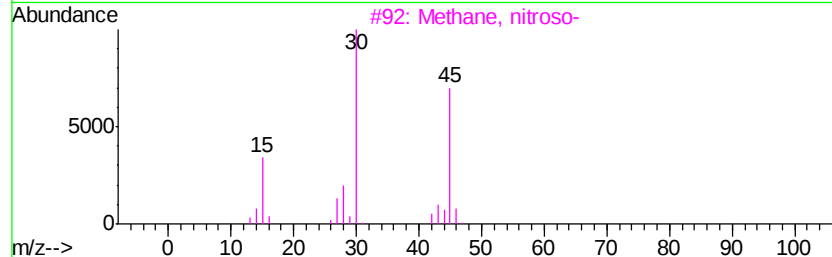
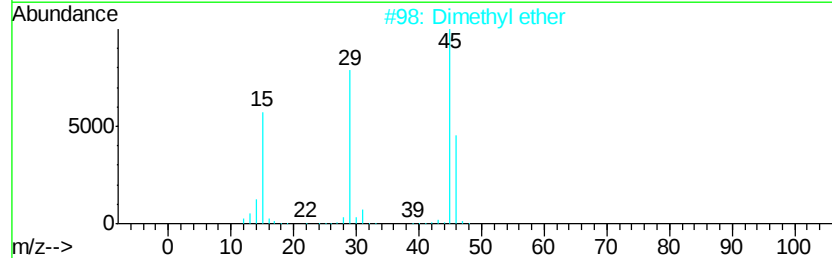
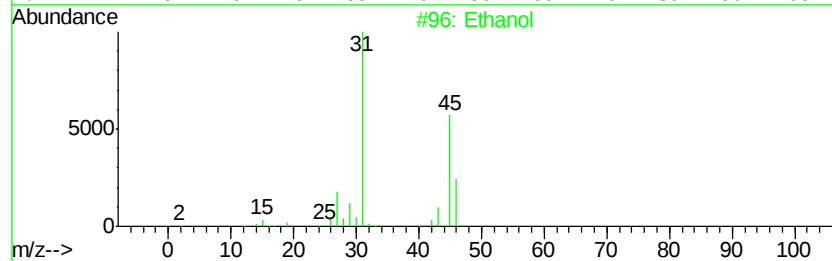
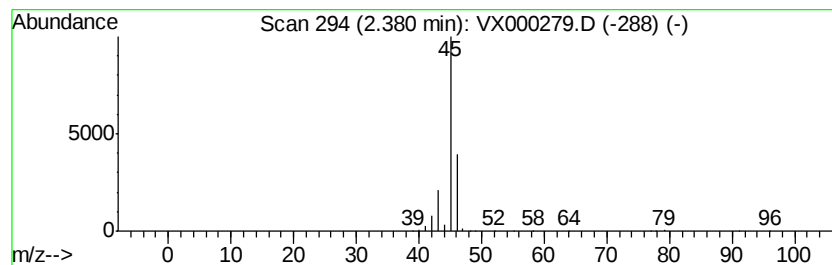
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X030718W.M
Quant Title : SW846 8260

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

Peak Number 2 Ethanol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.38	20.36 ug/l	224642	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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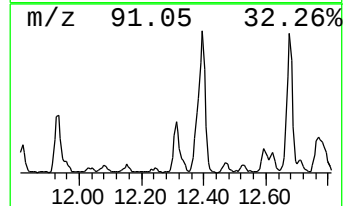
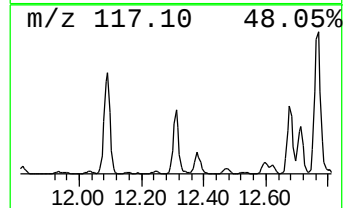
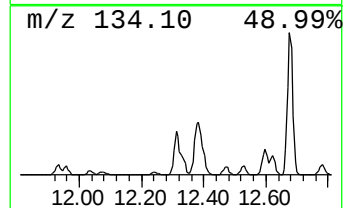
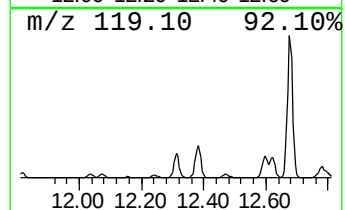
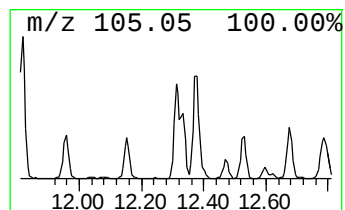
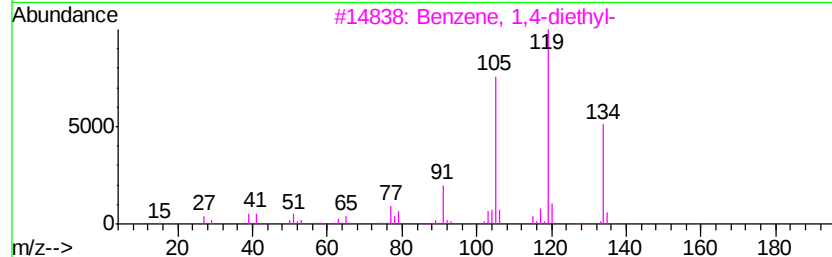
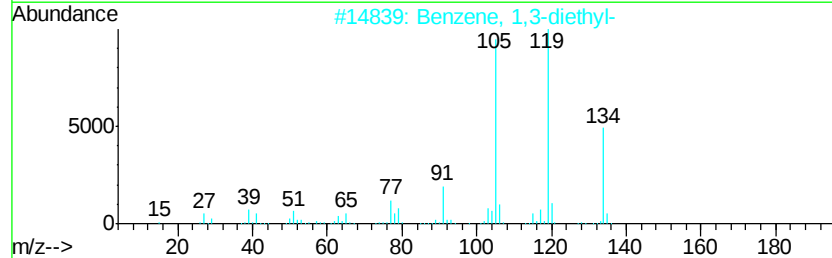
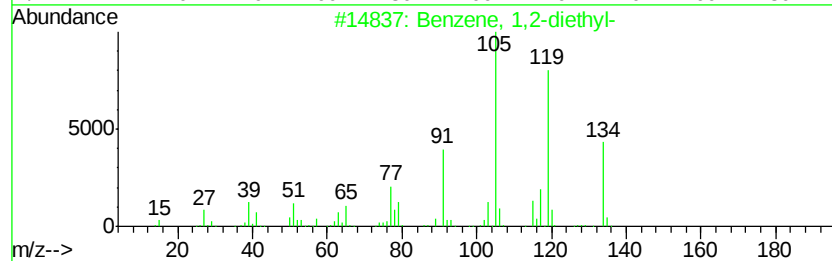
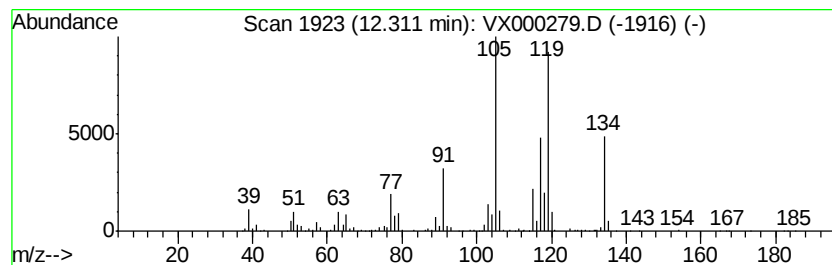
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1,2-diethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.31	13.87 ug/l	307491	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	89
2		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	76
3		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	76
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	53
5		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	49



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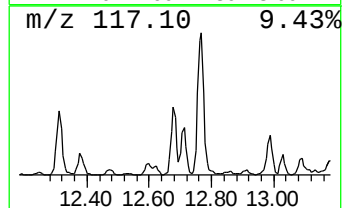
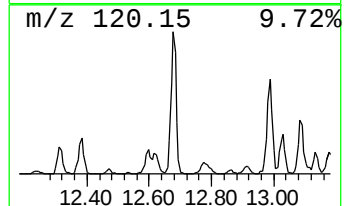
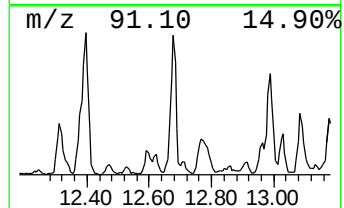
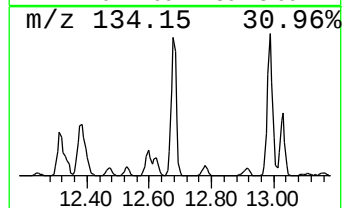
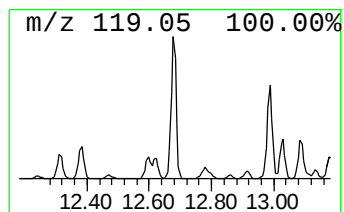
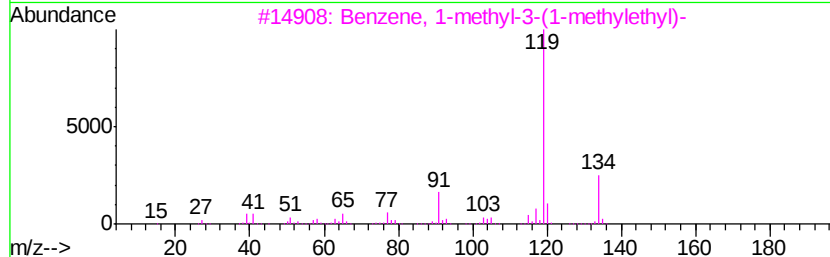
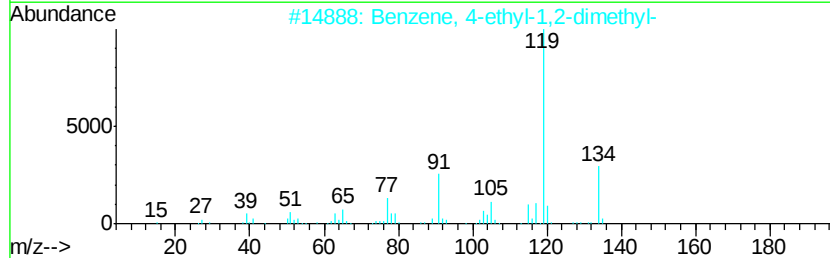
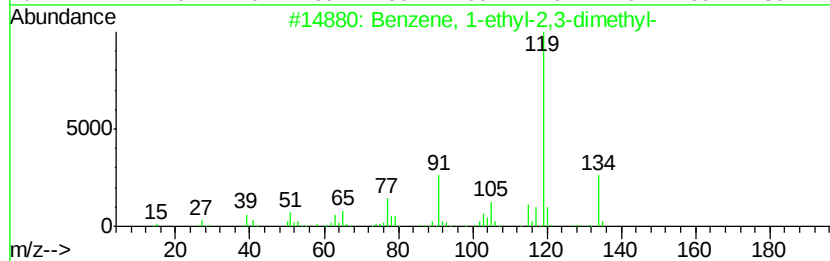
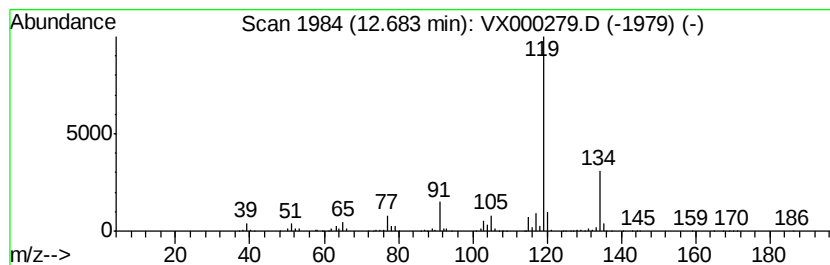
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X030718W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.68	20.57 ug/l	455961	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	94
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
Data File : VX000279.D
Acq On : 12 Mar 2018 16:54
Operator : JC/MD
Sample : J1727-02
Misc : 5.27g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T-2

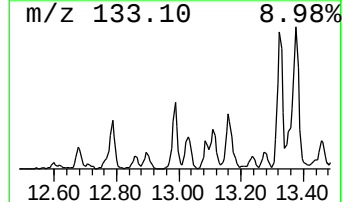
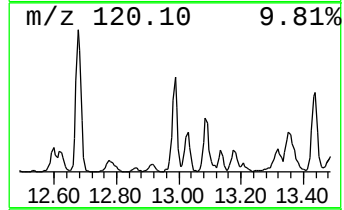
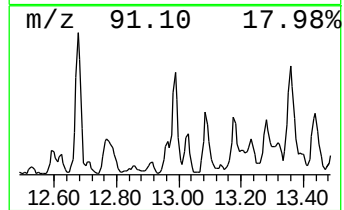
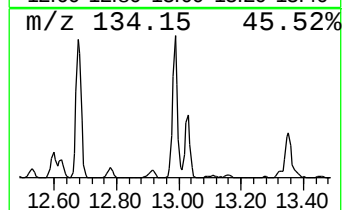
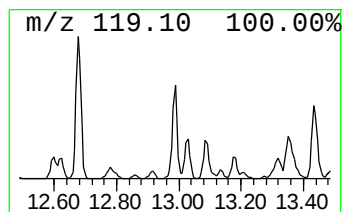
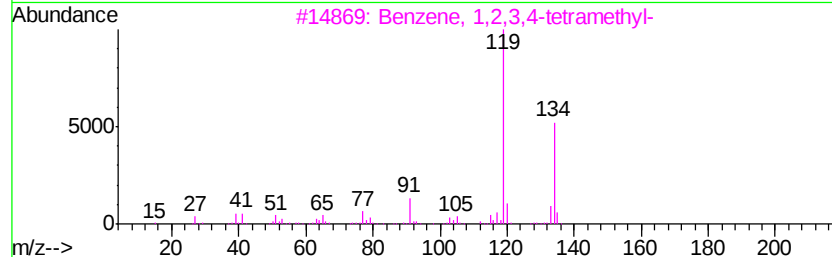
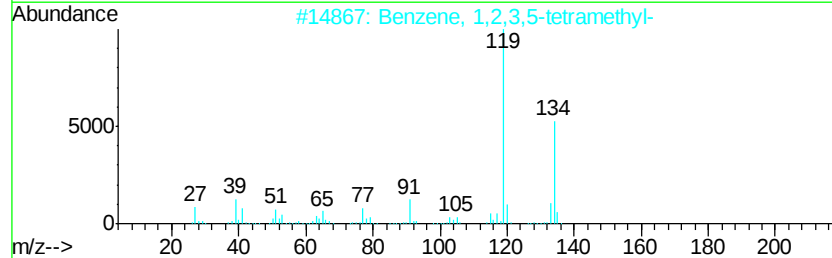
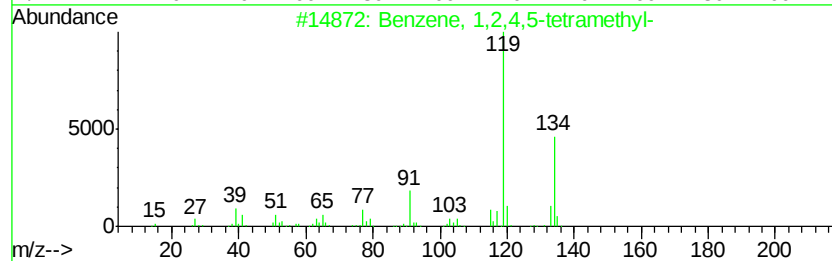
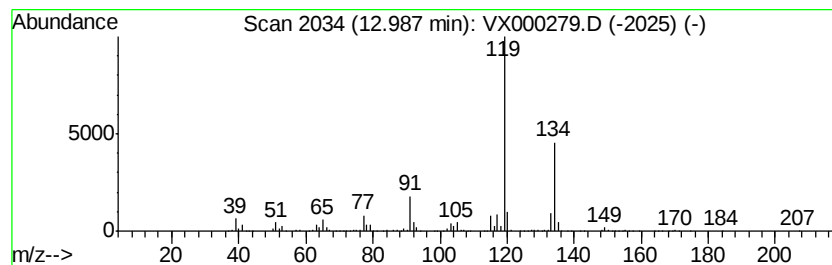
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.99	23.75 ug/l	526514	1,4-Dichlorobenzene-d4	12.09

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
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-2,4-dimethyl-	134	C10H14	000874-41-9	93
5	o-Cymene	134	C10H14	000527-84-4	91



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
Data File : VX000279.D
Acq On : 12 Mar 2018 16:54
Operator : JC/MD
Sample : J1727-02
Misc : 5.27g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T-2

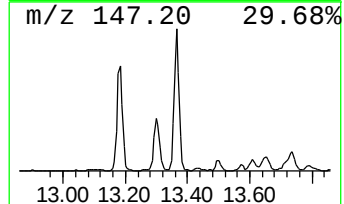
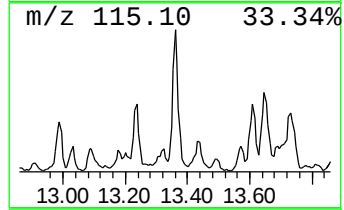
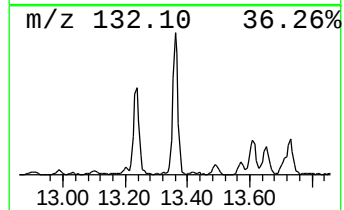
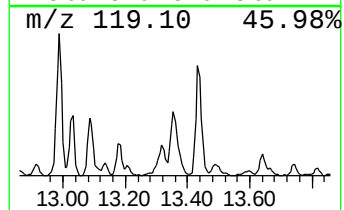
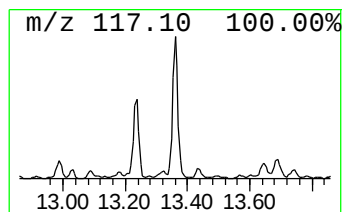
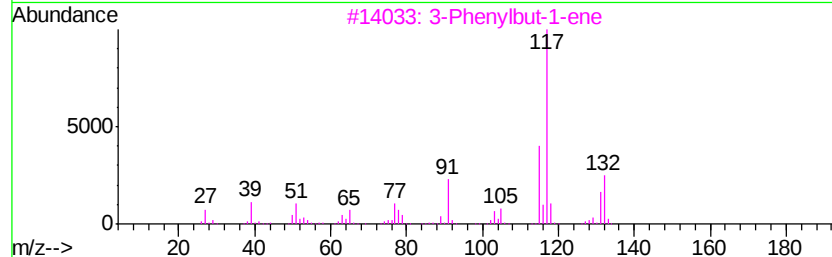
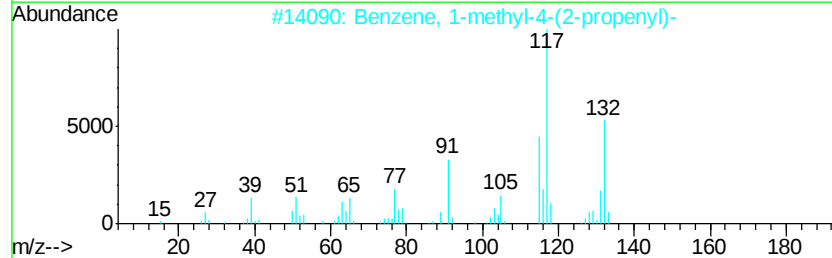
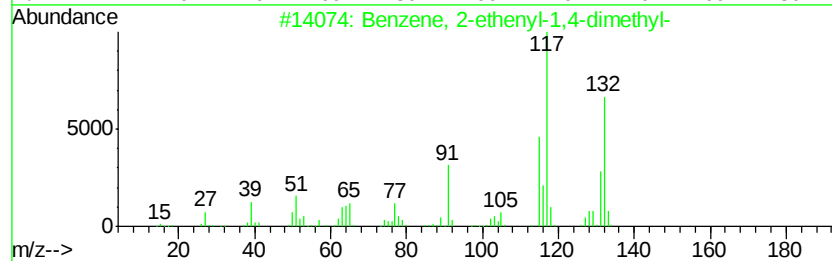
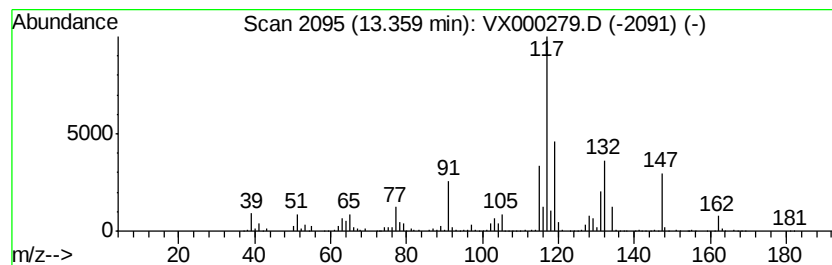
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.36	27.43 ug/l	607966	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	89
2		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	83
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	64
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	60
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	60



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
Data File : VX000279.D
Acq On : 12 Mar 2018 16:54
Operator : JC/MD
Sample : J1727-02
Misc : 5.27g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T-2

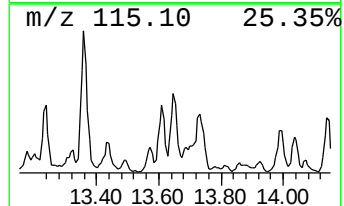
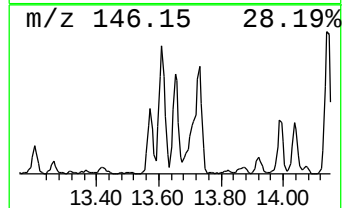
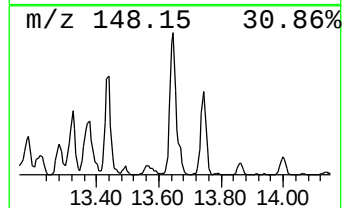
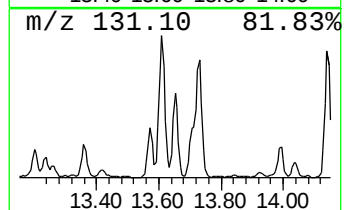
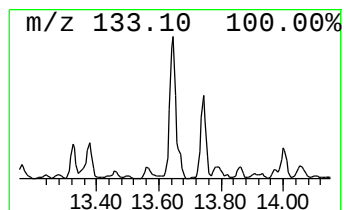
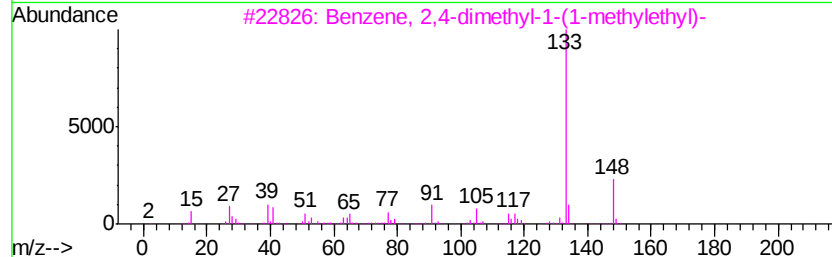
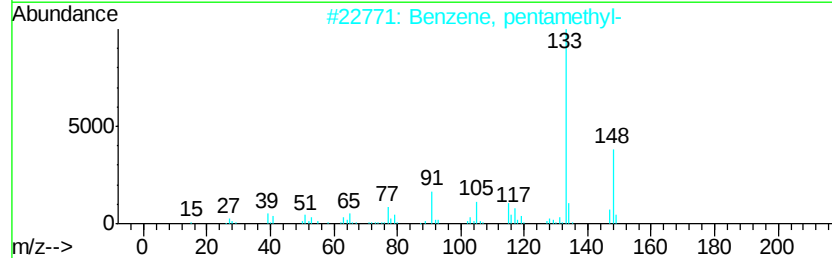
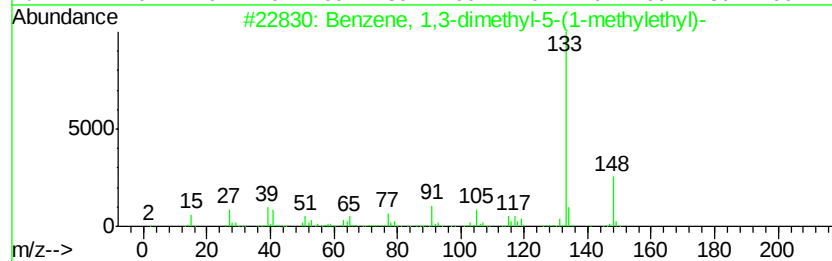
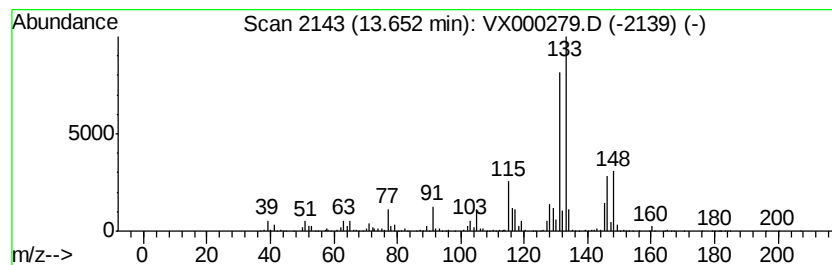
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,3-dimethyl-5-(1-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.65	20.13 ug/l	446150	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	76
2		Benzene, pentamethyl-	148	C11H16	000700-12-9	76
3		Benzene, 2,4-dimethyl-1-(1-methv...	148	C11H16	004706-89-2	70
4		3,4-Dimethylcumene	148	C11H16	1000370-34-1	68
5		Benzene, 1,4-dimethyl-2-(1-methy...	148	C11H16	004132-72-3	62



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
Data File : VX000279.D
Acq On : 12 Mar 2018 16:54
Operator : JC/MD
Sample : J1727-02
Misc : 5.27g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
T-2

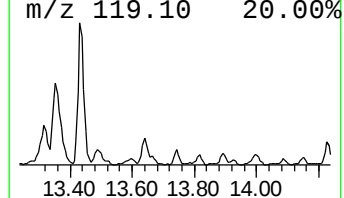
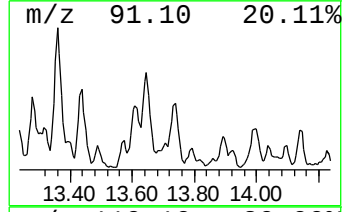
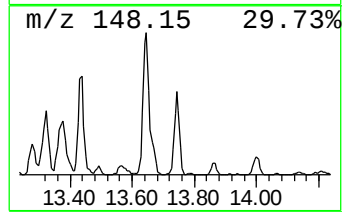
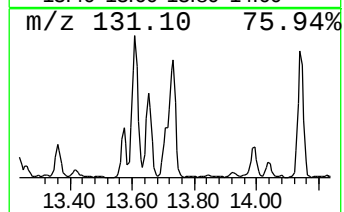
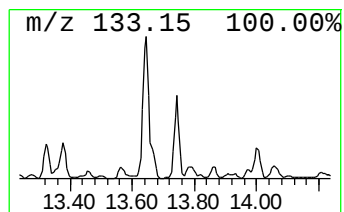
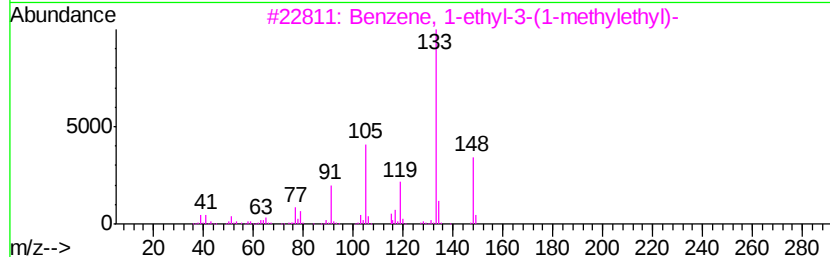
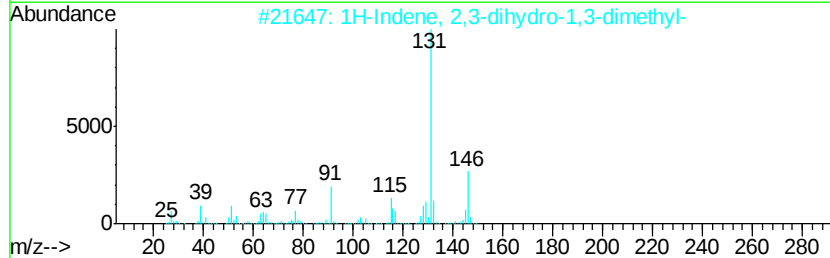
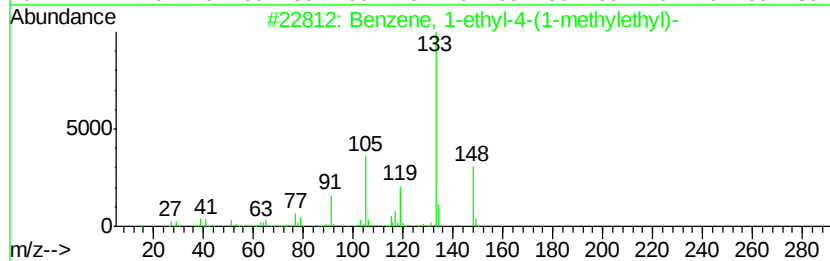
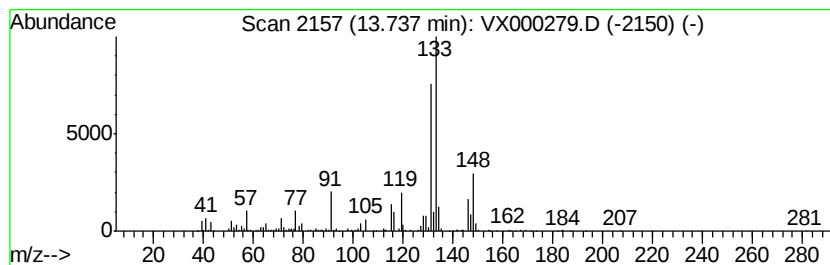
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Quant Title : SW846 8260

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

Peak Number 8 unknown13.74 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.74	16.80 ug/l	372384	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	49
2		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	46
3		Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	46
4		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	46
5		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	41



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
Data File : VX000279.D
Acq On : 12 Mar 2018 16:54
Operator : JC/MD
Sample : J1727-02
Misc : 5.27g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 1 Sample Multiplier: 1

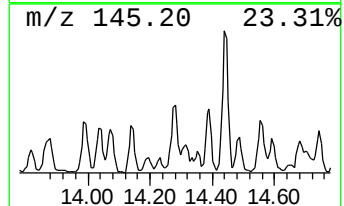
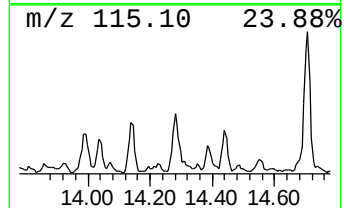
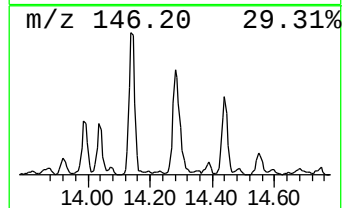
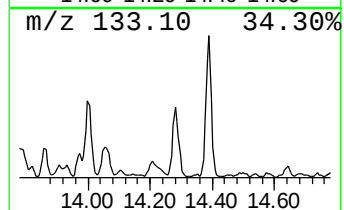
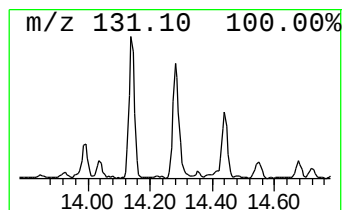
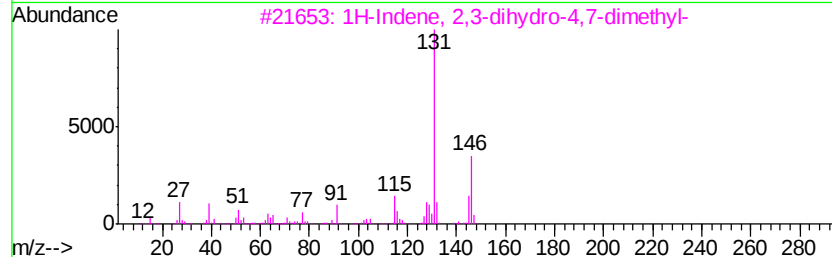
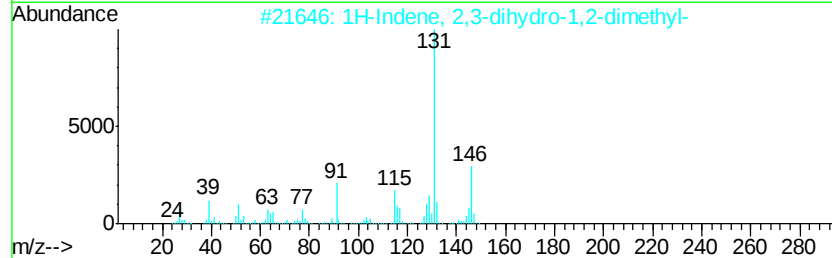
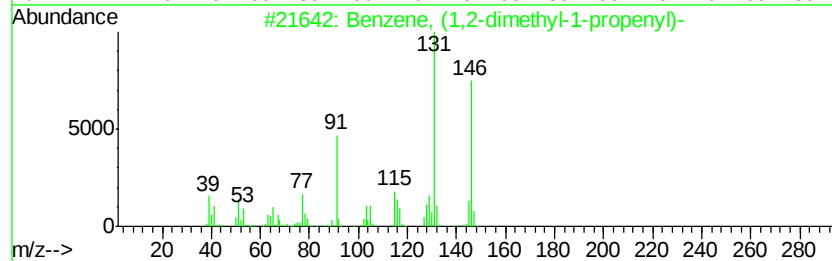
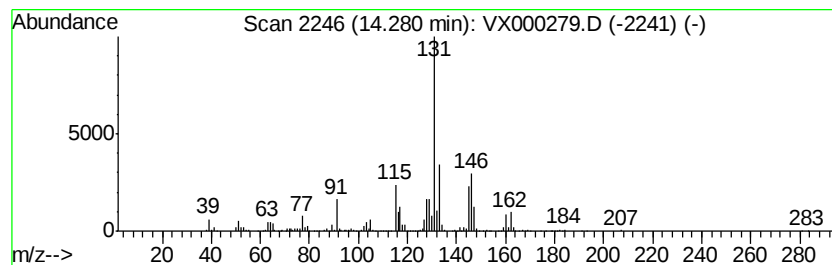
Instrument :
MSVOA_X
ClientSampleId :
T-2

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 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.28	15.98 ug/l	354239	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 92
2		1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8 76
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 70
4		Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2 70
5		1H-Indene, 2,3-dihydro-5,6-dimet...	146 C11H14	001075-22-5 64



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX031218\
Data File : VX000279.D
Acq On : 12 Mar 2018 16:54
Operator : JC/MD
Sample : J1727-02
Misc : 5.27g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
T-2

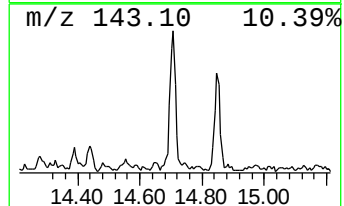
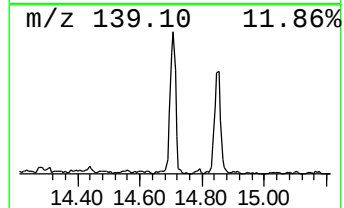
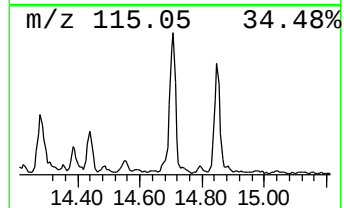
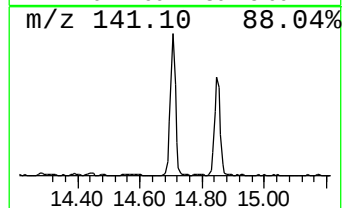
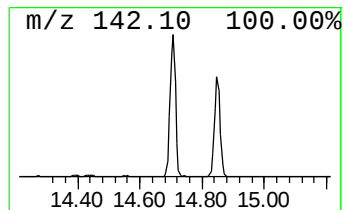
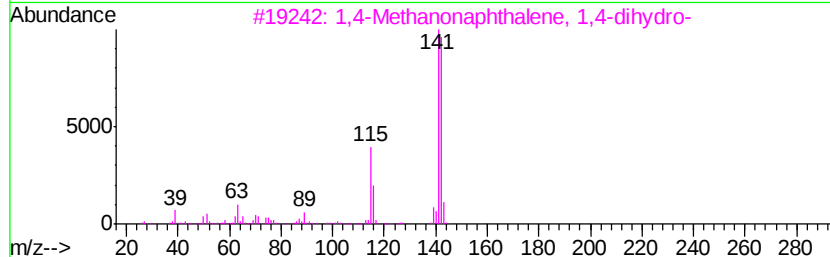
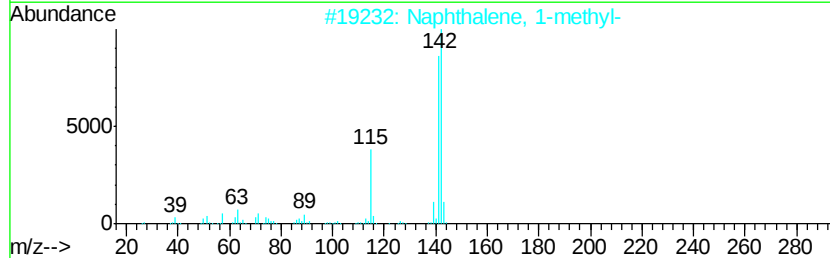
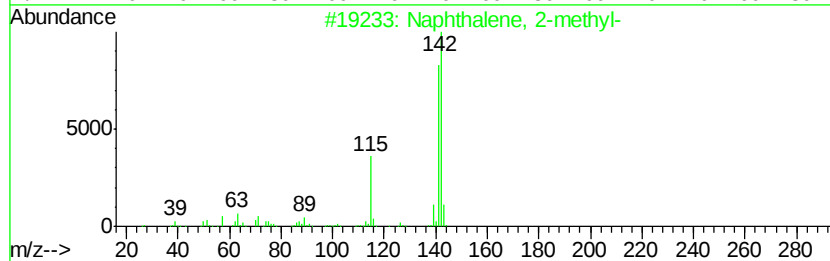
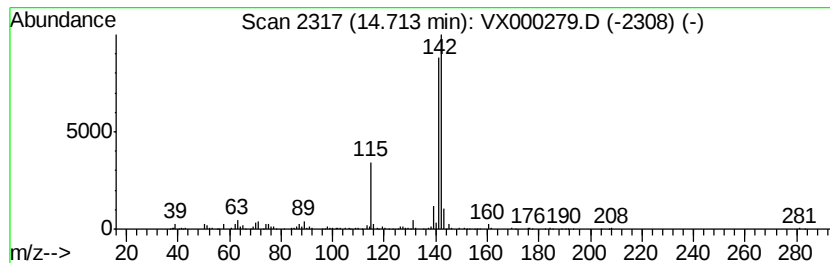
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.71	17.30 ug/l	383424	1,4-Dichlorobenzene-d4	12.09

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		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	91
4		Benzocycloheptatriene	142	C11H10	000264-09-5	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	87



Data Path : W:\HPCHEM1\MSVOA_X\DATA\VX031218\
Data File : VX000279.D
Acq On : 12 Mar 2018 16:54
Operator : JC/MD
Sample : J1727-02
Misc : 5.27g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T-2

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	2.38	20.4	ug/l	224642	1	5.67	551551	50.0
Benzene, 1,2-diet...	12.31	13.9	ug/l	307491	4	12.09	1108320	50.0
Benzene, 1-ethyl-...	12.68	20.6	ug/l	455961	4	12.09	1108320	50.0
Benzene, 1,2,4,5-...	12.99	23.8	ug/l	526514	4	12.09	1108320	50.0
Benzene, 2-etheny...	13.36	27.4	ug/l	607966	4	12.09	1108320	50.0
Benzene, 1,3-dime...	13.65	20.1	ug/l	446150	4	12.09	1108320	50.0
unknown13.74	13.74	16.8	ug/l	372384	4	12.09	1108320	50.0
Benzene, (1,2-dim...	14.28	16.0	ug/l	354239	4	12.09	1108320	50.0
Naphthalene, 1-me...	14.71	17.3	ug/l	383424	4	12.09	1108320	50.0