

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061318\
 Data File : VX002529.D
 Acq On : 14 Jun 2018 01:21
 Operator : JC/MD
 Sample : J3413-04
 Misc : 6.70g/5mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 B-5ARR

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\82X060418W.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	92	rBV	792424	998829	6.80%	0.578%
2	1.941	217	222	232	rVB	227738	338199	2.30%	0.196%
3	2.855	366	372	384	rVB	479807	1043950	7.11%	0.604%
4	3.136	409	418	438	rVB	352213	806530	5.49%	0.467%
5	3.489	466	476	492	rBV	523578	1176987	8.02%	0.681%
6	4.373	612	621	634	rVB	330845	954502	6.50%	0.552%
7	5.586	788	820	828	rBV5	592413	2854310	19.44%	1.652%
8	5.666	828	833	838	rBV	138007	375618	2.56%	0.217%
9	5.909	860	873	890	rBV2	638411	1916776	13.05%	1.109%
10	6.074	890	900	905	rVV	157165	409039	2.79%	0.237%
11	6.141	905	911	919	rVV	141656	382569	2.61%	0.221%
12	6.245	919	928	937	rVV2	197380	723429	4.93%	0.419%
13	6.342	937	944	953	rVV3	231411	679193	4.63%	0.393%
14	6.446	953	961	974	rVB2	239277	691357	4.71%	0.400%
15	6.696	990	1002	1020	rBV	1056328	2569695	17.50%	1.487%
16	6.867	1020	1030	1036	rVV2	336623	982179	6.69%	0.568%
17	6.934	1037	1041	1053	rVB4	224774	569159	3.88%	0.329%
18	7.458	1115	1127	1138	rBV	1459514	3355050	22.85%	1.941%
19	7.574	1138	1146	1150	rBV2	182749	376349	2.56%	0.218%
20	7.635	1150	1156	1165	rVB	239454	489290	3.33%	0.283%
21	7.732	1165	1172	1182	rVB	183215	368327	2.51%	0.213%
22	7.842	1182	1190	1198	rVB2	199923	419923	2.86%	0.243%
23	8.257	1254	1258	1262	rVB2	196867	324915	2.21%	0.188%
24	8.342	1267	1272	1275	rVV	973331	1625427	11.07%	0.940%
25	8.379	1275	1278	1284	rVV2	506064	823160	5.61%	0.476%
26	8.500	1293	1298	1303	rVV	965305	1685998	11.48%	0.976%
27	8.574	1307	1310	1319	rVB4	189529	356911	2.43%	0.207%
28	8.665	1319	1325	1329	rBV3	409233	818750	5.58%	0.474%
29	8.720	1329	1334	1341	rVB	808925	1465292	9.98%	0.848%
30	8.976	1370	1376	1382	rBV2	1373347	2107351	14.35%	1.219%
31	9.043	1382	1387	1398	rVB4	233536	570168	3.88%	0.330%
32	9.165	1398	1407	1413	rBV3	188217	439070	2.99%	0.254%
33	9.226	1413	1417	1421	rBV2	232816	340867	2.32%	0.197%
34	9.445	1448	1453	1458	rBV2	244511	363934	2.48%	0.211%

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35	9.555	1467	1471	1478	rBV3	399876	767753	5.23%	0.444%
36	9.622	1478	1482	1486	rVB2	241033	328247	2.24%	0.190%
37	9.823	1508	1515	1520	rBV2	189343	378779	2.58%	0.219%
38	9.878	1520	1524	1528	rVV2	191331	366960	2.50%	0.212%
39	9.927	1528	1532	1534	rVV2	208222	348822	2.38%	0.202%
40	9.964	1534	1538	1543	rVV	1195678	1652095	11.25%	0.956%
41	10.067	1547	1555	1559	rVB	997368	1315308	8.96%	0.761%
42	10.116	1559	1563	1568	rBV	719399	1012189	6.89%	0.586%
43	10.256	1581	1586	1591	rBV	3017574	3827758	26.07%	2.215%
44	10.366	1598	1604	1609	rVV	9230032	12624999	85.98%	7.305%
45	10.415	1609	1612	1623	rVB	1377755	1962048	13.36%	1.135%
46	10.646	1645	1650	1653	rBV	333415	440171	3.00%	0.255%
47	10.835	1673	1681	1685	rBV	371749	652864	4.45%	0.378%
48	10.915	1685	1694	1698	rBV2	274173	493220	3.36%	0.285%
49	11.024	1705	1712	1718	rBV	402635	668118	4.55%	0.387%
50	11.085	1718	1722	1725	rBV2	272848	352031	2.40%	0.204%
51	11.140	1725	1731	1734	rBV2	1101515	1731534	11.79%	1.002%
52	11.250	1742	1749	1752	rBV	548820	758201	5.16%	0.439%
53	11.366	1763	1768	1775	rVB	1618957	1943121	13.23%	1.124%
54	11.439	1775	1780	1787	rVV	6129790	10805928	73.59%	6.252%
55	11.512	1787	1792	1803	rVB2	3629016	5841136	39.78%	3.380%
56	11.677	1814	1819	1826	rBV	1863576	2587855	17.62%	1.497%
57	11.817	1837	1842	1851	rVB	11667938	14683551	100.00%	8.496%
58	11.994	1867	1871	1874	rBV2	252347	383257	2.61%	0.222%
59	12.030	1874	1877	1880	rVB	370322	403668	2.75%	0.234%
60	12.085	1880	1886	1891	rBV2	1127586	1619525	11.03%	0.937%
61	12.146	1891	1896	1901	rBV	2630113	3637296	24.77%	2.105%
62	12.231	1908	1910	1917	rVB2	305157	345783	2.35%	0.200%
63	12.305	1917	1922	1924	rBV	2487687	3302535	22.49%	1.911%
64	12.329	1924	1926	1929	rVV	2512544	2636977	17.96%	1.526%
65	12.378	1929	1934	1941	rVV3	4191939	6330749	43.11%	3.663%
66	12.481	1945	1951	1954	rVV2	950612	1317913	8.98%	0.763%
67	12.524	1954	1958	1964	rVV	911213	1163294	7.92%	0.673%
68	12.591	1964	1969	1971	rVV	1853890	2313656	15.76%	1.339%
69	12.616	1971	1973	1978	rVV	1845666	2138490	14.56%	1.237%
70	12.677	1978	1983	1986	rVV	3112823	3947658	26.88%	2.284%
71	12.707	1986	1988	1992	rVV	590968	721453	4.91%	0.417%

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72	12.762	1992	1997	2005	rVV2	1527021	2540339	17.30%	1.470%
73	12.908	2017	2021	2025	rVV2	699727	892448	6.08%	0.516%
74	12.981	2025	2033	2037	rVV	1838605	2827482	19.26%	1.636%
75	13.024	2037	2040	2046	rVV	2642195	3168183	21.58%	1.833%
76	13.103	2046	2053	2056	rVV4	867487	1965610	13.39%	1.137%
77	13.128	2056	2057	2061	rVV	689570	708975	4.83%	0.410%
78	13.231	2066	2074	2077	rVV3	1856223	3132611	21.33%	1.813%
79	13.274	2077	2081	2084	rVV2	659252	878614	5.98%	0.508%
80	13.323	2084	2089	2091	rVV4	1131621	1734999	11.82%	1.004%
81	13.353	2091	2094	2100	rVV2	2990310	4479589	30.51%	2.592%
82	13.432	2104	2107	2112	rVV2	739717	1033043	7.04%	0.598%
83	13.487	2112	2116	2124	rVB3	568360	972961	6.63%	0.563%
84	13.567	2124	2129	2132	rBV2	595574	838957	5.71%	0.485%
85	13.609	2132	2136	2138	rVV	1009368	1398130	9.52%	0.809%
86	13.646	2138	2142	2148	rVV3	1160063	2356897	16.05%	1.364%
87	13.725	2152	2155	2161	rVV2	982019	1745429	11.89%	1.010%
88	13.780	2161	2164	2169	rVV2	323559	605604	4.12%	0.350%
89	13.841	2169	2174	2179	rVV	1807742	2493968	16.98%	1.443%
90	13.890	2179	2182	2184	rVV2	231352	339007	2.31%	0.196%
91	13.920	2185	2187	2192	rVB3	361728	499349	3.40%	0.289%
92	13.987	2192	2198	2203	rBV4	610070	1135894	7.74%	0.657%
93	14.036	2203	2206	2209	rVV2	345783	433956	2.96%	0.251%
94	14.097	2213	2216	2219	rVB3	489906	545273	3.71%	0.315%
95	14.140	2219	2223	2227	rBV2	842836	1006643	6.86%	0.582%
96	14.280	2241	2246	2255	rBV2	747963	1280804	8.72%	0.741%
97	14.384	2260	2263	2268	rVV	312416	412998	2.81%	0.239%
98	14.438	2268	2272	2276	rVB	490758	589991	4.02%	0.341%
99	14.701	2311	2315	2325	rVB	2026160	2643316	18.00%	1.529%
100	14.847	2335	2339	2348	rVB	702731	961489	6.55%	0.556%

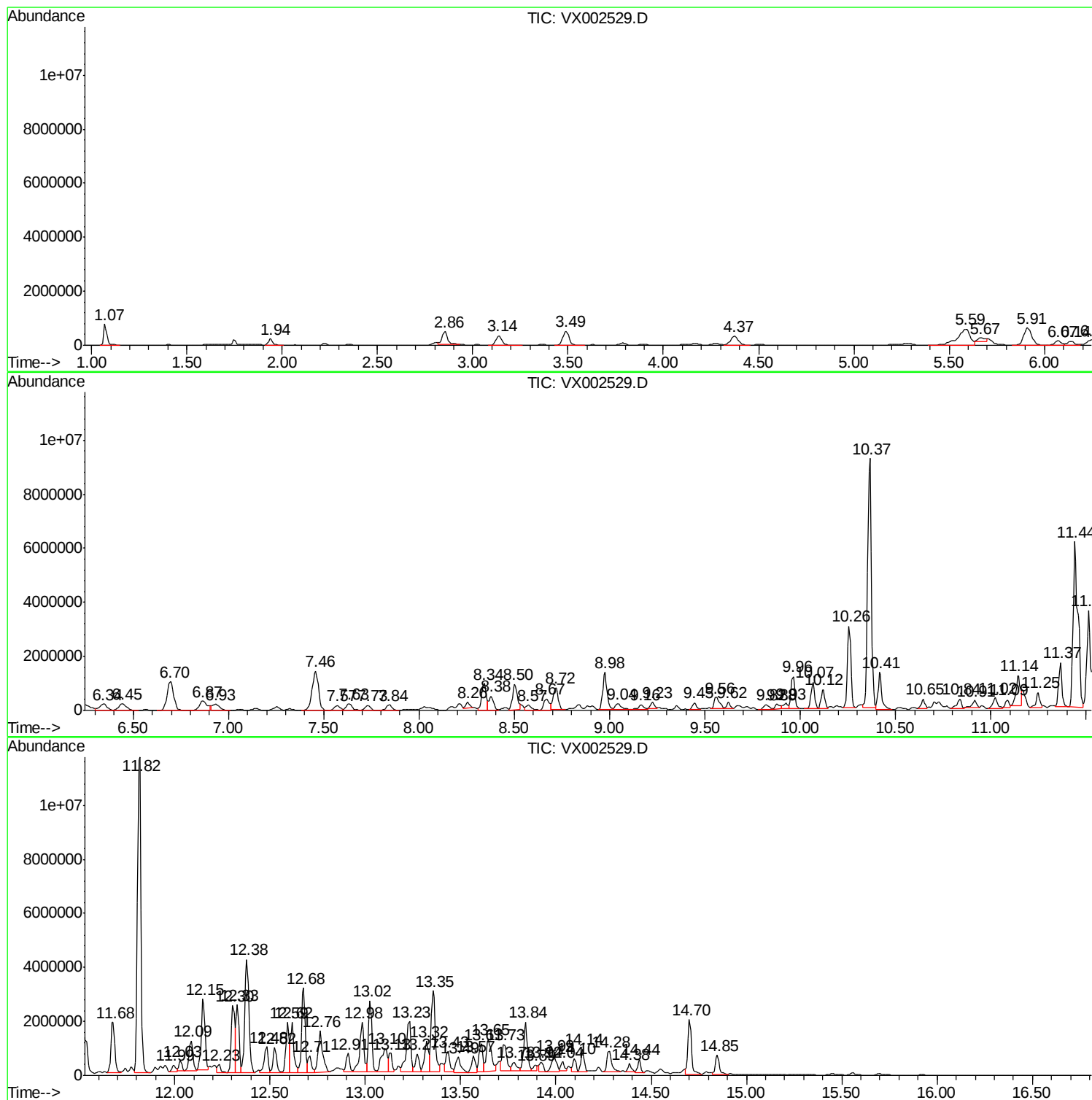
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Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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



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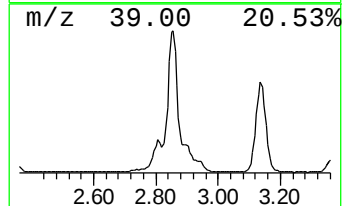
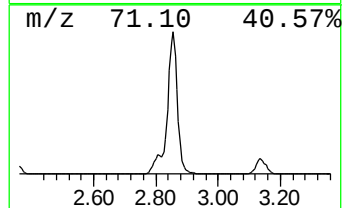
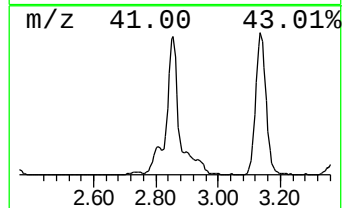
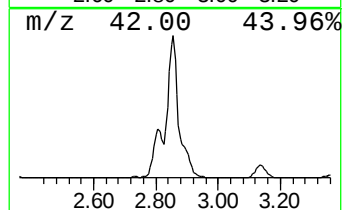
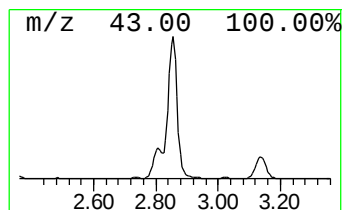
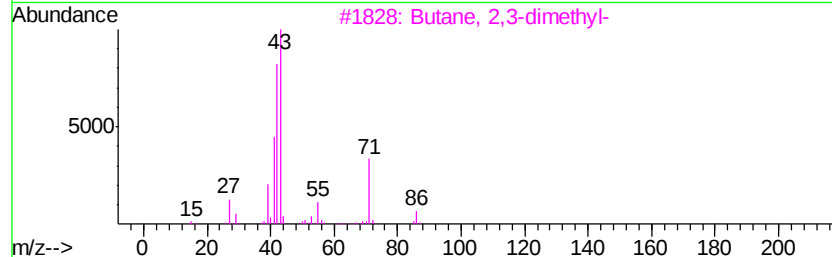
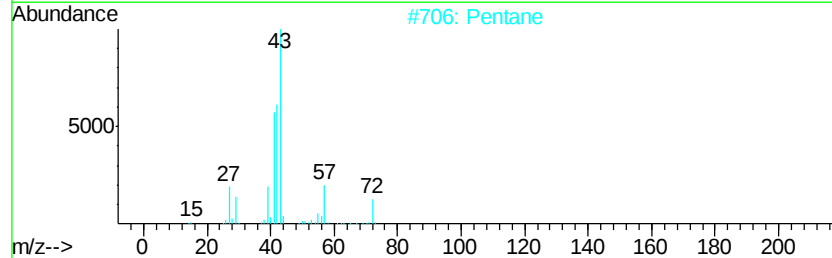
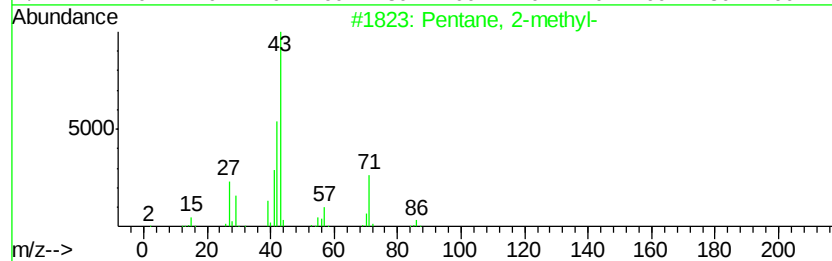
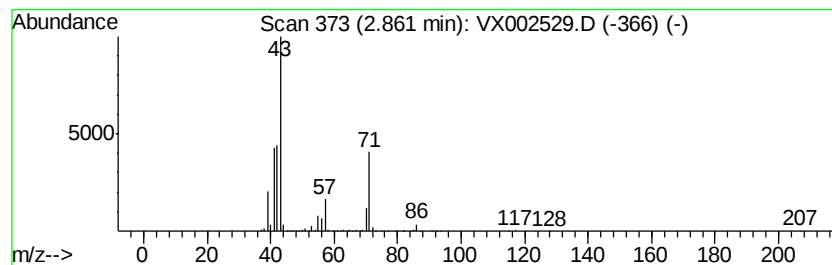
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

Peak Number 2 Pentane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.86	138.96 ug/l	1043950	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2		Pentane	72	C5H12	000109-66-0	46
3		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	43
4		Heptane, 4-methyl-	114	C8H18	000589-53-7	28
5		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	28



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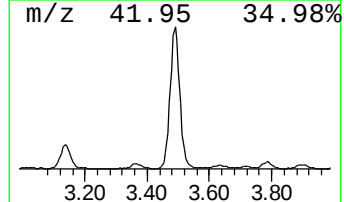
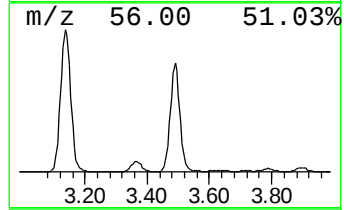
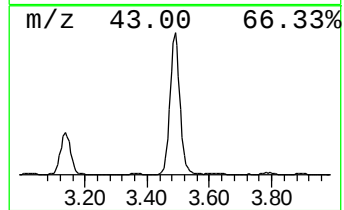
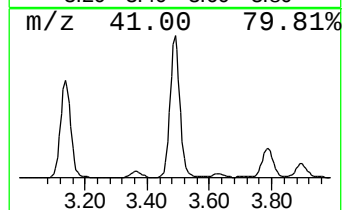
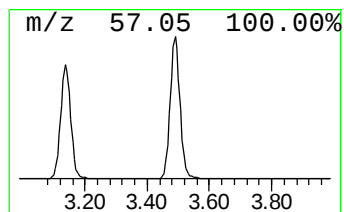
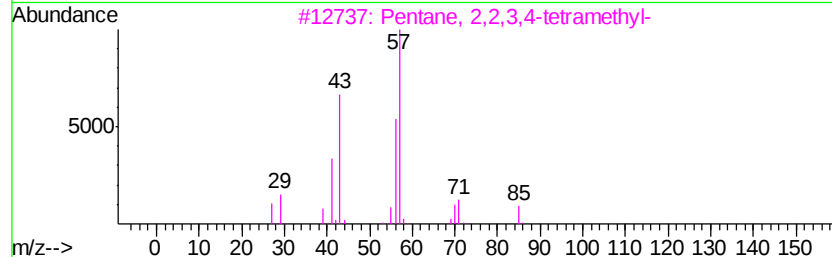
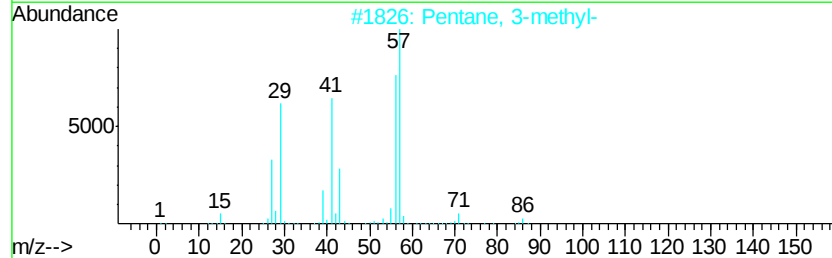
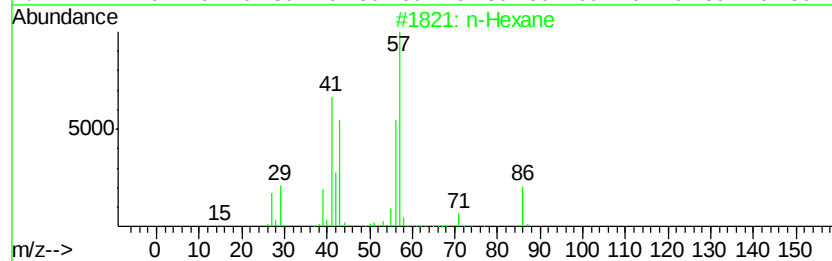
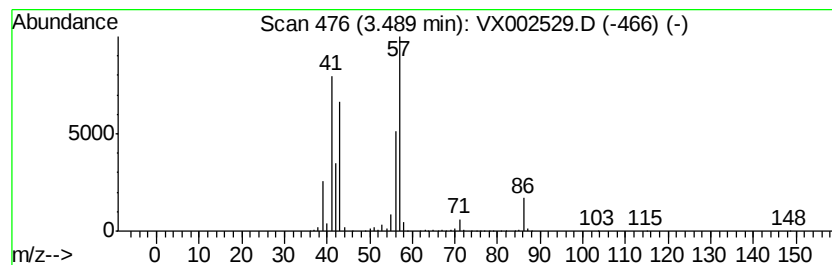
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.49	156.67 ug/l	1176990	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexane	86	C6H14	000110-54-3	91
2		Pentane, 3-methyl-	86	C6H14	000096-14-0	40
3		Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	40
4		Propanal, 2,2-dimethyl-	86	C5H10O	000630-19-3	38
5		Butanal, 2-methyl-	86	C5H10O	000096-17-3	38



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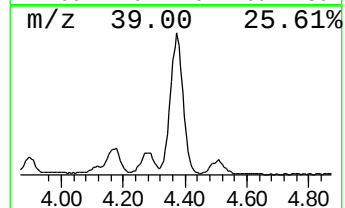
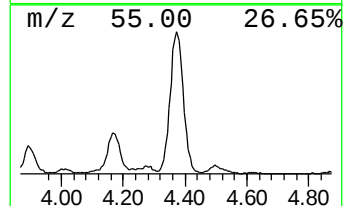
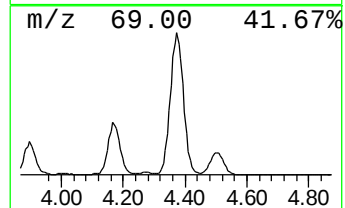
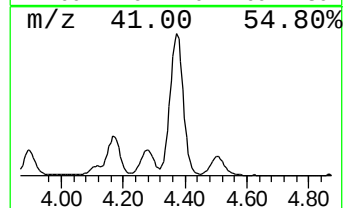
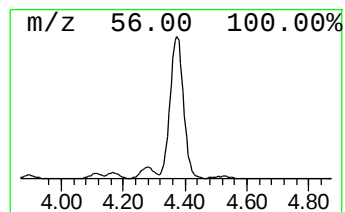
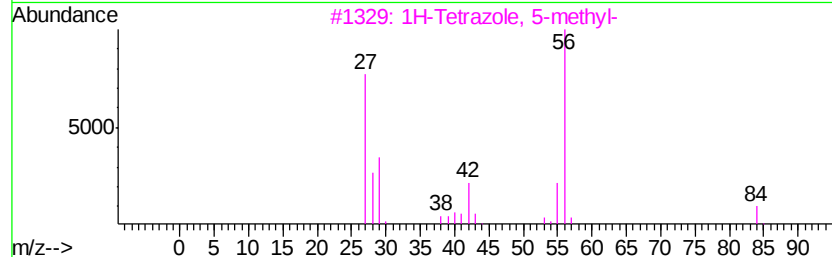
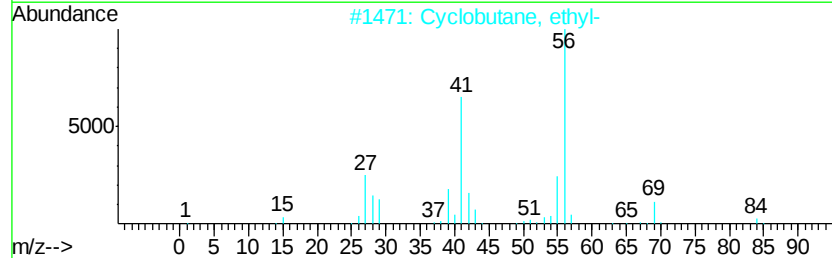
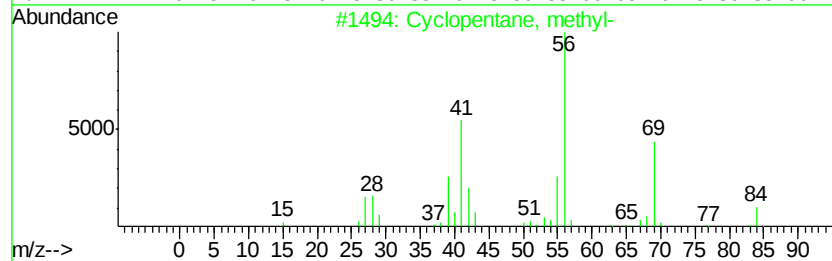
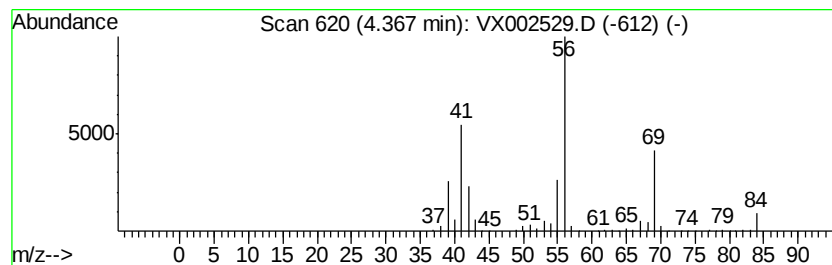
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

Peak Number 4 Cyclopentane, methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.37	127.06 ug/l	954502	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2		Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
3		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	56
4		Cyclobutane	56	C4H8	000287-23-0	47
5		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002529.D
Acq On : 14 Jun 2018 01:21
Operator : JC/MD
Sample : J3413-04
Misc : 6.70uL/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-5ARR

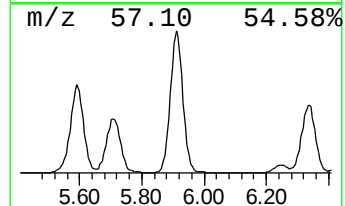
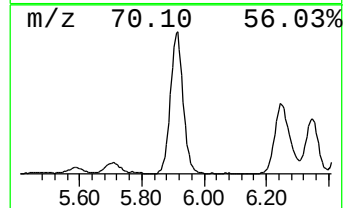
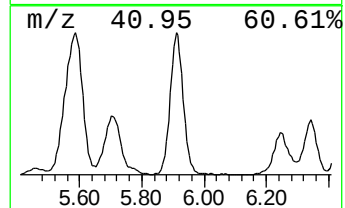
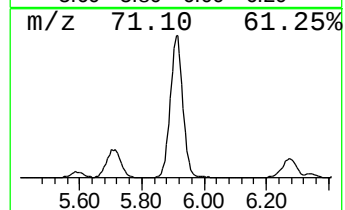
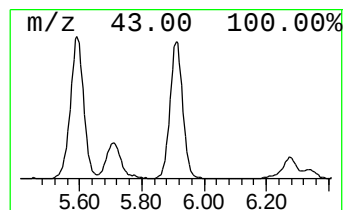
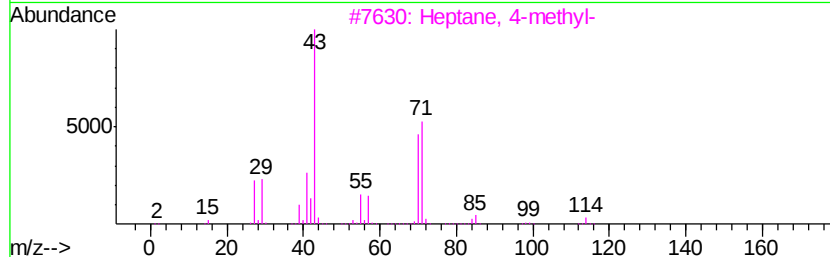
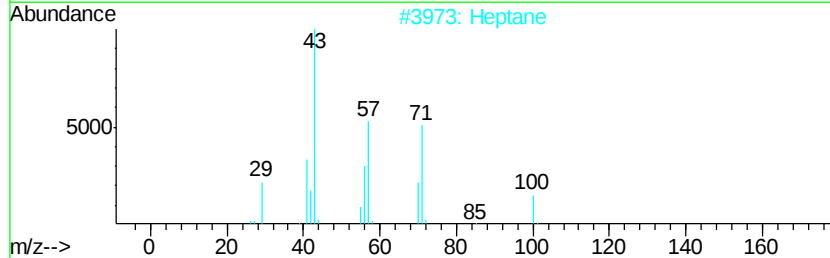
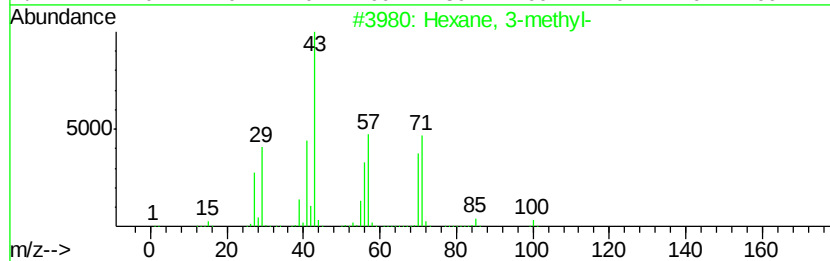
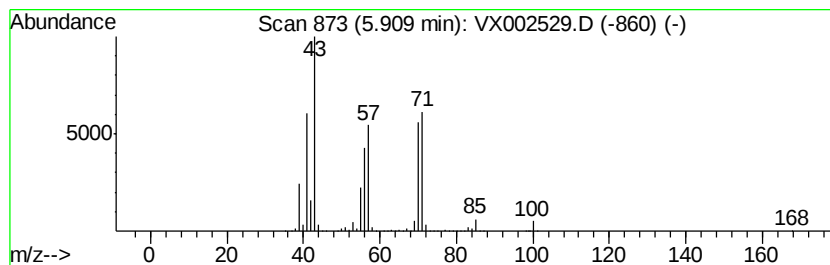
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Quant Title : SW846 8260

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

Peak Number 5 Hexane, 3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.91	255.15 ug/l	1916780	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexane, 3-methyl-	100	C7H16	000589-34-4	91
2		Heptane	100	C7H16	000142-82-5	80
3		Heptane, 4-methyl-	114	C8H18	000589-53-7	59
4		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	53
5		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002529.D
Acq On : 14 Jun 2018 01:21
Operator : JC/MD
Sample : J3413-04
Misc : 6.70uL/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-5ARR

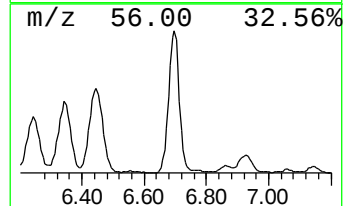
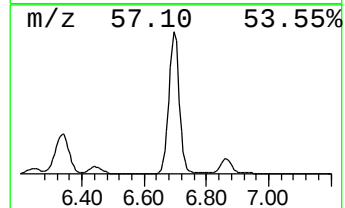
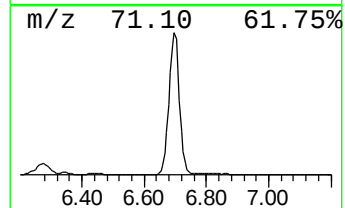
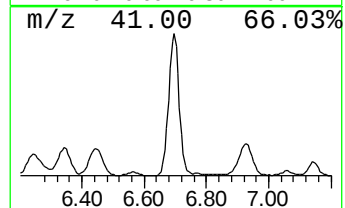
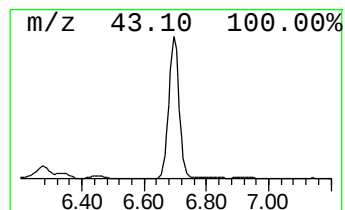
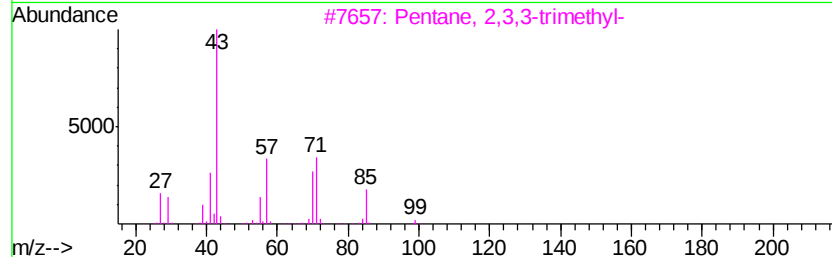
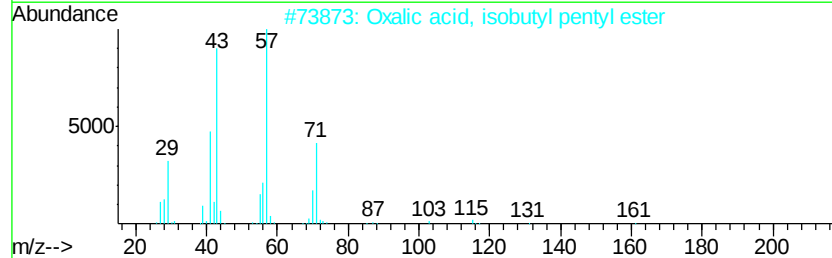
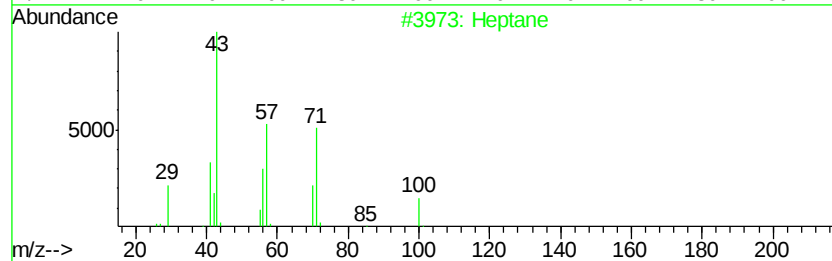
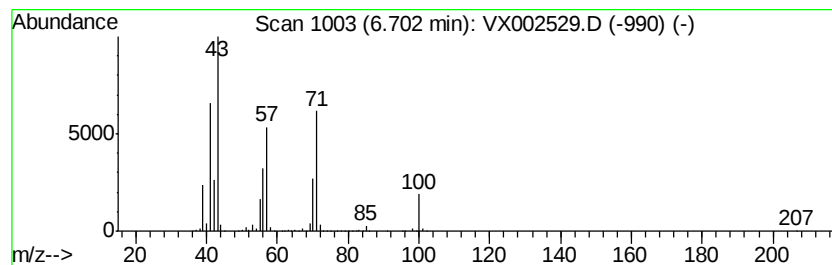
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Quant Title : SW846 8260

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

Peak Number 6 Heptane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.70	130.82 ug/l	2569700	1,4-Difluorobenzene	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane	100	C7H16	000142-82-5	91
2		Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	50
3		Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	43
4		Hexane, 3-methyl-	100	C7H16	000589-34-4	40
5		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	40



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002529.D
Acq On : 14 Jun 2018 01:21
Operator : JC/MD
Sample : J3413-04
Misc : 6.70uL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-5ARR

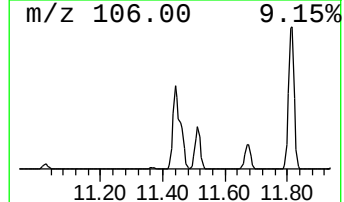
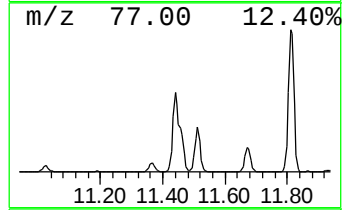
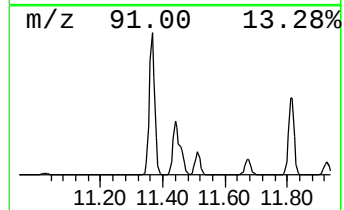
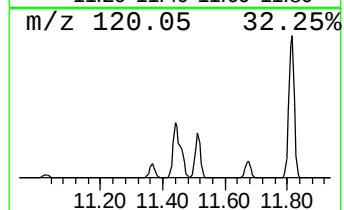
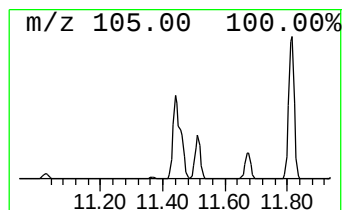
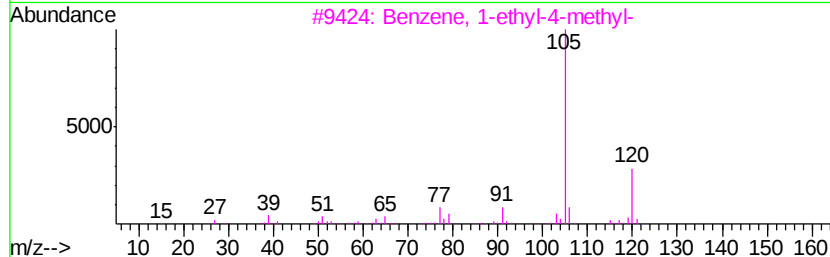
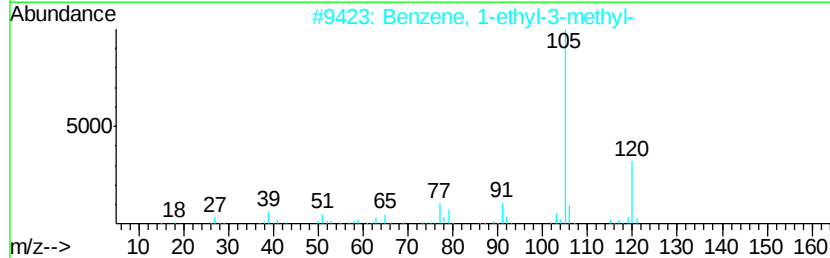
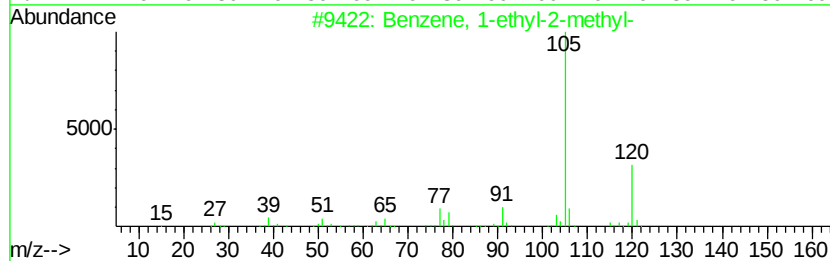
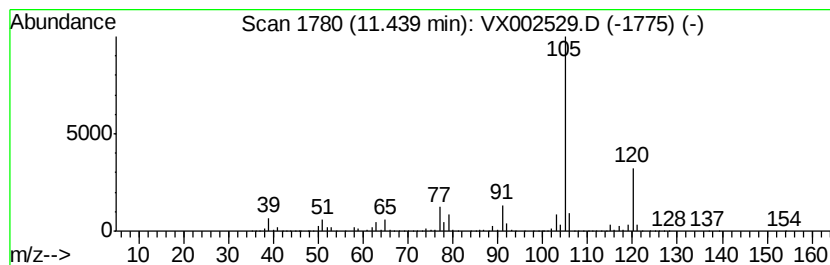
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	333.61 ug/l	10805900	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002529.D
Acq On : 14 Jun 2018 01:21
Operator : JC/MD
Sample : J3413-04
Misc : 6.70uL/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 32 Sample Multiplier: 1

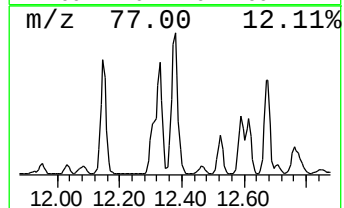
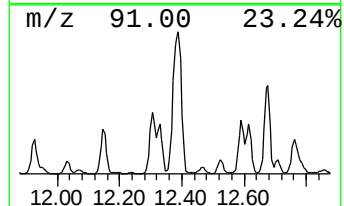
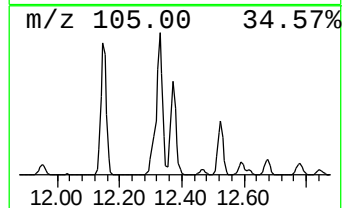
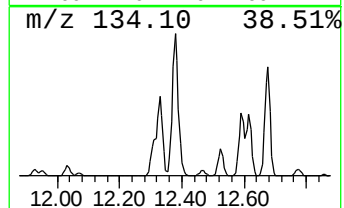
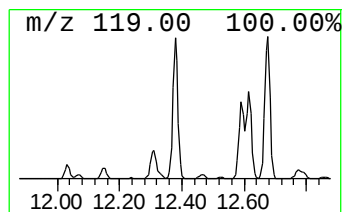
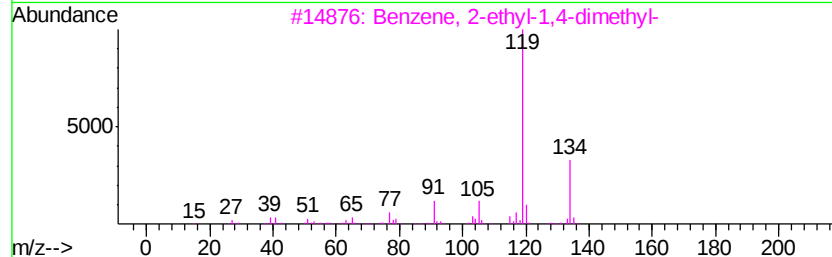
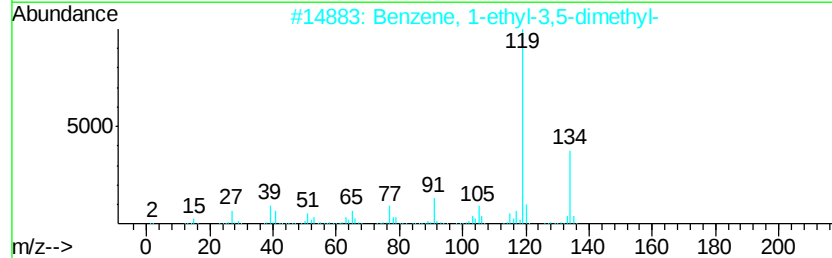
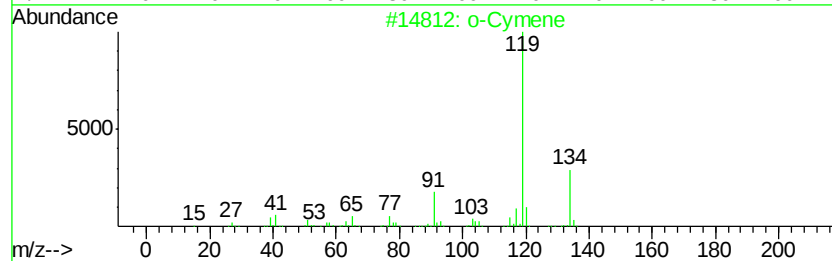
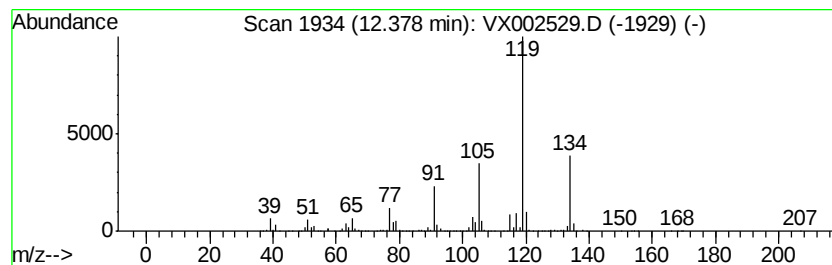
Instrument :
MSVOA_X
ClientSampleId :
B-5ARR

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

Peak Number 8 o-Cymene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.38	195.45 ug/l	6330750	1,4-Dichlorobenzene-d4	12.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		o-Cymene	134 C10H14	000527-84-4 95
2		Benzene, 1-ethyl-3,5-dimethyl-	134 C10H14	000934-74-7 94
3		Benzene, 2-ethyl-1,4-dimethyl-	134 C10H14	001758-88-9 94
4		Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2 93
5		Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9 93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002529.D
Acq On : 14 Jun 2018 01:21
Operator : JC/MD
Sample : J3413-04
Misc : 6.70µL/100µL/5.0mL/MSVOA_X/MEOH
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-5ARR

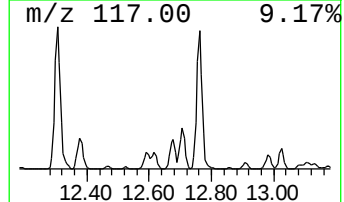
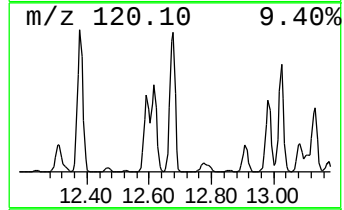
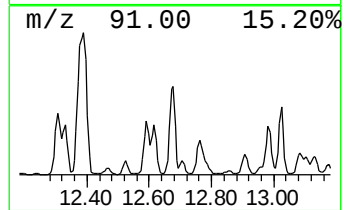
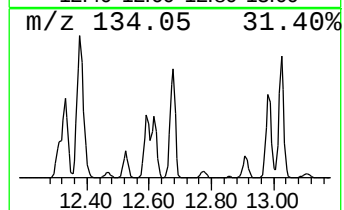
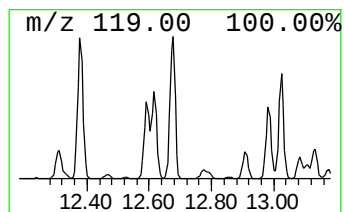
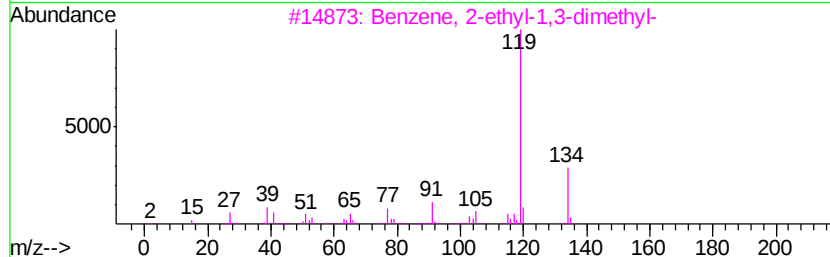
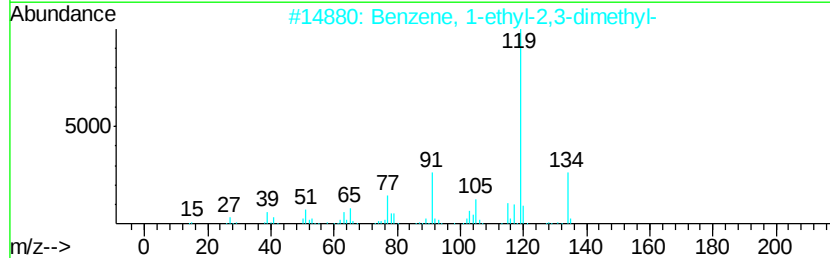
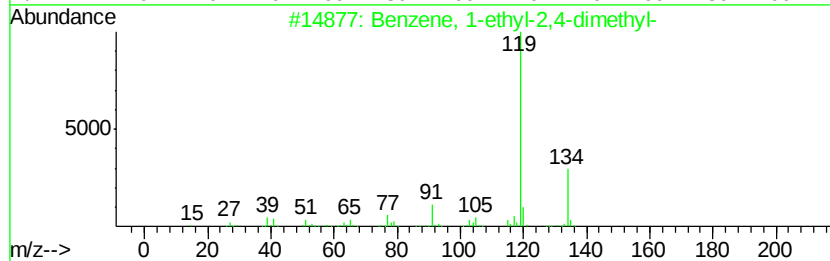
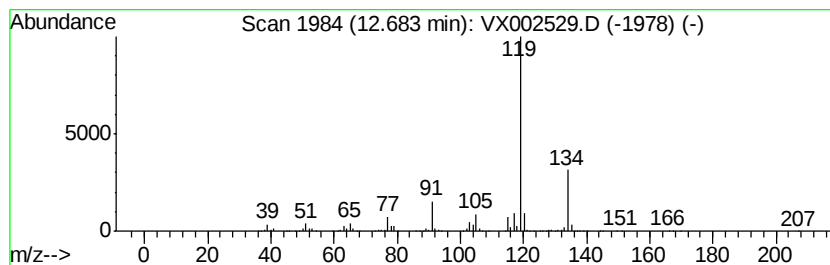
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.68	121.88 ug/l	3947660	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002529.D
Acq On : 14 Jun 2018 01:21
Operator : JC/MD
Sample : J3413-04
Misc : 6.70uL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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ClientSampleId :
B-5ARR

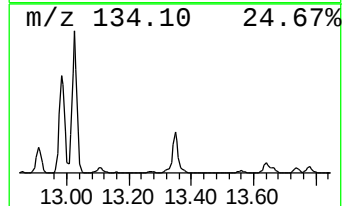
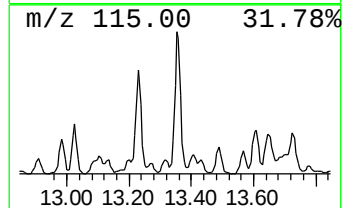
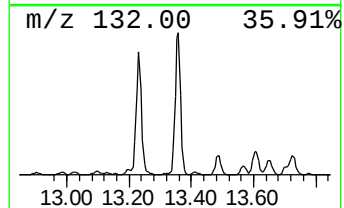
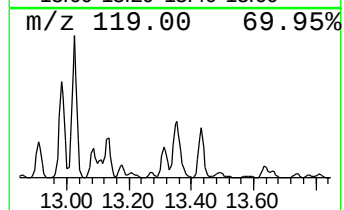
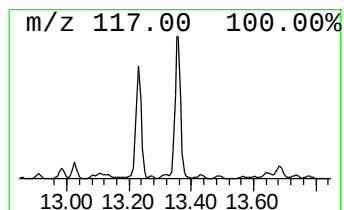
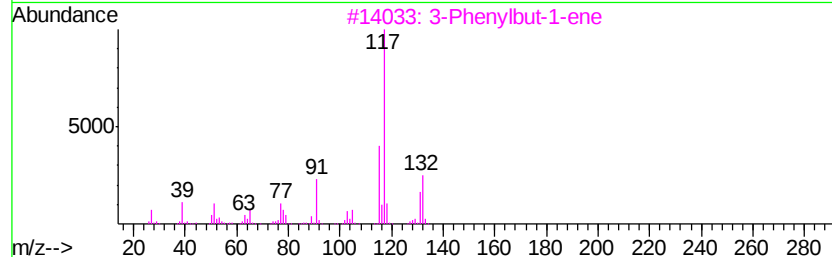
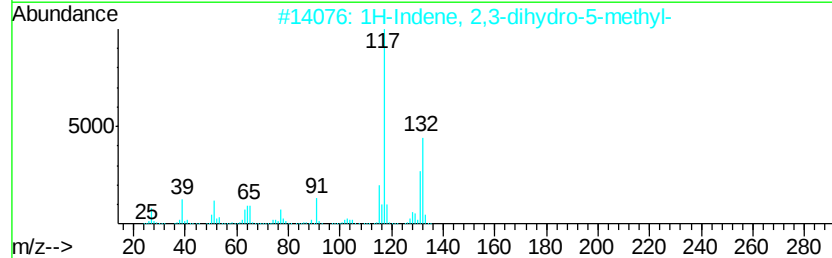
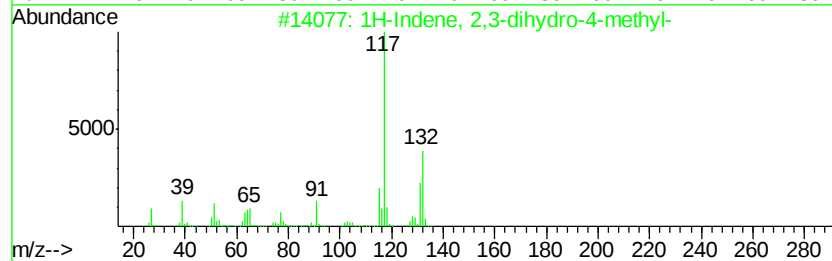
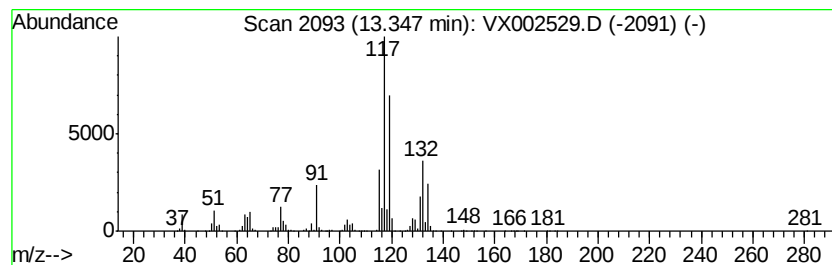
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Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	138.30 ug/l	4479590	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	86
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
4		Indan, 1-methyl-	132	C10H12	000767-58-8	76
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX061318\
Data File : VX002529.D
Acq On : 14 Jun 2018 01:21
Operator : JC/MD
Sample : J3413-04
Misc : 6.70g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-5ARR

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.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
Pentane, 2-methyl-	2.86	139.0	ug/l	1043950	1	5.66	375618	50.0
n-Hexane	3.49	156.7	ug/l	1176990	1	5.66	375618	50.0
Cyclopentane, met...	4.37	127.1	ug/l	954502	1	5.66	375618	50.0
Hexane, 3-methyl-	5.91	255.2	ug/l	1916780	1	5.66	375618	50.0
Heptane	6.70	130.8	ug/l	2569700	2	6.87	982179	50.0
Benzene, 1-ethyl-...	11.44	333.6	ug/l	10805900	4	12.09	1619530	50.0
o-Cymene	12.38	195.4	ug/l	6330750	4	12.09	1619530	50.0
Benzene, 1-ethyl-...	12.68	121.9	ug/l	3947660	4	12.09	1619530	50.0
1H-Indene, 2,3-di...	13.35	138.3	ug/l	4479590	4	12.09	1619530	50.0