

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061318\
 Data File : VX002532.D
 Acq On : 14 Jun 2018 02:41
 Operator : JC/MD
 Sample : J3383-14
 Misc : 6.28g/5mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 B-42R(I-2R)R

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	891018	1110488	8.35%	0.650%
2	1.940	216	222	236	rVB	574035	866131	6.52%	0.507%
3	2.855	366	372	384	rVB	1019210	2235204	16.82%	1.308%
4	3.135	409	418	436	rVB	711446	1628856	12.25%	0.953%
5	3.489	467	476	491	rBV	1166312	2618965	19.70%	1.533%
6	3.788	517	525	534	rVB	202246	481272	3.62%	0.282%
7	4.172	581	588	596	rVV	162015	457501	3.44%	0.268%
8	4.373	612	621	633	rVV	756791	2204249	16.58%	1.290%
9	5.281	744	770	787	rBV2	204746	851518	6.41%	0.498%
10	5.586	787	820	828	rBV5	988236	4739265	35.66%	2.774%
11	5.915	861	874	889	rBV2	986931	2984558	22.45%	1.747%
12	6.074	889	900	906	rVV	166912	455803	3.43%	0.267%
13	6.245	918	928	937	rVV3	330302	1200042	9.03%	0.702%
14	6.348	938	945	952	rVV2	344341	972258	7.31%	0.569%
15	6.446	952	961	973	rVB2	410945	1181658	8.89%	0.692%
16	6.696	989	1002	1020	rBV	1538273	3745283	28.18%	2.192%
17	6.866	1021	1030	1036	rBV2	378790	1200068	9.03%	0.702%
18	6.934	1037	1041	1054	rVB3	386675	1005286	7.56%	0.588%
19	7.251	1084	1093	1100	rBV2	220417	490013	3.69%	0.287%
20	7.458	1114	1127	1138	rBV	2305651	5395268	40.59%	3.158%
21	7.568	1138	1145	1150	rBV2	241506	496022	3.73%	0.290%
22	7.635	1150	1156	1165	rVB	320305	649598	4.89%	0.380%
23	7.726	1165	1171	1182	rVB	287536	574827	4.32%	0.336%
24	7.842	1182	1190	1198	rVB2	277371	593134	4.46%	0.347%
25	8.025	1215	1220	1236	rVB4	186267	676733	5.09%	0.396%
26	8.171	1236	1244	1247	rBV3	188904	404164	3.04%	0.237%
27	8.214	1247	1251	1254	rVV	278895	408079	3.07%	0.239%
28	8.256	1254	1258	1262	rVB2	253283	392008	2.95%	0.229%
29	8.342	1267	1272	1275	rVV	1217394	2028809	15.26%	1.188%
30	8.378	1275	1278	1284	rVV2	661842	1084762	8.16%	0.635%
31	8.500	1293	1298	1306	rVV	1206181	2350006	17.68%	1.376%
32	8.573	1306	1310	1319	rVB3	304747	604911	4.55%	0.354%
33	8.665	1319	1325	1329	rBV3	577443	1170245	8.80%	0.685%
34	8.720	1329	1334	1341	rVV	902704	1736462	13.06%	1.016%

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Integration Parameters: RTEINT.P

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35	8.836	1349	1353	1357	rVV2	281585	526467	3.96%	0.308%
36	8.976	1370	1376	1383	rVV2	1590460	2643879	19.89%	1.548%
37	9.043	1383	1387	1398	rVB3	337209	801601	6.03%	0.469%
38	9.226	1412	1417	1421	rBV2	352497	529524	3.98%	0.310%
39	9.445	1448	1453	1457	rBV	272448	393806	2.96%	0.231%
40	9.561	1467	1472	1478	rVB2	453140	881610	6.63%	0.516%
41	9.622	1478	1482	1486	rVB2	337764	444176	3.34%	0.260%
42	9.823	1508	1515	1520	rBV2	242193	479609	3.61%	0.281%
43	9.878	1520	1524	1528	rVV3	211955	436193	3.28%	0.255%
44	9.921	1528	1531	1534	rVV2	232422	399352	3.00%	0.234%
45	9.963	1534	1538	1543	rVV	1129299	1645756	12.38%	0.963%
46	10.067	1547	1555	1559	rVB	940990	1294613	9.74%	0.758%
47	10.116	1559	1563	1568	rBV	797278	1104253	8.31%	0.646%
48	10.256	1581	1586	1591	rVV	3915503	5125090	38.56%	3.000%
49	10.366	1591	1604	1609	rVV	4190829	6326466	47.60%	3.703%
50	10.415	1609	1612	1623	rVB	1062527	1616825	12.16%	0.946%
51	10.646	1644	1650	1653	rBV2	331022	468049	3.52%	0.274%
52	10.835	1673	1681	1685	rBV2	359768	641930	4.83%	0.376%
53	11.024	1706	1712	1718	rBV	492099	770465	5.80%	0.451%
54	11.140	1725	1731	1735	rBV	1075258	1793300	13.49%	1.050%
55	11.250	1742	1749	1752	rBV	410419	588050	4.42%	0.344%
56	11.366	1763	1768	1775	rVB	2094282	2517066	18.94%	1.473%
57	11.457	1775	1783	1787	rVV	2838872	4546510	34.21%	2.661%
58	11.512	1787	1792	1803	rVB2	2791740	4408708	33.17%	2.581%
59	11.670	1814	1818	1825	rBV	1034122	1460853	10.99%	0.855%
60	11.811	1837	1841	1850	rVB	10696743	13291705	100.00%	7.781%
61	12.085	1880	1886	1891	rBV	1143848	1606217	12.08%	0.940%
62	12.146	1891	1896	1901	rBV	3052464	3985393	29.98%	2.333%
63	12.304	1917	1922	1925	rBV	2642841	3686749	27.74%	2.158%
64	12.378	1929	1934	1940	rVB3	2949274	4768222	35.87%	2.791%
65	12.481	1940	1951	1954	rBV3	509445	927359	6.98%	0.543%
66	12.524	1954	1958	1964	rVV	884958	1096958	8.25%	0.642%
67	12.591	1964	1969	1971	rVV	1605986	2060796	15.50%	1.206%
68	12.615	1971	1973	1978	rVV	1537812	1738278	13.08%	1.018%
69	12.670	1978	1982	1986	rVV	3009474	3788883	28.51%	2.218%
70	12.707	1986	1988	1992	rVV	582905	662976	4.99%	0.388%
71	12.762	1992	1997	2005	rVV3	1494159	2401974	18.07%	1.406%

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Integration Parameters: RTEINT.P

Integrator: RTE
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 Stop Thrs : 0 Peak Location: TOP

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72	12.908	2017	2021	2025	rVV2	699033	851007	6.40%	0.498%
73	12.981	2025	2033	2037	rVV	1886343	2650203	19.94%	1.551%
74	13.024	2037	2040	2046	rVV	2639115	3090056	23.25%	1.809%
75	13.085	2046	2050	2051	rVV	493227	579677	4.36%	0.339%
76	13.103	2051	2053	2056	rVV3	643262	899062	6.76%	0.526%
77	13.128	2056	2057	2061	rVV	483542	474175	3.57%	0.278%
78	13.231	2069	2074	2077	rVV2	1843784	2577150	19.39%	1.509%
79	13.274	2077	2081	2084	rVV2	515204	684065	5.15%	0.400%
80	13.323	2084	2089	2091	rVV2	739581	1270055	9.56%	0.743%
81	13.353	2091	2094	2100	rVV2	3290847	4555696	34.27%	2.667%
82	13.432	2104	2107	2112	rVV2	637508	879457	6.62%	0.515%
83	13.487	2112	2116	2124	rVB2	558613	899830	6.77%	0.527%
84	13.566	2124	2129	2132	rBV2	562742	777327	5.85%	0.455%
85	13.603	2132	2135	2138	rVV	955427	1297413	9.76%	0.759%
86	13.646	2138	2142	2147	rVV3	1098773	2129567	16.02%	1.247%
87	13.701	2148	2151	2152	rVV3	381489	468743	3.53%	0.274%
88	13.725	2152	2155	2161	rVV2	965772	1616143	12.16%	0.946%
89	13.780	2161	2164	2169	rVV2	230736	433982	3.27%	0.254%
90	13.841	2169	2174	2179	rVV	2416387	3213425	24.18%	1.881%
91	13.920	2185	2187	2192	rVB3	317284	423675	3.19%	0.248%
92	13.987	2192	2198	2203	rBV3	569024	997775	7.51%	0.584%
93	14.036	2203	2206	2209	rVV	322348	416420	3.13%	0.244%
94	14.097	2213	2216	2219	rVB3	347693	386567	2.91%	0.226%
95	14.140	2219	2223	2227	rBV	817612	1024282	7.71%	0.600%
96	14.274	2241	2245	2255	rVV2	738236	1298387	9.77%	0.760%
97	14.383	2259	2263	2268	rVV	319439	413008	3.11%	0.242%
98	14.438	2268	2272	2276	rVB	482652	592550	4.46%	0.347%
99	14.700	2311	2315	2325	rVB	2848283	3639659	27.38%	2.131%
100	14.847	2335	2339	2348	rVB	1055937	1440521	10.84%	0.843%

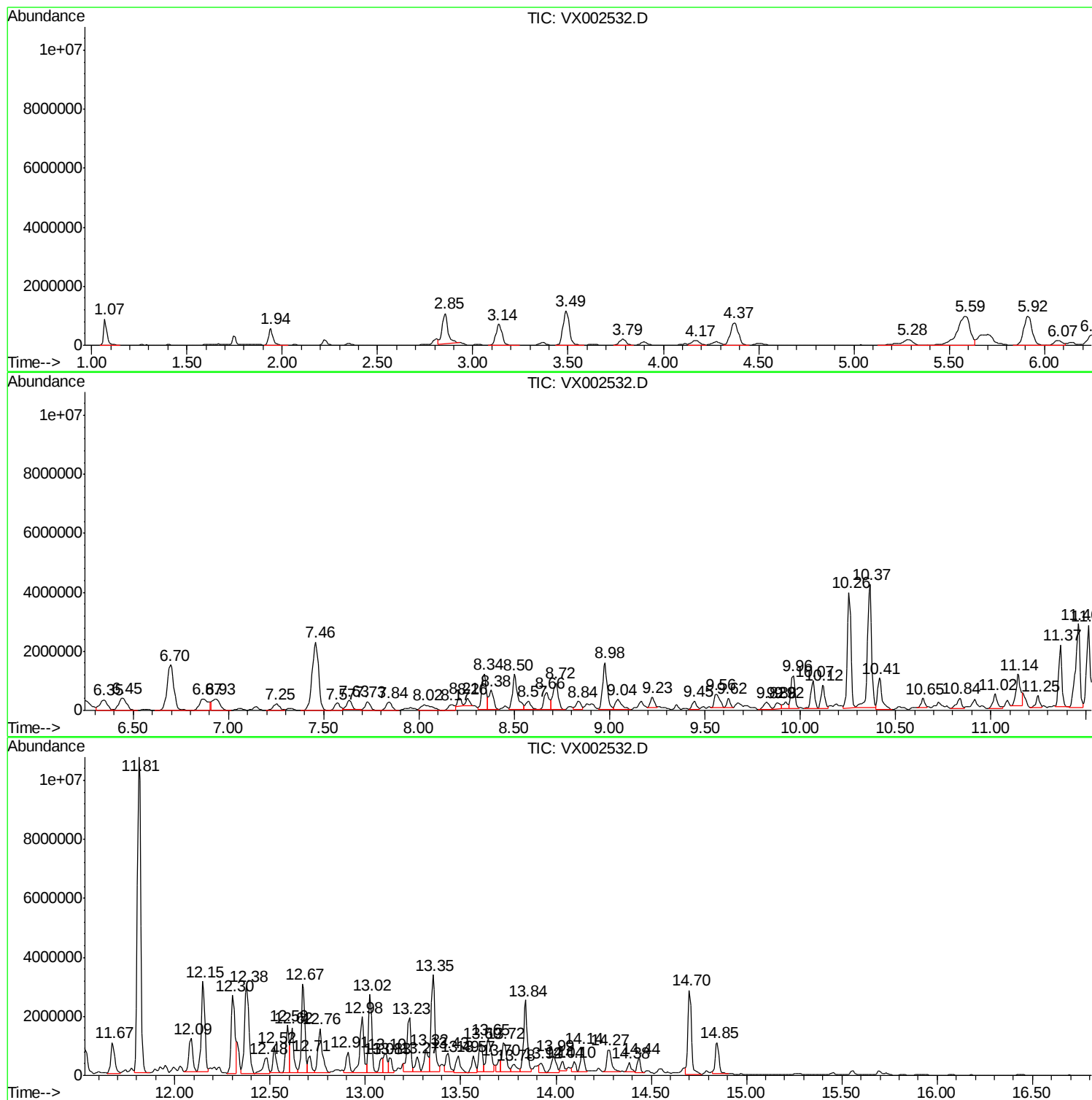
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Instrument :
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B-42R(I-2R)R

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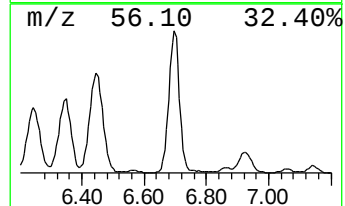
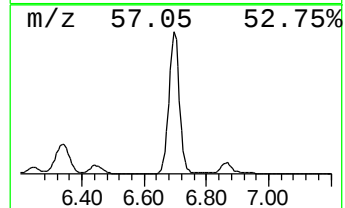
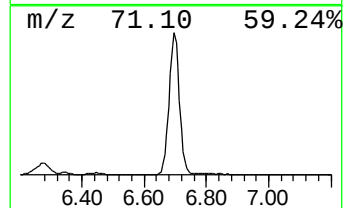
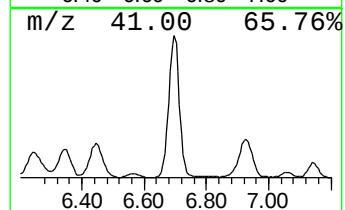
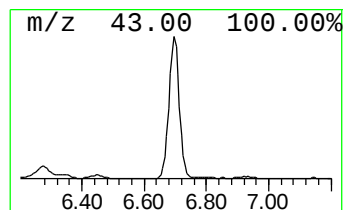
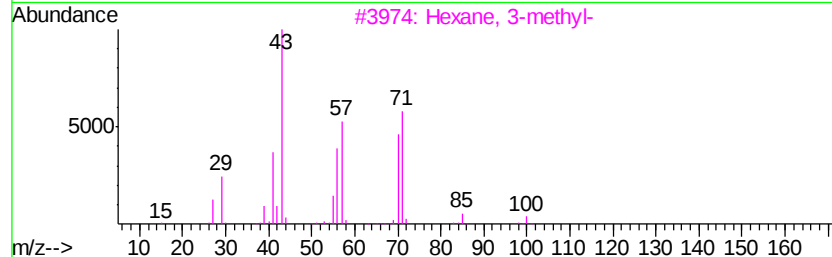
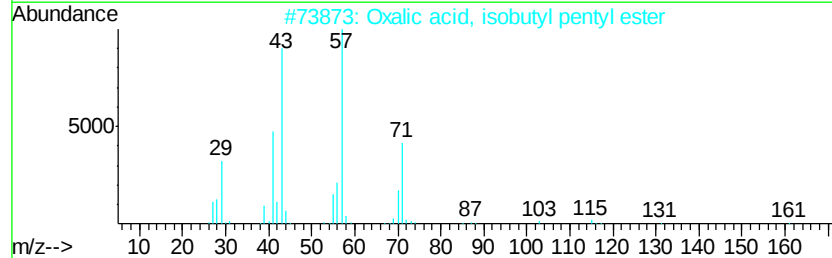
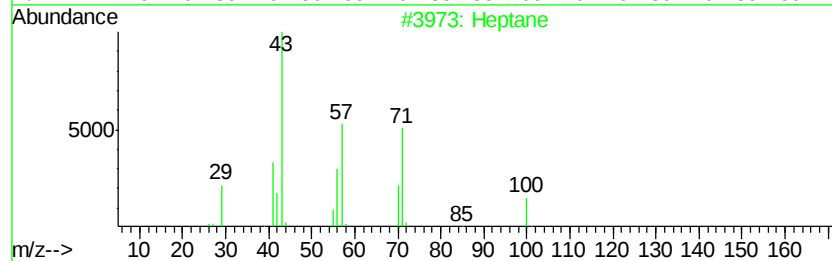
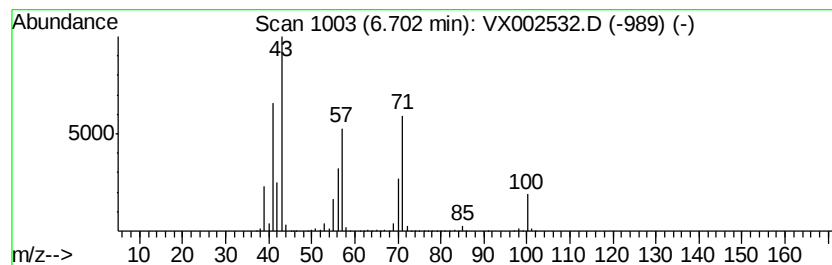
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

Peak Number 1 Heptane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.70	156.05 ug/l	3745280	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		Hexane, 3-methyl-	100	C7H16	000589-34-4	40
4		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	40
5		Acetaldehyde, propylhydrazone	100	C5H12N2	007422-88-0	38



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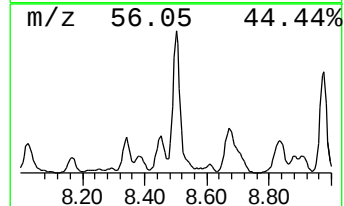
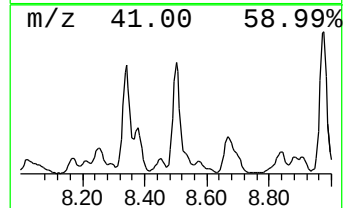
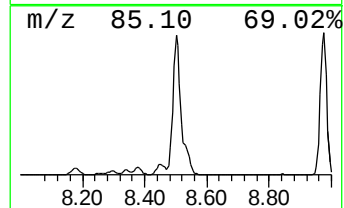
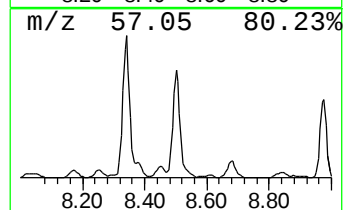
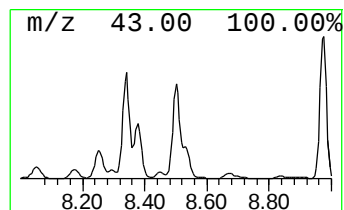
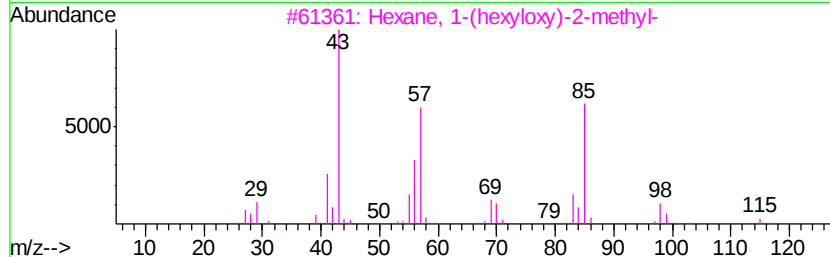
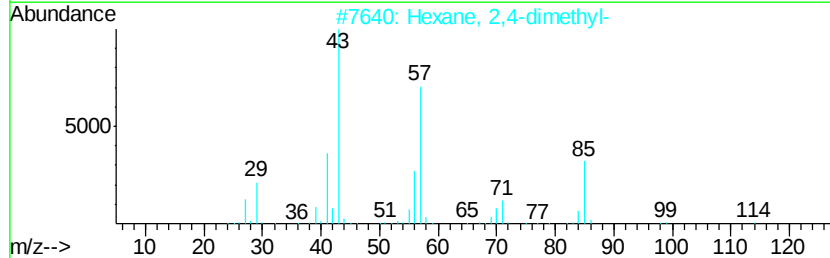
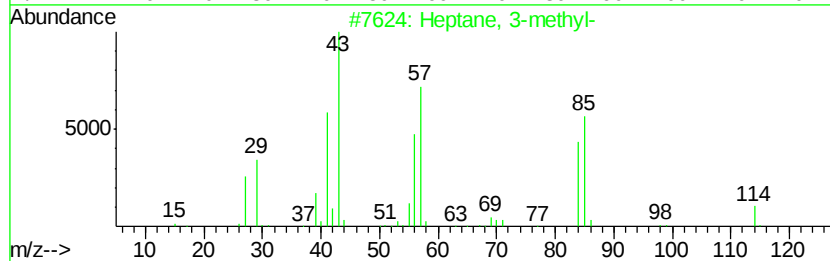
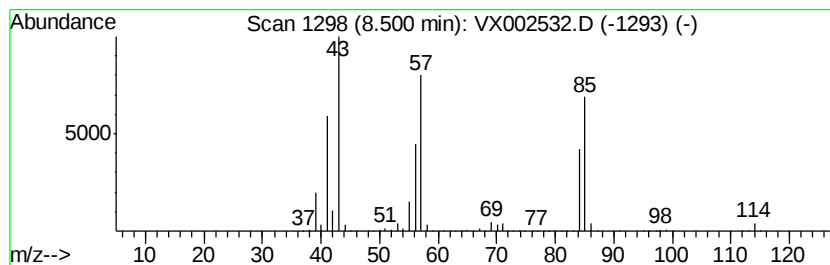
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Quant Title : SW846 8260

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

Peak Number 2 Heptane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.50	106.41 ug/l	2350010	Chlorobenzene-d5	10.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane, 3-methyl-	114	C8H18	000589-81-1	94
2		Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	59
3		Hexane, 1-(hexyloxy)-2-methyl-	200	C13H28O	074421-17-3	59
4		Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	59
5		Hexane, 2-methyl-	100	C7H16	000591-76-4	53



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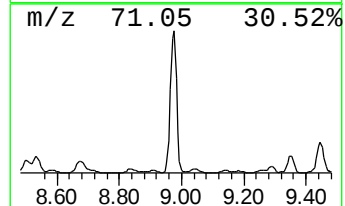
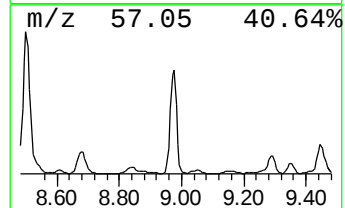
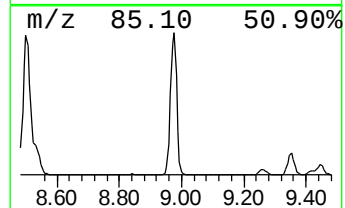
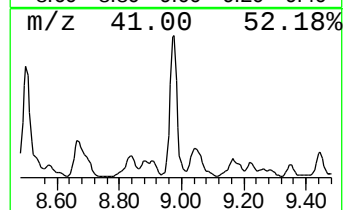
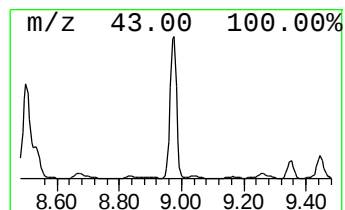
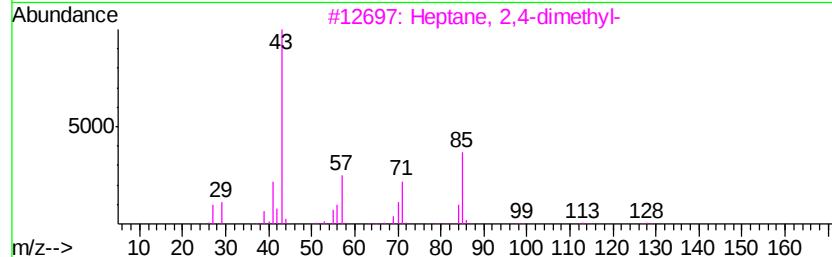
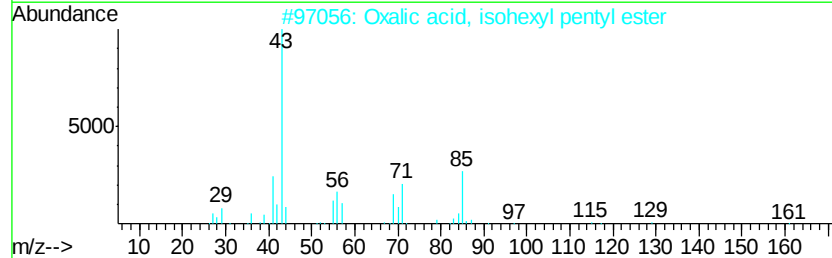
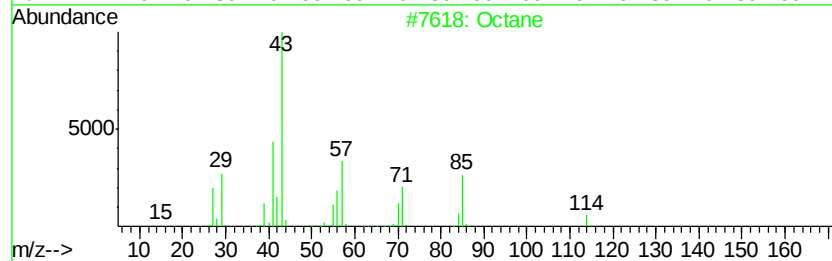
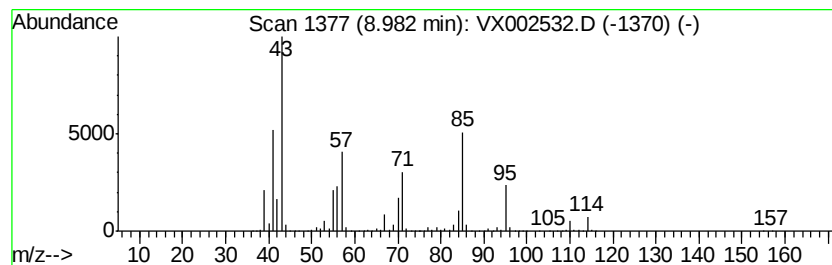
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Quant Title : SW846 8260

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

Peak Number 3 Octane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.98	119.71 ug/l	2643880	Chlorobenzene-d5	10.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane		114	C8H18	000111-65-9	93
2	Oxalic acid, isohexyl pentyl ester		244	C13H24O4	1000309-32-8	78
3	Heptane, 2,4-dimethyl-		128	C9H20	002213-23-2	64
4	Undecane, 2,4-dimethyl-		184	C13H28	017312-80-0	59
5	Heptane, 2,3-dimethyl-		128	C9H20	003074-71-3	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002532.D
Acq On : 14 Jun 2018 02:41
Operator : JC/MD
Sample : J3383-14
Misc : 6.28µ/5mL/100µL/5.0mL/MSVOA_X/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
B-42R(I-2R)R

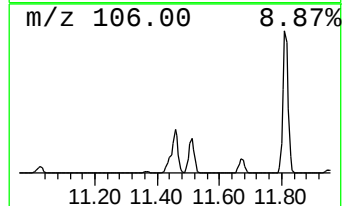
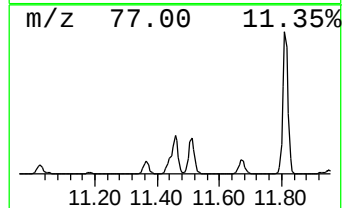
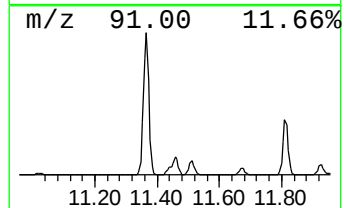
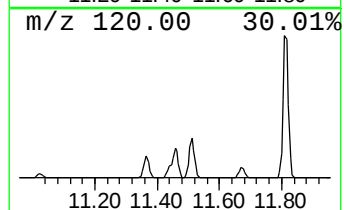
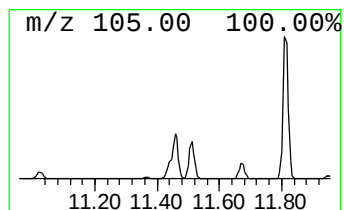
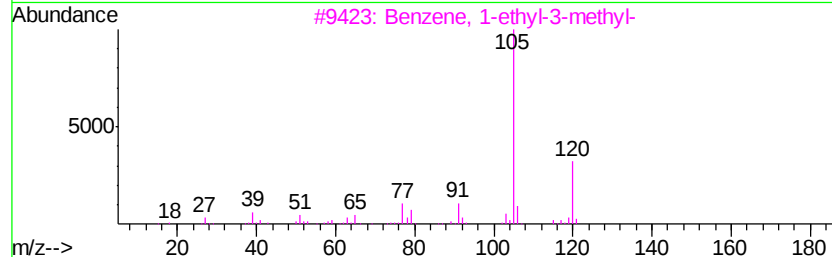
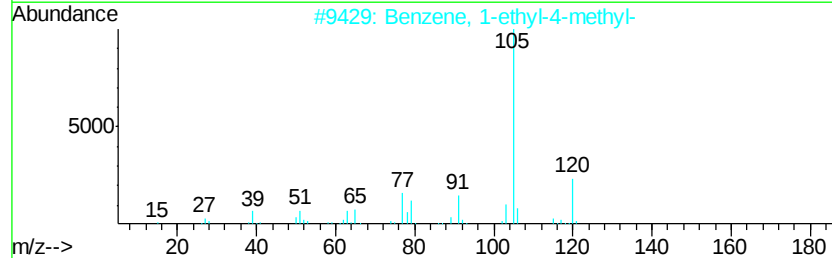
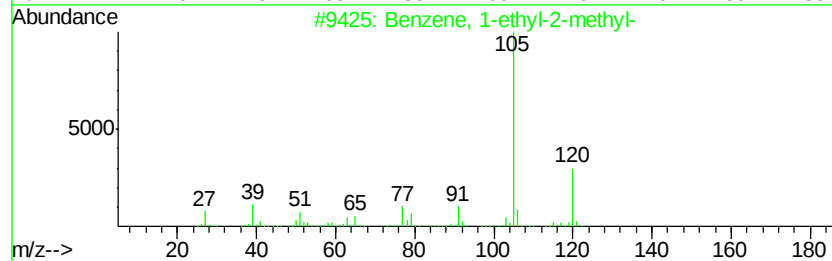
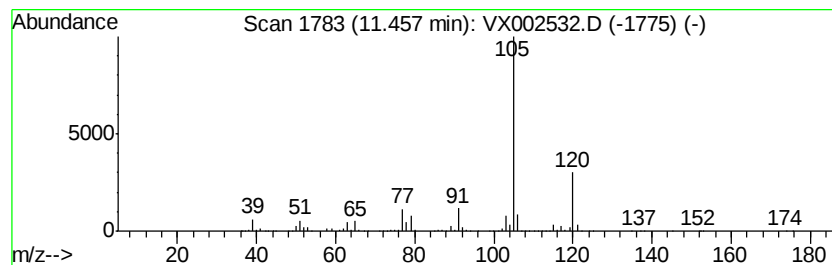
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.46	141.53 ug/l	4546510	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-4-methyl-	120	C9H12	000622-96-8	93
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002532.D
Acq On : 14 Jun 2018 02:41
Operator : JC/MD
Sample : J3383-14
Misc : 6.28g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-42R(I-2R)R

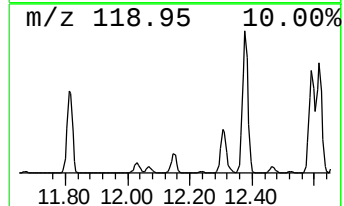
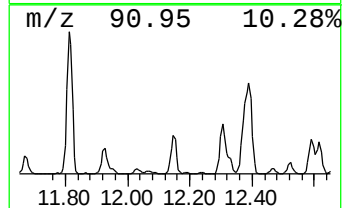
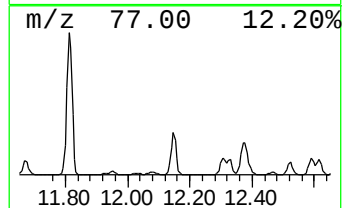
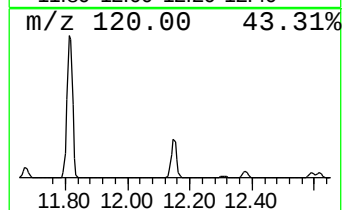
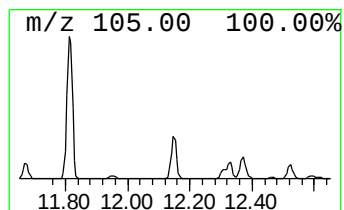
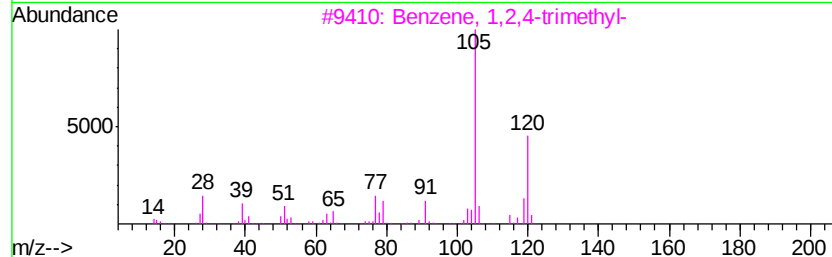
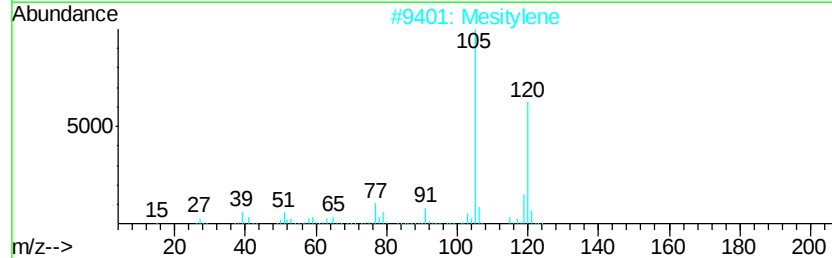
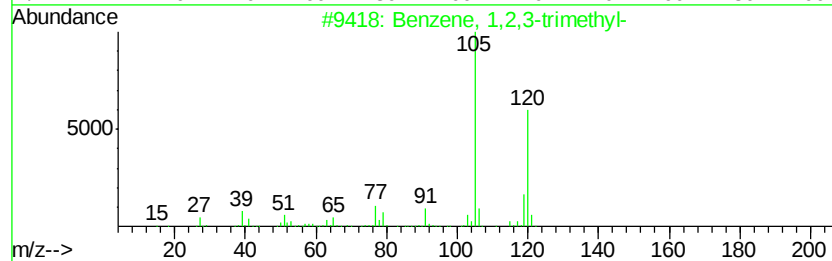
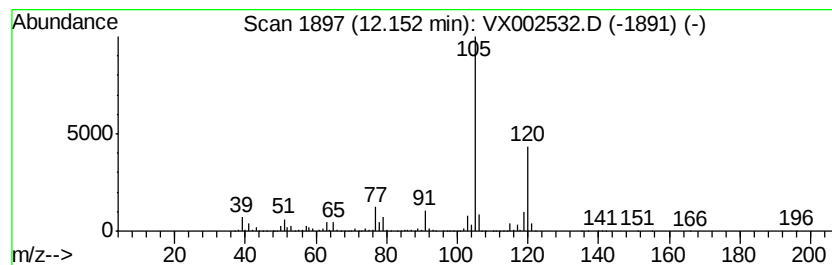
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	124.06 ug/l	3985390	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Mesitylene	120	C9H12	000108-67-8	94
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002532.D
Acq On : 14 Jun 2018 02:41
Operator : JC/MD
Sample : J3383-14
Misc : 6.28µ/5mL/100µL/5.0mL/MSVOA_X/MEOH
ALS Vial : 35 Sample Multiplier: 1

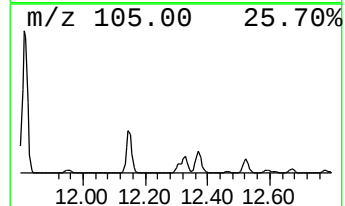
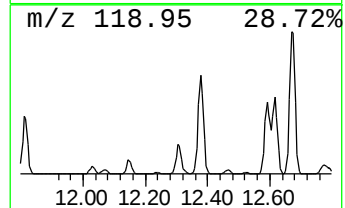
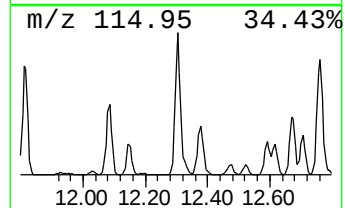
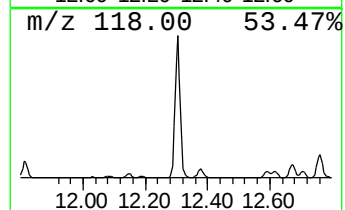
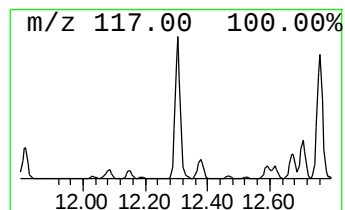
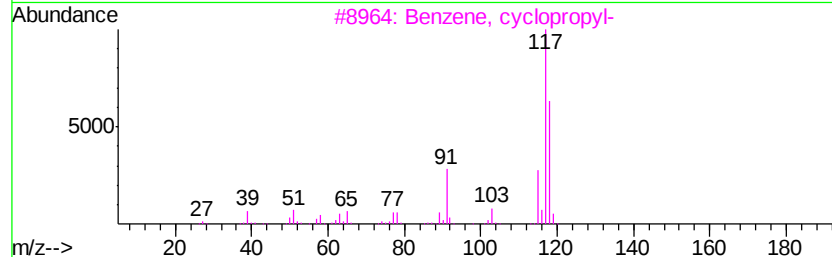
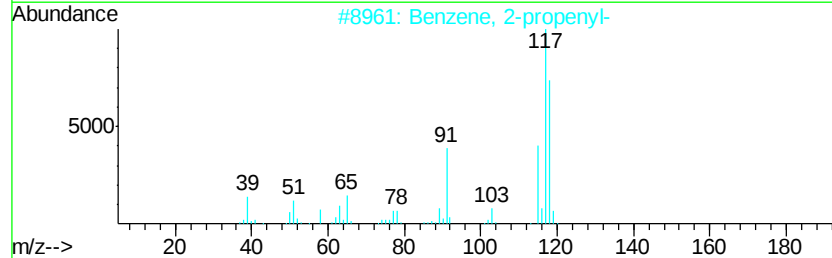
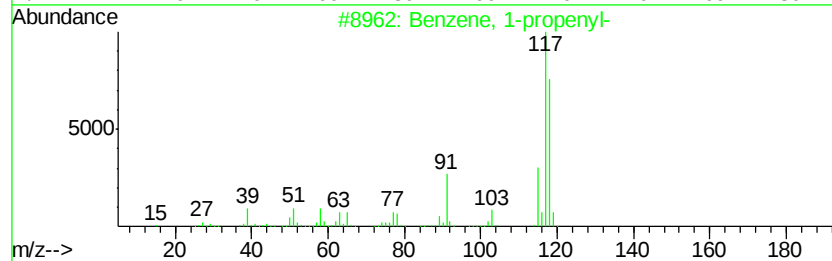
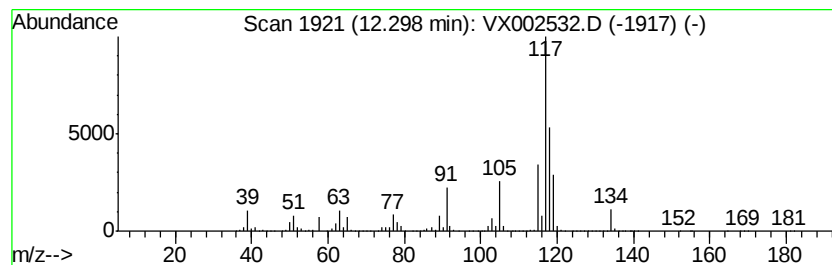
Instrument :
MSVOA_X
ClientSampleId :
B-42R(I-2R)R

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 6 Benzene, 1-propenyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	114.77 ug/l	3686750	1,4-Dichlorobenzene-d4	12.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-propenyl-	118 C9H10	000637-50-3 91
2		Benzene, 2-propenyl-	118 C9H10	000300-57-2 60
3		Benzene, cyclopropyl-	118 C9H10	000873-49-4 60
4		Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 60
5		Indane	118 C9H10	000496-11-7 55



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002532.D
Acq On : 14 Jun 2018 02:41
Operator : JC/MD
Sample : J3383-14
Misc : 6.28g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 35 Sample Multiplier: 1

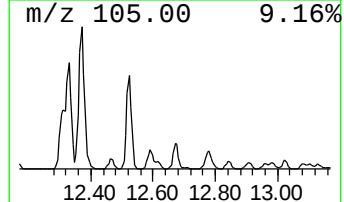
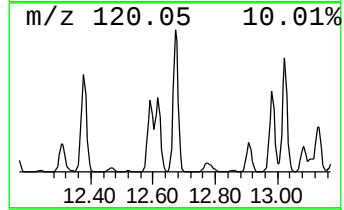
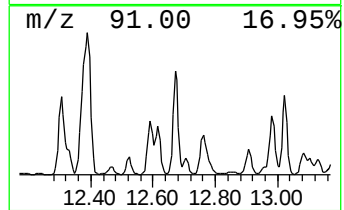
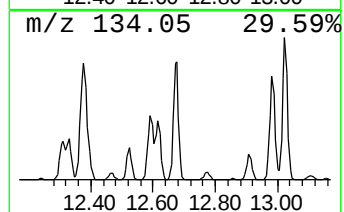
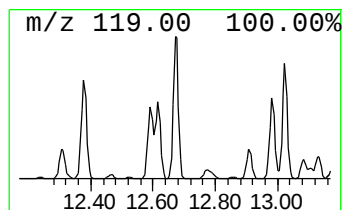
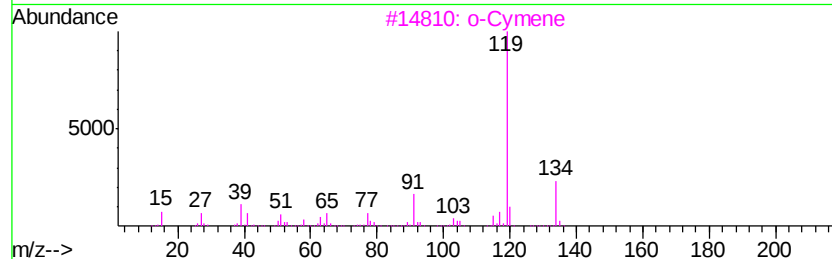
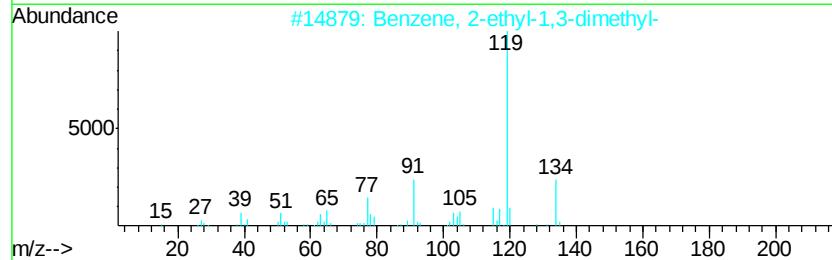
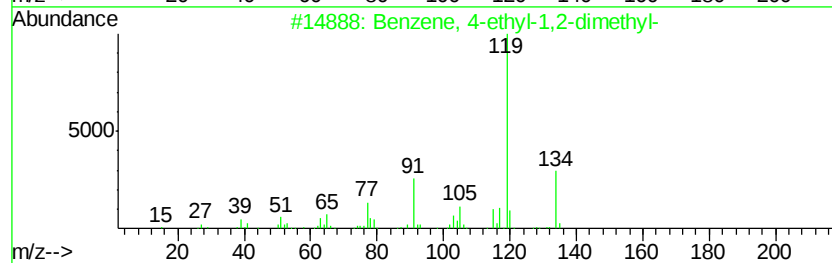
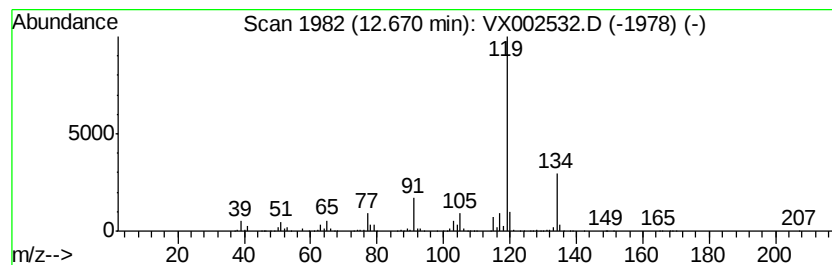
Instrument :
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ClientSampleId :
B-42R(I-2R)R

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 7 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	117.94 ug/l	3788880	1,4-Dichlorobenzene-d4	12.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5 95
2		Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4 95
3		o-Cymene	134 C10H14	000527-84-4 95
4		Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2 95
5		Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9 94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002532.D
Acq On : 14 Jun 2018 02:41
Operator : JC/MD
Sample : J3383-14
Misc : 6.28µ/5mL/100µL/5.0mL/MSVOA_X/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-42R(I-2R)R

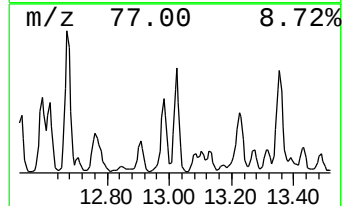
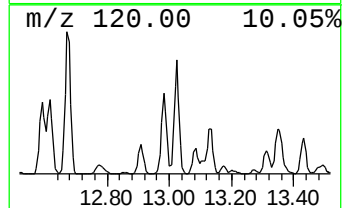
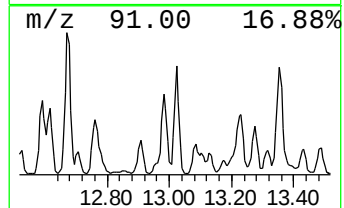
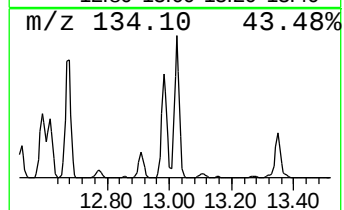
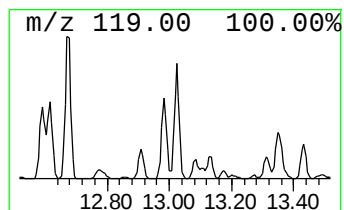
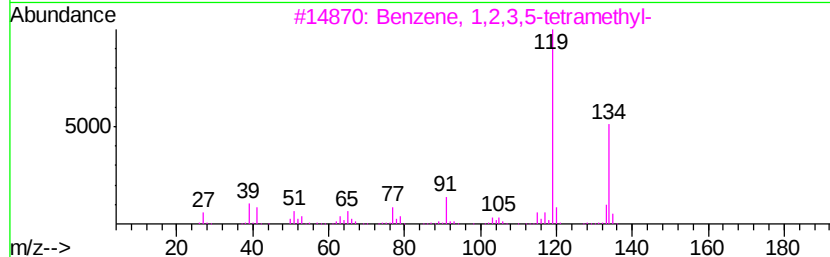
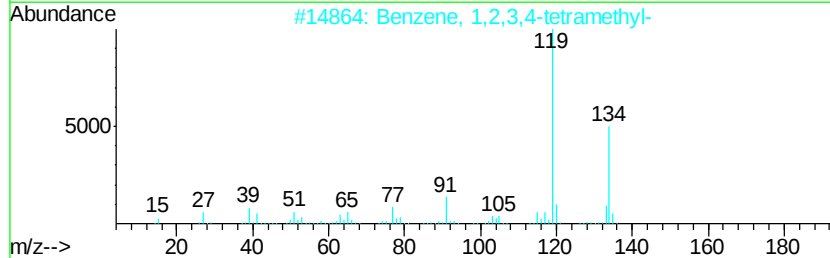
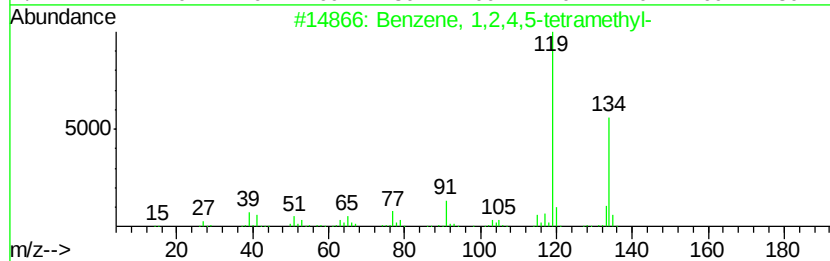
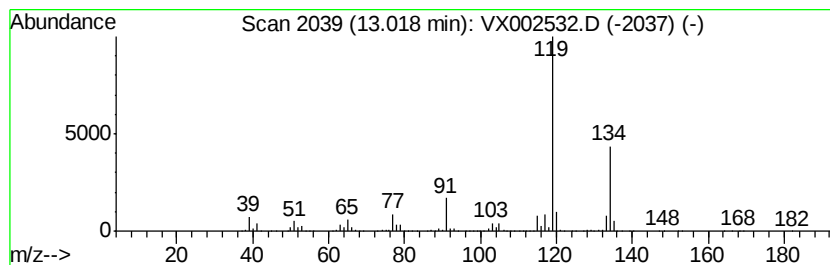
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	96.19 ug/l	3090060	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	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002532.D
Acq On : 14 Jun 2018 02:41
Operator : JC/MD
Sample : J3383-14
Misc : 6.28g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-42R(I-2R)R

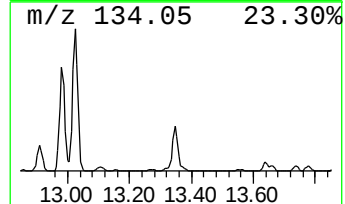
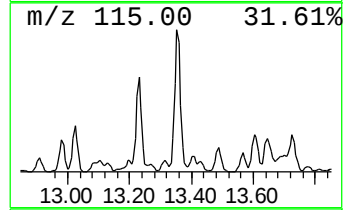
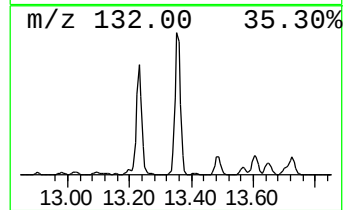
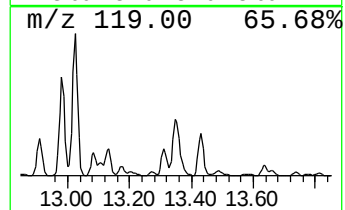
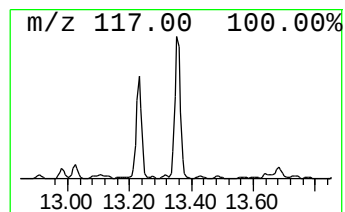
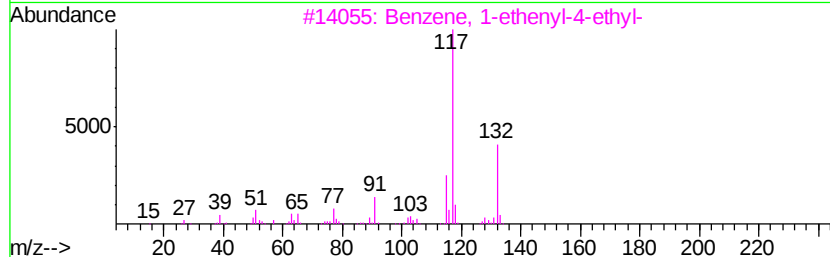
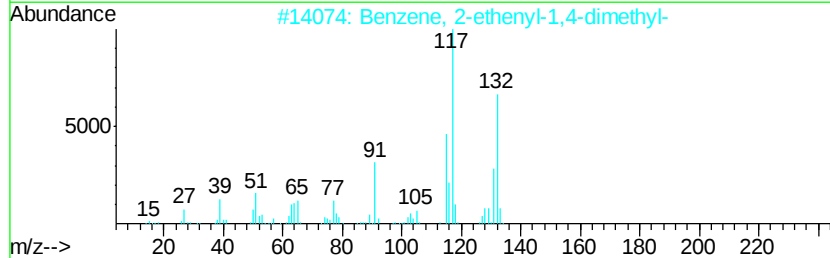
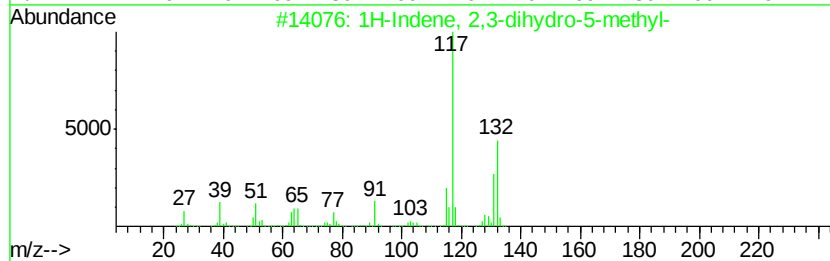
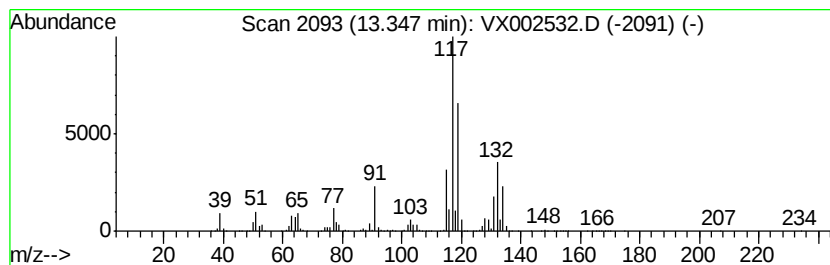
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Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	141.81 ug/l	4555700	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	95
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002532.D
Acq On : 14 Jun 2018 02:41
Operator : JC/MD
Sample : J3383-14
Misc : 6.28g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-42R(I-2R)R

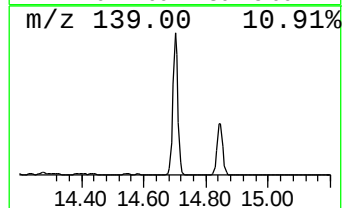
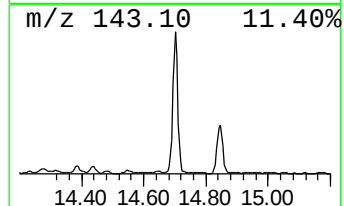
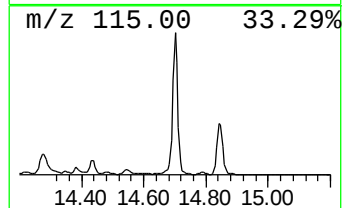
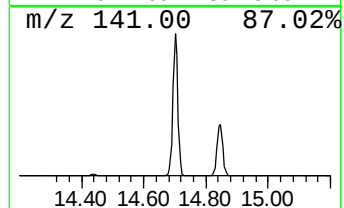
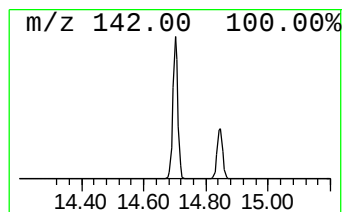
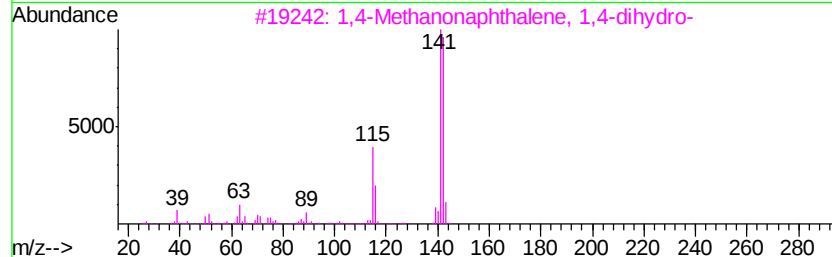
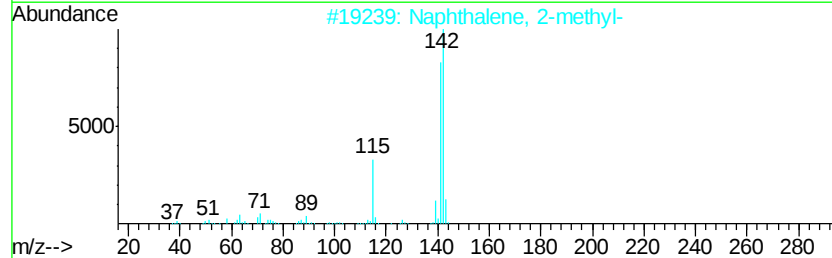
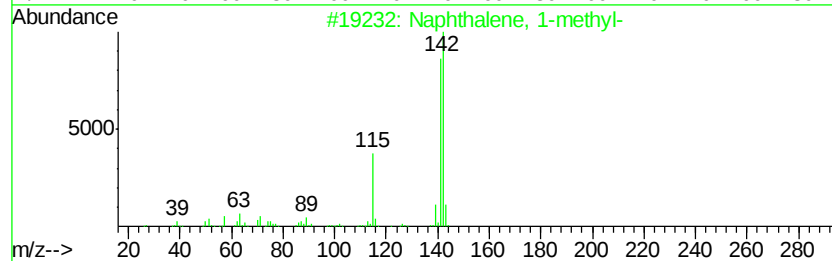
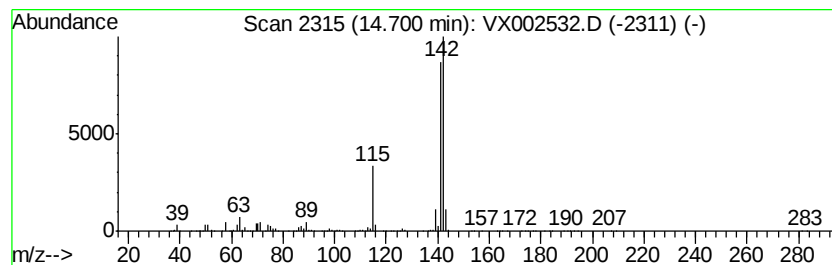
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.70	113.30 ug/l	3639660	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	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	86



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX061318\
Data File : VX002532.D
Acq On : 14 Jun 2018 02:41
Operator : JC/MD
Sample : J3383-14
Misc : 6.28g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-42R(I-2R)R

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
Heptane	6.70	156.1	ug/l	3745280	2	6.87	1200070	50.0
Heptane, 3-methyl-	8.50	106.4	ug/l	2350010	3	10.12	1104250	50.0
Octane	8.98	119.7	ug/l	2643880	3	10.12	1104250	50.0
Benzene, 1-ethyl-...	11.46	141.5	ug/l	4546510	4	12.09	1606220	50.0
Benzene, 1,2,3-tr...	12.15	124.1	ug/l	3985390	4	12.09	1606220	50.0
Benzene, 1-propenyl-	12.30	114.8	ug/l	3686750	4	12.09	1606220	50.0
Benzene, 4-ethyl-...	12.67	117.9	ug/l	3788880	4	12.09	1606220	50.0
Benzene, 1,2,4,5-...	13.02	96.2	ug/l	3090060	4	12.09	1606220	50.0
1H-Indene, 2,3-di...	13.35	141.8	ug/l	4555700	4	12.09	1606220	50.0
Naphthalene, 1-me...	14.70	113.3	ug/l	3639660	4	12.09	1606220	50.0