

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061318\
 Data File : VX002528.D
 Acq On : 14 Jun 2018 00:55
 Operator : JC/MD
 Sample : J3413-06
 Misc : 5.81g/5mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SW-1RR

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	1.746	186	190	201	rBV	2679477	2926635	9.17%	0.626%
2	1.935	216	221	236	rVB	4013721	5636129	17.66%	1.206%
3	2.215	262	267	281	rVB	711401	1120503	3.51%	0.240%
4	2.800	357	363	366	rBV	877752	1806135	5.66%	0.387%
5	2.849	366	371	383	rVB	4816307	10095521	31.63%	2.161%
6	3.136	408	418	440	rVB	3097047	7045147	22.08%	1.508%
7	3.489	465	476	491	rBV	5255809	11714757	36.71%	2.508%
8	3.782	518	524	534	rVB	772101	1836980	5.76%	0.393%
9	4.166	581	587	596	rVV	581387	1539544	4.82%	0.330%
10	4.276	596	605	612	rVV	457019	1362215	4.27%	0.292%
11	4.373	612	621	633	rVV	2447524	7080136	22.19%	1.516%
12	5.281	762	770	787	rVB	656511	2125058	6.66%	0.455%
13	5.592	803	821	830	rVV4	3555245	14878133	46.62%	3.185%
14	5.708	830	840	861	rVB3	1231137	5054025	15.84%	1.082%
15	5.915	861	874	890	rBV	3753383	11057562	34.65%	2.367%
16	6.245	917	928	938	rBV2	1007891	3661810	11.47%	0.784%
17	6.342	938	944	952	rVV2	1076098	2907362	9.11%	0.622%
18	6.446	952	961	973	rVB2	1284085	3723365	11.67%	0.797%
19	6.702	989	1003	1011	rBV	6607159	15646893	49.03%	3.349%
20	6.934	1022	1041	1055	rVB5	1165872	5097536	15.97%	1.091%
21	7.251	1083	1093	1100	rBV	599902	1392583	4.36%	0.298%
22	7.458	1114	1127	1138	rBV	6989006	16294243	51.06%	3.488%
23	7.574	1138	1146	1151	rBV2	947079	1962505	6.15%	0.420%
24	7.635	1151	1156	1165	rVB	1211164	2399153	7.52%	0.514%
25	7.732	1165	1172	1182	rVB	866920	1711163	5.36%	0.366%
26	7.842	1182	1190	1198	rVB2	991686	2081522	6.52%	0.446%
27	8.025	1215	1220	1237	rVB3	613492	2253032	7.06%	0.482%
28	8.171	1237	1244	1247	rBV2	629050	1291271	4.05%	0.276%
29	8.257	1254	1258	1262	rVB2	934422	1393292	4.37%	0.298%
30	8.342	1267	1272	1276	rBV	4611281	7837391	24.56%	1.678%
31	8.378	1276	1278	1285	rVB	2367986	3362605	10.54%	0.720%
32	8.507	1293	1299	1303	rBV	4531379	7605189	23.83%	1.628%
33	8.574	1307	1310	1319	rVB4	856159	1709105	5.36%	0.366%
34	8.671	1319	1326	1329	rBV3	1979948	3824085	11.98%	0.819%

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35	8.836	1348	1353	1357	rVV	941311	1867672	5.85%	0.400%
36	8.884	1357	1361	1363	rVV	687998	1162180	3.64%	0.249%
37	8.976	1370	1376	1383	rVV	7231919	10949656	34.31%	2.344%
38	9.043	1383	1387	1398	rVB4	1120955	2637936	8.27%	0.565%
39	9.226	1413	1417	1421	rBV2	1002158	1452888	4.55%	0.311%
40	9.445	1448	1453	1458	rBV	1076401	1548631	4.85%	0.331%
41	9.561	1467	1472	1479	rBV2	1731853	3270563	10.25%	0.700%
42	9.622	1479	1482	1486	rVB2	1115801	1482580	4.65%	0.317%
43	9.677	1486	1491	1495	rBV3	572837	1166764	3.66%	0.250%
44	9.823	1508	1515	1520	rBV2	783199	1565424	4.91%	0.335%
45	9.878	1520	1524	1529	rVV2	809555	1726415	5.41%	0.370%
46	9.927	1529	1532	1534	rVV2	785503	1166791	3.66%	0.250%
47	9.964	1534	1538	1543	rVB	4561859	6317362	19.80%	1.352%
48	10.067	1550	1555	1560	rVB	3837576	5173449	16.21%	1.107%
49	10.122	1560	1564	1567	rBV	834149	1135206	3.56%	0.243%
50	10.256	1581	1586	1591	rVV	9174807	12428344	38.94%	2.660%
51	10.366	1599	1604	1609	rVB	12511166	16448430	51.54%	3.521%
52	10.421	1609	1613	1624	rVB	5386281	7557111	23.68%	1.618%
53	10.646	1644	1650	1653	rBV2	1239617	1665187	5.22%	0.356%
54	10.726	1657	1663	1669	rBV2	729110	1328605	4.16%	0.284%
55	10.835	1674	1681	1685	rBV	1289666	2305702	7.22%	0.494%
56	10.915	1685	1694	1698	rBV2	1036135	1898362	5.95%	0.406%
57	11.024	1706	1712	1719	rBV	1213864	2115242	6.63%	0.453%
58	11.146	1726	1732	1734	rVV3	1849670	3213030	10.07%	0.688%
59	11.171	1734	1736	1742	rVB2	1917524	2687931	8.42%	0.575%
60	11.250	1742	1749	1752	rBV	1698908	2310455	7.24%	0.495%
61	11.366	1764	1768	1777	rVB	4565146	5544907	17.37%	1.187%
62	11.463	1777	1784	1787	rBV	7825373	9735983	30.51%	2.084%
63	11.512	1787	1792	1803	rVB2	8261738	14802974	46.39%	3.169%
64	11.677	1814	1819	1826	rBV	3084434	4444418	13.93%	0.951%
65	11.817	1837	1842	1850	rVB	26136407	31913270	100.00%	6.831%
66	12.085	1880	1886	1891	rBV3	1225332	2147931	6.73%	0.460%
67	12.152	1891	1897	1901	rBV	6656837	9099590	28.51%	1.948%
68	12.305	1917	1922	1929	rBV	6189245	8671006	27.17%	1.856%
69	12.378	1929	1934	1941	rVB4	8238323	12742462	39.93%	2.728%
70	12.481	1946	1951	1955	rVV2	2114038	2606374	8.17%	0.558%
71	12.524	1955	1958	1965	rVB	1869457	2383383	7.47%	0.510%

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72	12.591	1965	1969	1971	rBV	3688368	4529309	14.19%	0.970%
73	12.616	1971	1973	1979	rVV	3314013	3845513	12.05%	0.823%
74	12.676	1979	1983	1986	rVV	6662745	7641123	23.94%	1.636%
75	12.707	1986	1988	1992	rVB	1076047	1122413	3.52%	0.240%
76	12.762	1992	1997	2005	rVB2	3217355	5260576	16.48%	1.126%
77	12.908	2018	2021	2025	rVB2	1340114	1572645	4.93%	0.337%
78	12.981	2025	2033	2037	rBV	3743060	5585777	17.50%	1.196%
79	13.024	2037	2040	2046	rVB	5485351	6553975	20.54%	1.403%
80	13.085	2046	2050	2051	rBV	1215722	1327603	4.16%	0.284%
81	13.109	2051	2054	2056	rVV3	1308309	1787754	5.60%	0.383%
82	13.134	2056	2058	2061	rVB	1225535	1141134	3.58%	0.244%
83	13.231	2067	2074	2078	rBV3	3659210	5250268	16.45%	1.124%
84	13.274	2078	2081	2084	rVB2	1258676	1424192	4.46%	0.305%
85	13.323	2084	2089	2091	rBV3	1926475	2848122	8.92%	0.610%
86	13.359	2091	2095	2100	rVV2	6025989	8301752	26.01%	1.777%
87	13.432	2104	2107	2112	rVB	1459965	1732512	5.43%	0.371%
88	13.487	2112	2116	2124	rVB2	1282065	2148784	6.73%	0.460%
89	13.567	2124	2129	2132	rBV2	1265607	1754557	5.50%	0.376%
90	13.609	2132	2136	2139	rBV	2001777	2604954	8.16%	0.558%
91	13.646	2139	2142	2147	rVB4	1983741	3155167	9.89%	0.675%
92	13.731	2152	2156	2161	rVV2	1832005	2932992	9.19%	0.628%
93	13.841	2170	2174	2179	rVB	3422677	4150233	13.00%	0.888%
94	13.926	2184	2188	2192	rVB3	765257	1199567	3.76%	0.257%
95	13.987	2192	2198	2203	rBV4	1324607	2559951	8.02%	0.548%
96	14.140	2219	2223	2227	rBV2	1855272	2265991	7.10%	0.485%
97	14.280	2241	2246	2255	rBV2	1698948	3036354	9.51%	0.650%
98	14.438	2268	2272	2276	rVB	1114203	1343789	4.21%	0.288%
99	14.701	2306	2315	2325	rVB	4001006	5979549	18.74%	1.280%
100	14.847	2335	2339	2348	rVB	1454430	2011300	6.30%	0.431%

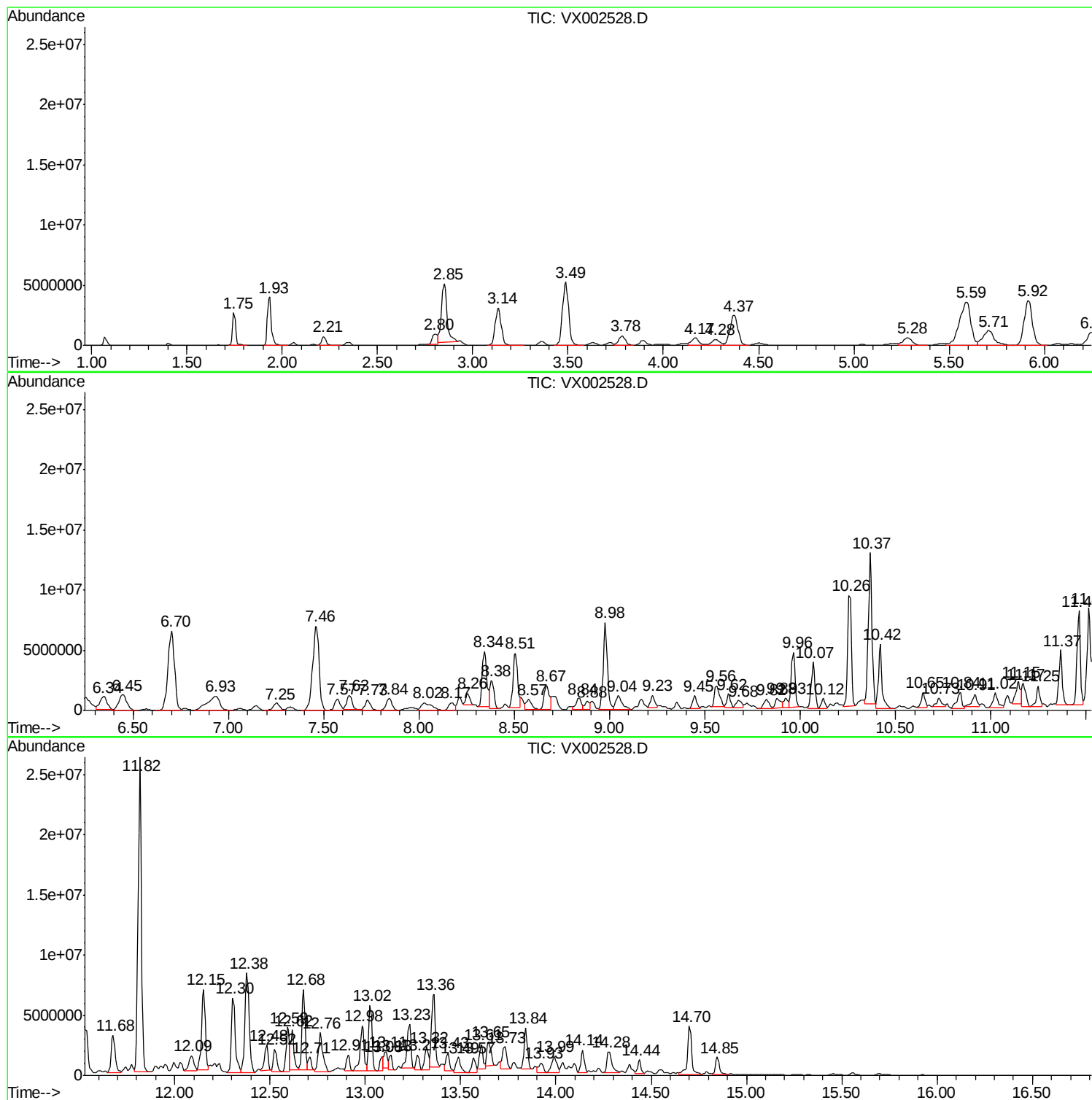
Sum of corrected areas: 467174655

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Instrument :
MSVOA_X
ClientSampleId :
SW-1RR

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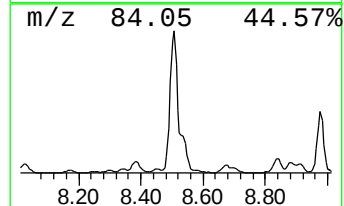
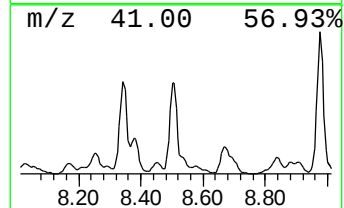
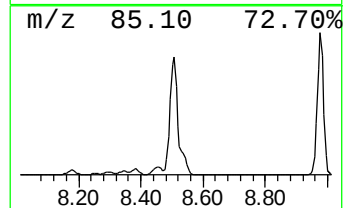
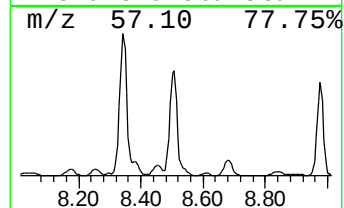
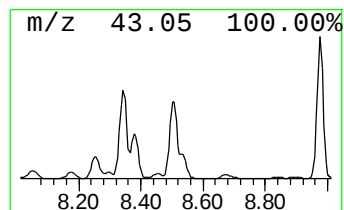
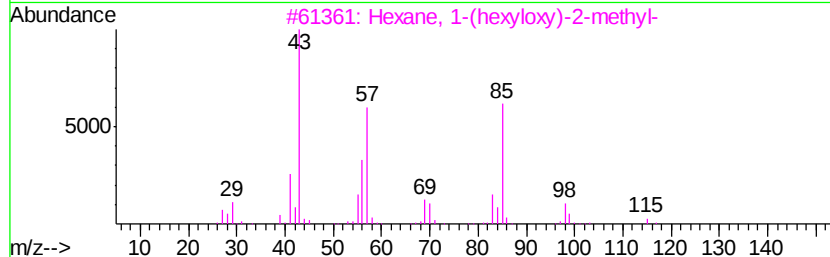
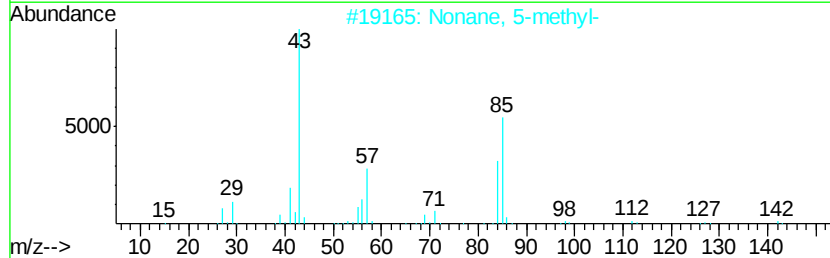
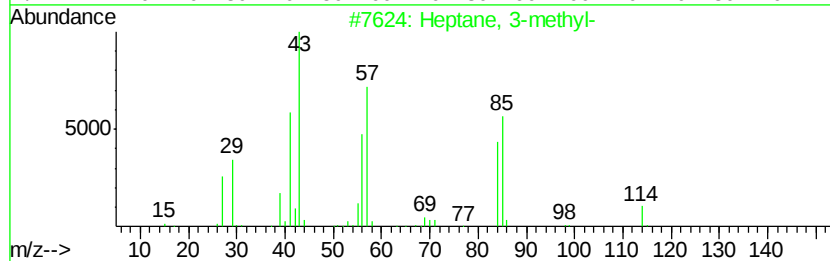
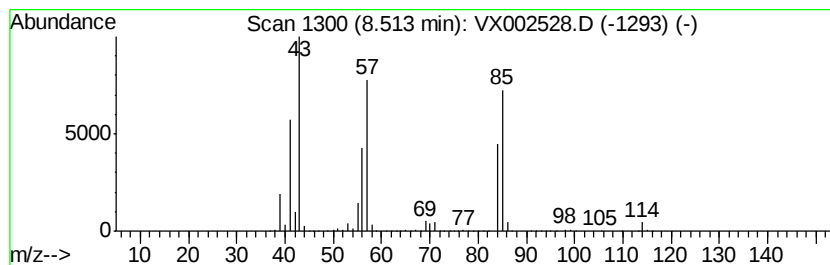
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

Peak Number 1 Heptane, 3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.51	334.97 ug/l	7605190	Chlorobenzene-d5	10.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane, 3-methyl-	114	C8H18	000589-81-1	94
2		Nonane, 5-methyl-	142	C10H22	015869-85-9	64
3		Hexane, 1-(hexyloxy)-2-methyl-	200	C13H28O	074421-17-3	59
4		2,2-Dimethylpropionic acid, 4-me...	186	C11H22O2	150588-62-8	59
5		Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	59



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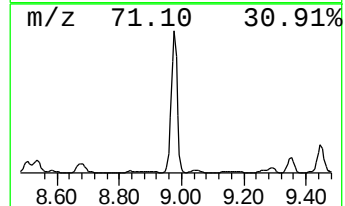
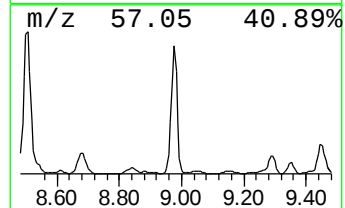
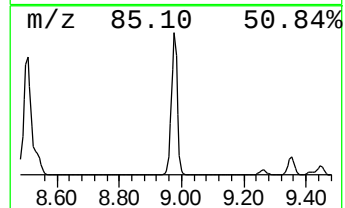
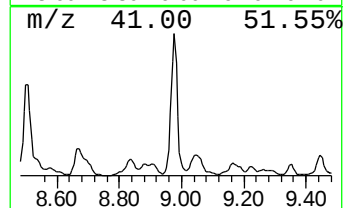
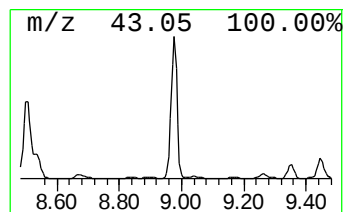
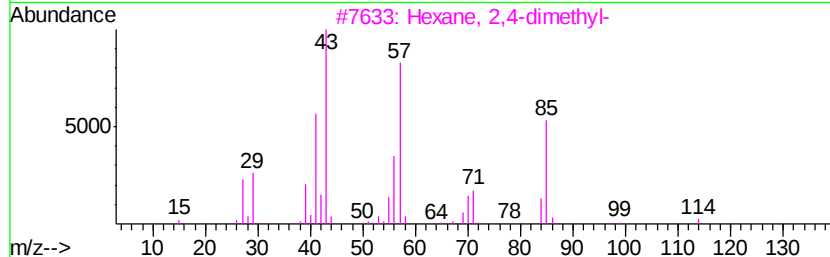
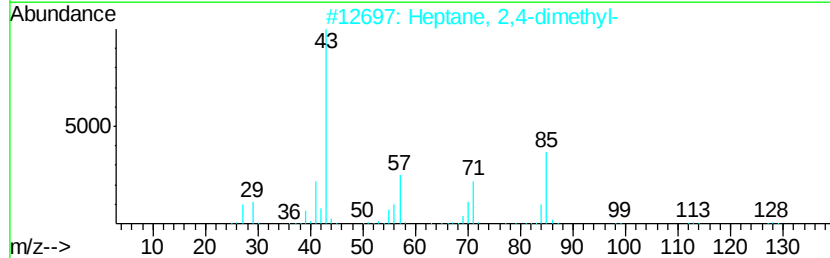
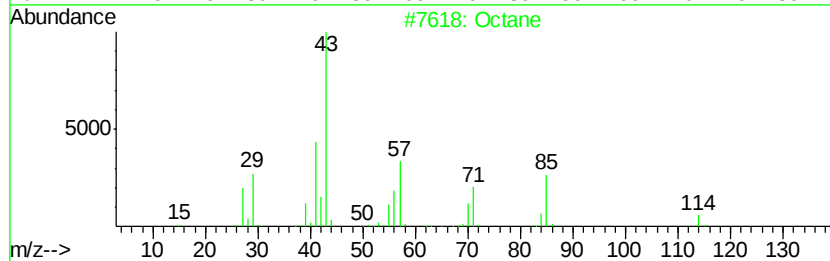
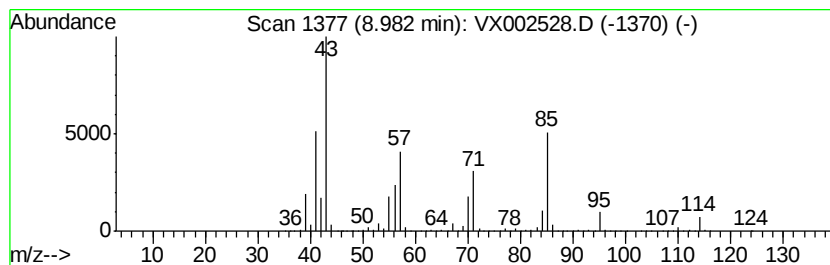
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Octane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.98	482.28 ug/l	10949700	Chlorobenzene-d5	10.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane		114	C8H18	000111-65-9	94
2	Heptane, 2,4-dimethyl-		128	C9H20	002213-23-2	64
3	Hexane, 2,4-dimethyl-		114	C8H18	000589-43-5	64
4	Undecane, 2,4-dimethyl-		184	C13H28	017312-80-0	59
5	Ethanol, 2-(hexyloxy)-		146	C8H18O2	000112-25-4	53



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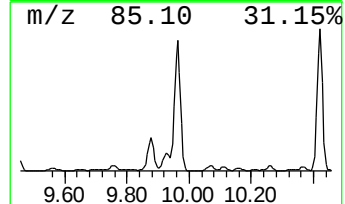
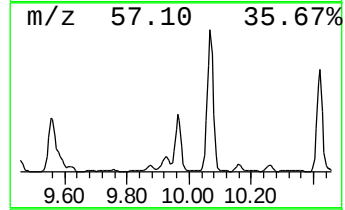
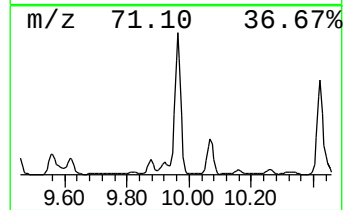
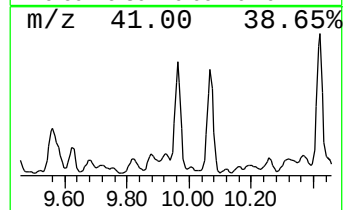
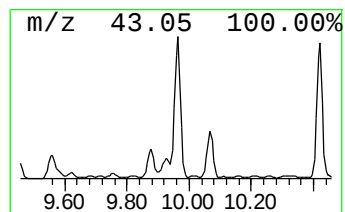
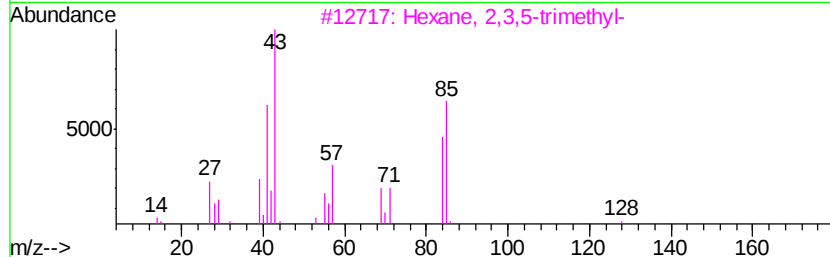
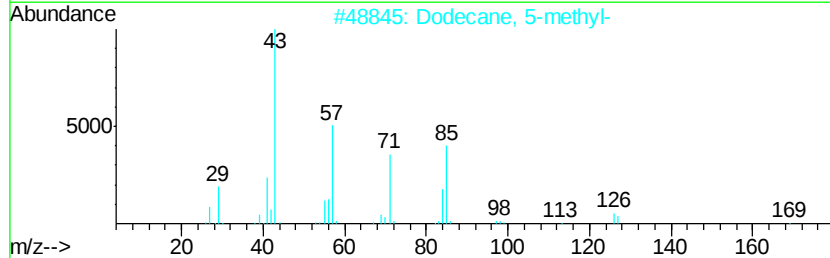
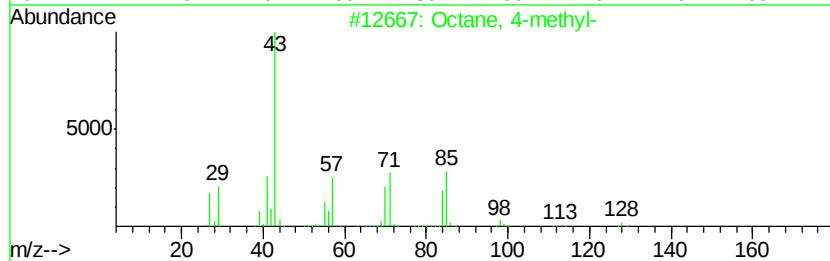
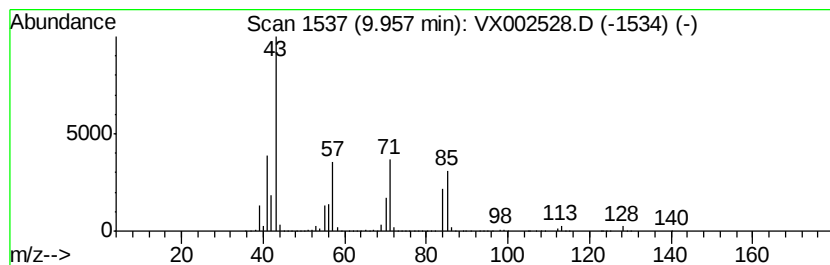
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Quant Title : SW846 8260

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

Peak Number 3 Octane, 4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.96	278.25 ug/l	6317360	Chlorobenzene-d5	10.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 4-methyl-	128	C9H20	002216-34-4	72
2		Dodecane, 5-methyl-	184	C13H28	017453-93-9	59
3		Hexane, 2,3,5-trimethyl-	128	C9H20	001069-53-0	53
4		Sulfurous acid, 2-propyl tetra...	320	C17H36O3S	1000309-12-5	53
5		Decane	142	C10H22	000124-18-5	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002528.D
Acq On : 14 Jun 2018 00:55
Operator : JC/MD
Sample : J3413-06
Misc : 5.81µ/5mL/100µL/5.0mL/MSVOA_X/MEOH
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SW-1RR

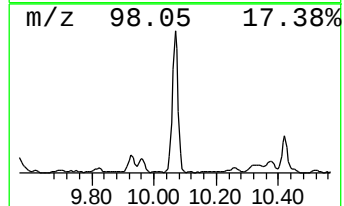
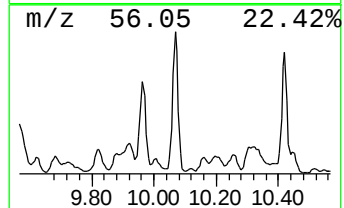
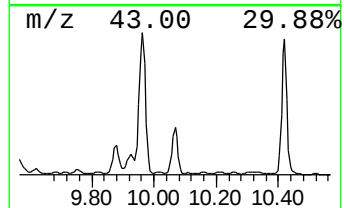
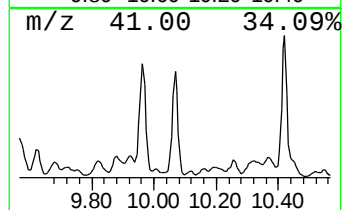
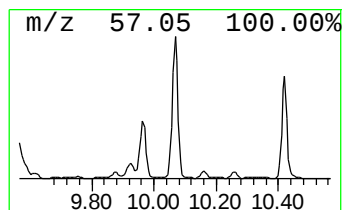
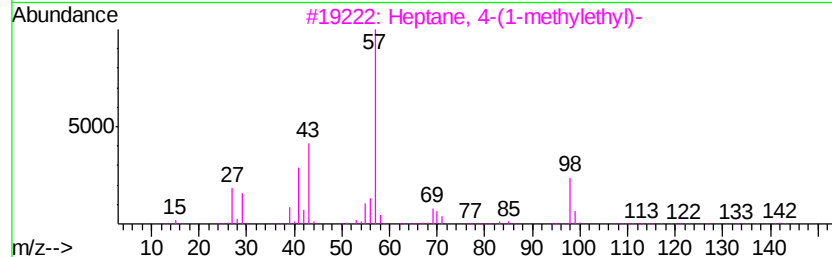
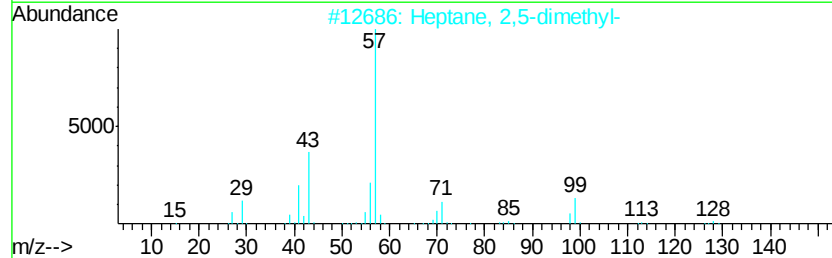
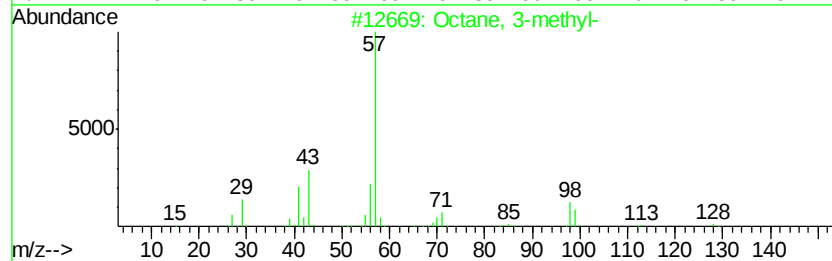
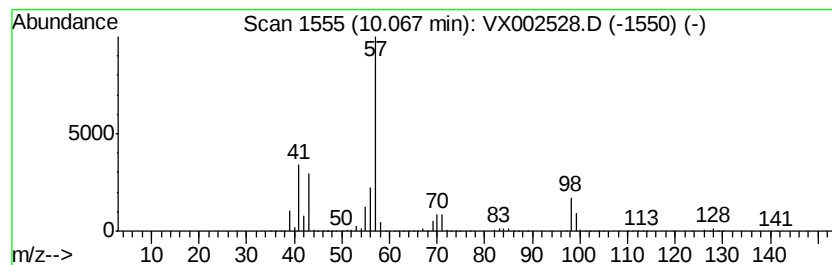
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Quant Title : SW846 8260

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

Peak Number 4 Octane, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.07	227.86 ug/l	5173450	Chlorobenzene-d5	10.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 3-methyl-	128	C9H20	002216-33-3	87
2		Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	64
3		Heptane, 4-(1-methylethyl)-	142	C10H22	052896-87-4	59
4		Heptane, 4-propyl-	142	C10H22	003178-29-8	59
5		Oxalic acid, isobutyl heptyl ester	244	C13H24O4	1000309-37-2	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002528.D
Acq On : 14 Jun 2018 00:55
Operator : JC/MD
Sample : J3413-06
Misc : 5.81µ/5mL/100µL/5.0mL/MSVOA_X/MEOH
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SW-1RR

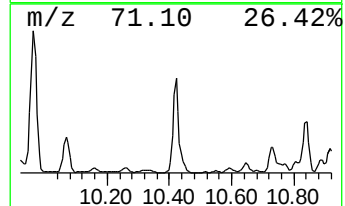
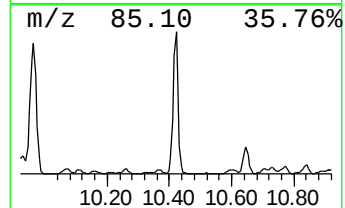
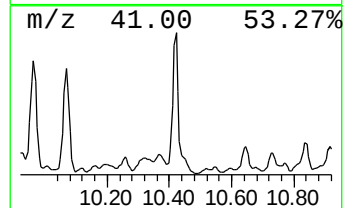
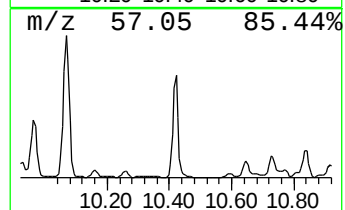
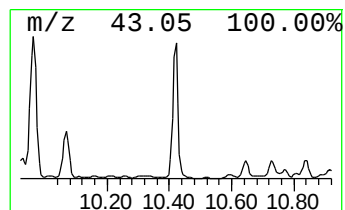
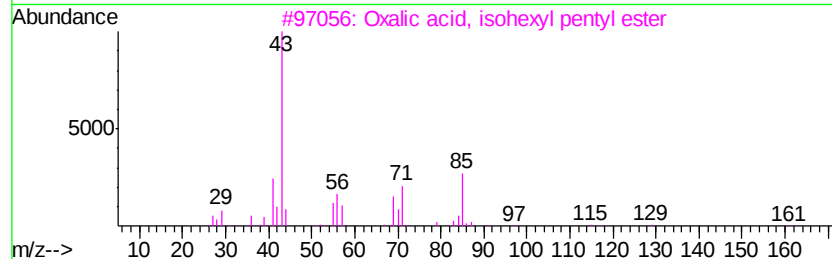
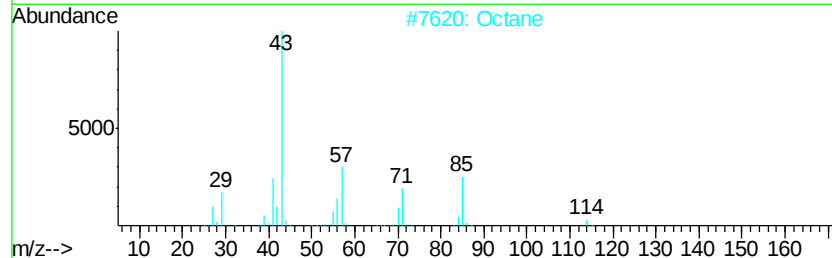
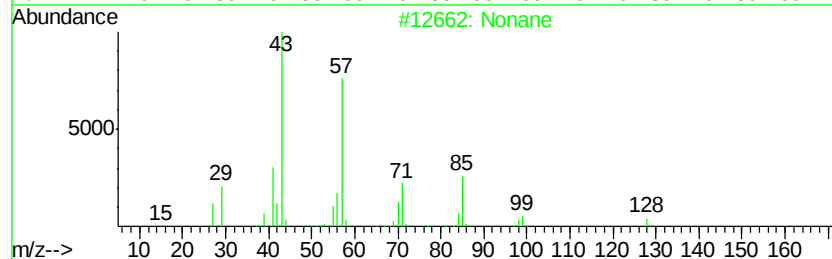
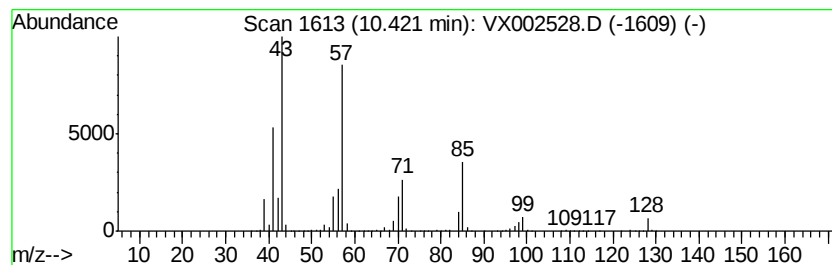
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Quant Title : SW846 8260

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

Peak Number 5 Nonane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.42	332.85 ug/l	7557110	Chlorobenzene-d5	10.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane		128	C9H20	000111-84-2	97
2	Octane		114	C8H18	000111-65-9	64
3	Oxalic acid, isohexyl pentyl ester		244	C13H24O4	1000309-32-8	64
4	Octane, 3,5-dimethyl-		142	C10H22	015869-93-9	53
5	Oxalic acid, isobutyl pentyl ester		216	C11H20O4	1000309-37-0	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002528.D
Acq On : 14 Jun 2018 00:55
Operator : JC/MD
Sample : J3413-06
Misc : 5.81µ/5mL/100µL/5.0mL/MSVOA_X/MEOH
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
SW-1RR

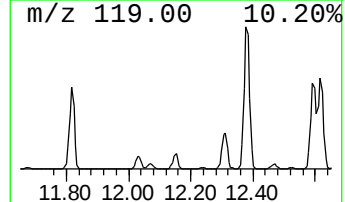
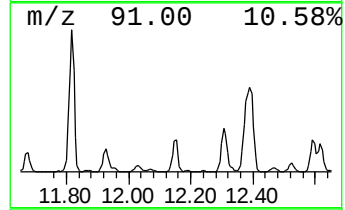
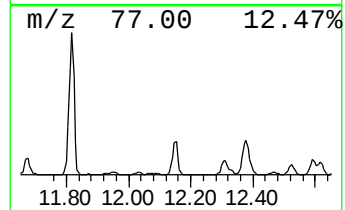
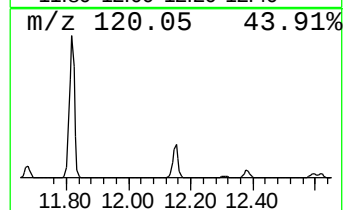
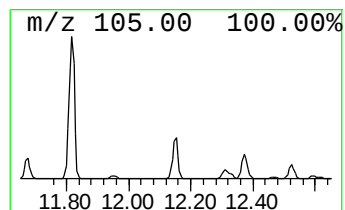
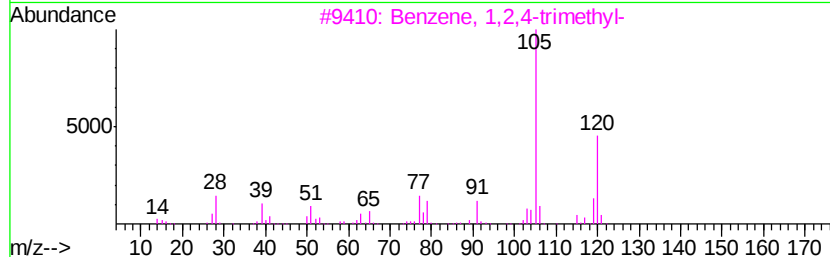
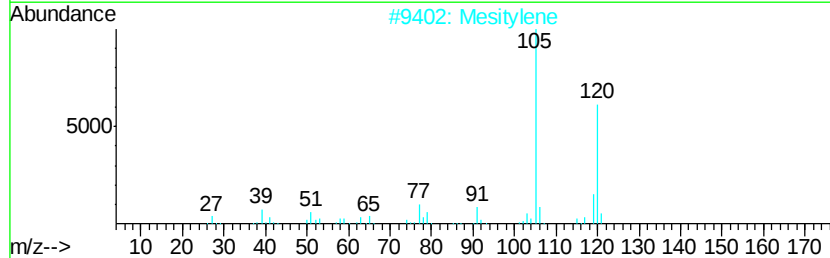
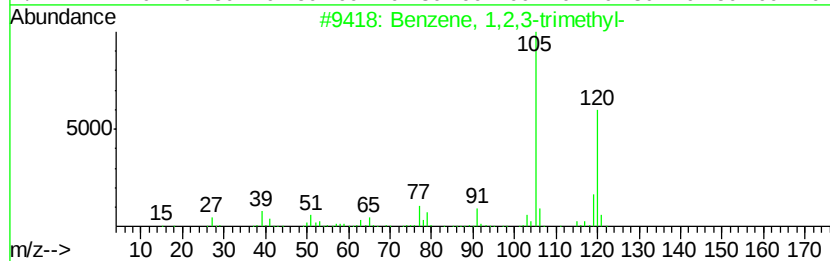
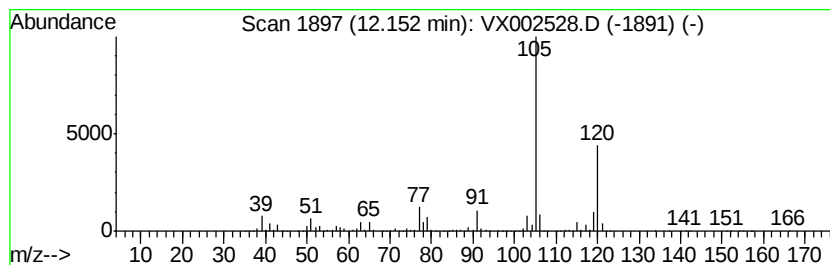
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	211.82 ug/l	9099590	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	93
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
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 : VX002528.D
Acq On : 14 Jun 2018 00:55
Operator : JC/MD
Sample : J3413-06
Misc : 5.81µ/5mL/100µL/5.0mL/MSVOA_X/MEOH
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SW-1RR

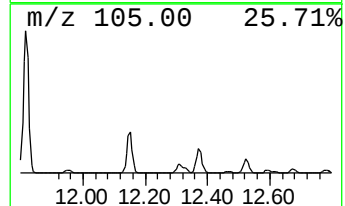
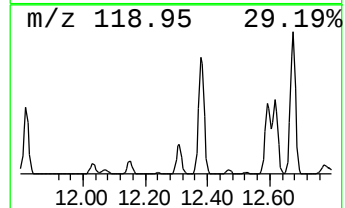
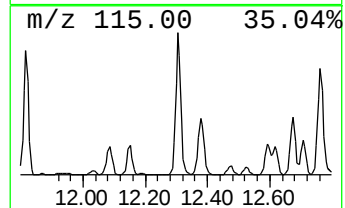
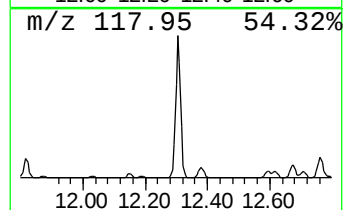
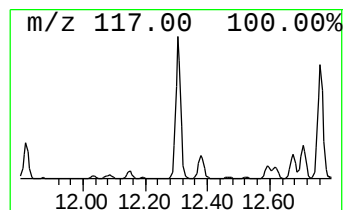
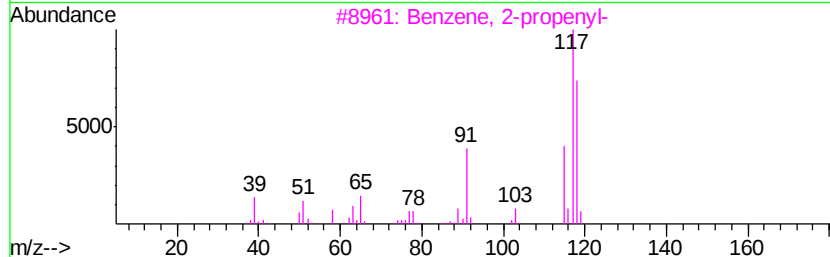
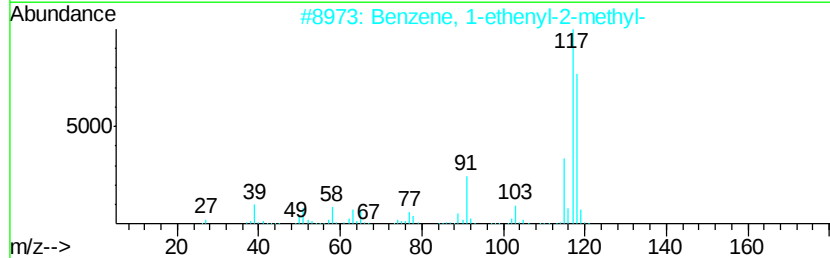
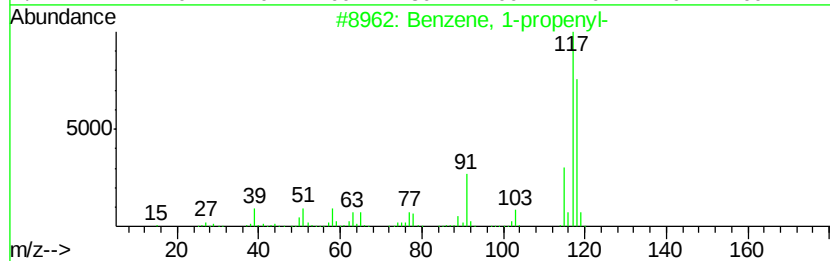
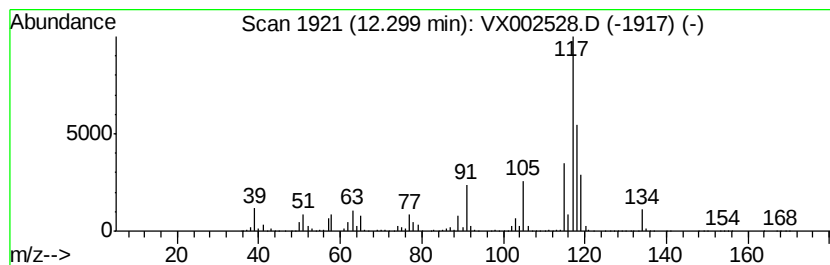
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-propenyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	201.85 ug/l	8671010	1,4-Dichlorobenzene-d4	12.09

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propenyl-	118	C9H10	000637-50-3	92	
2	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	91	
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	64	
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64	
5	Indane	118	C9H10	000496-11-7	60	



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002528.D
Acq On : 14 Jun 2018 00:55
Operator : JC/MD
Sample : J3413-06
Misc : 5.81µ/5mL/100µL/5.0mL/MSVOA_X/MEOH
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SW-1RR

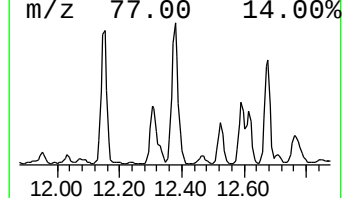
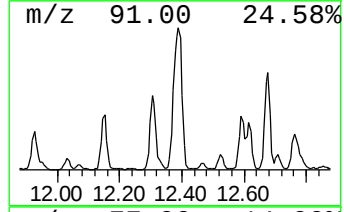
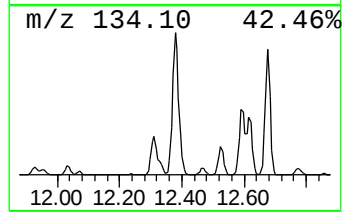
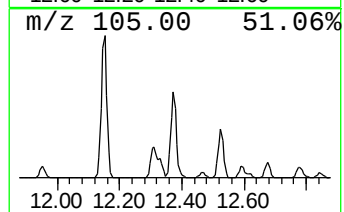
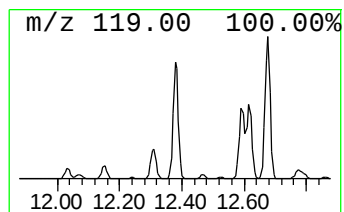
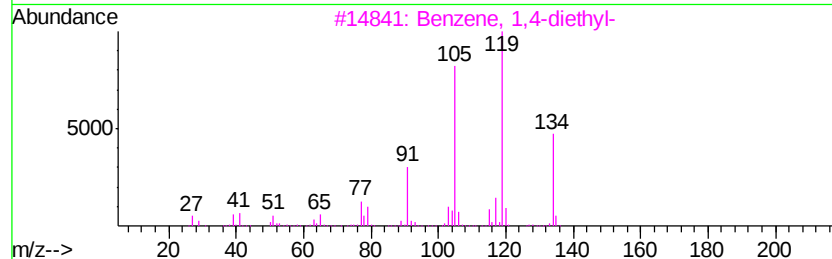
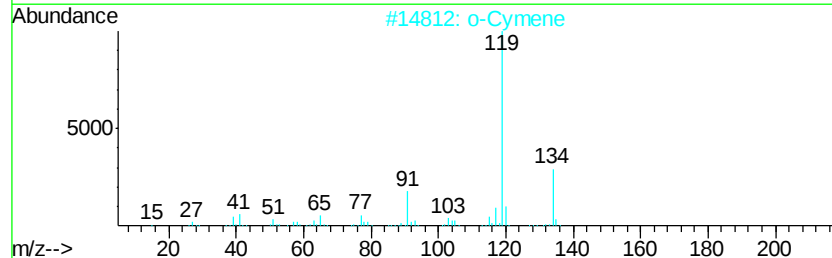
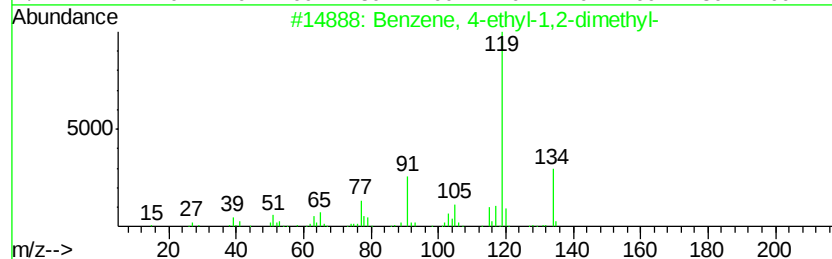
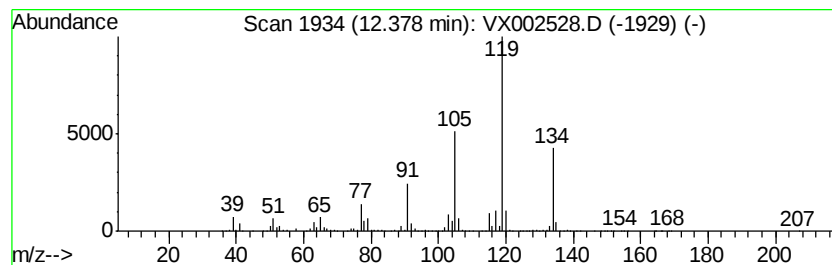
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.38	296.62 ug/l	12742500	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	94
2		o-Cymene	134	C10H14	000527-84-4	94
3		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
4		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061318\
Data File : VX002528.D
Acq On : 14 Jun 2018 00:55
Operator : JC/MD
Sample : J3413-06
Misc : 5.81µ/5mL/100µL/5.0mL/MSVOA_X/MEOH
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SW-1RR

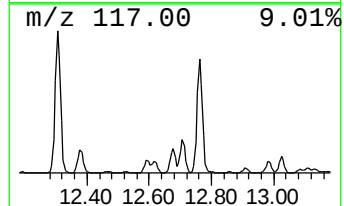
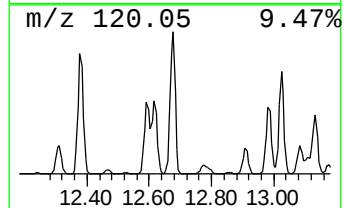
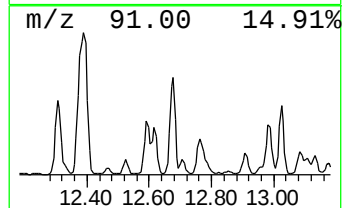
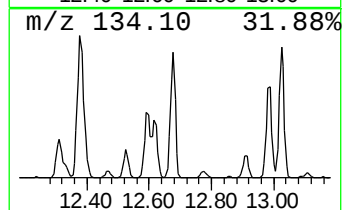
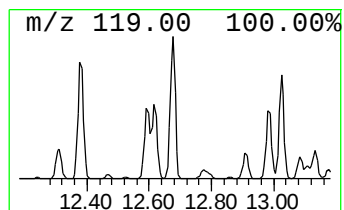
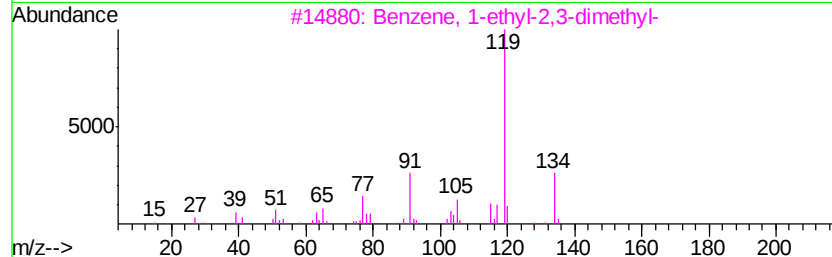
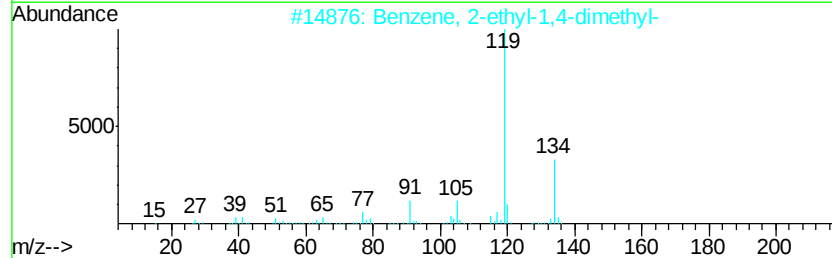
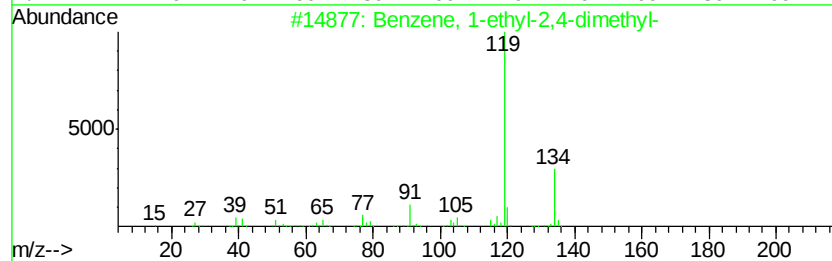
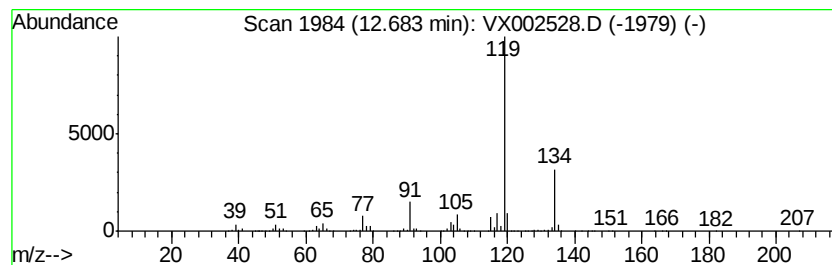
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Quant Title : SW846 8260

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	177.87 ug/l	7641120	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, 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		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 : VX002528.D
Acq On : 14 Jun 2018 00:55
Operator : JC/MD
Sample : J3413-06
Misc : 5.81µ/5mL/100µL/5.0mL/MSVOA_X/MEOH
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SW-1RR

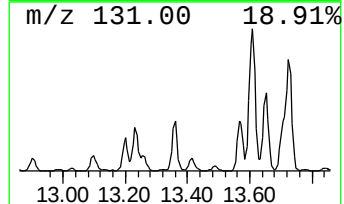
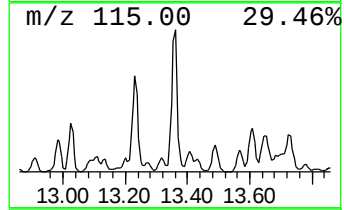
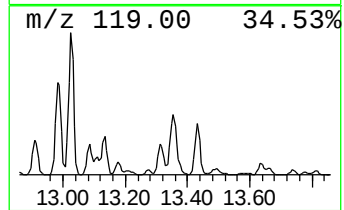
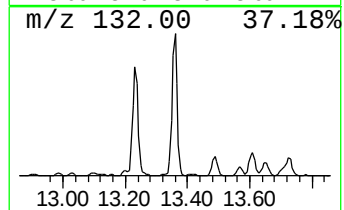
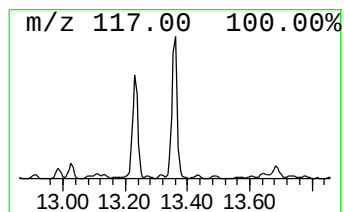
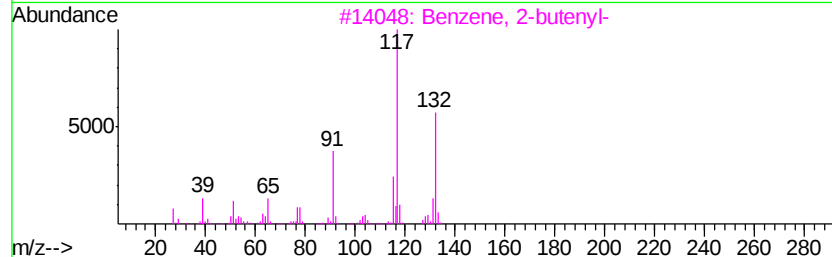
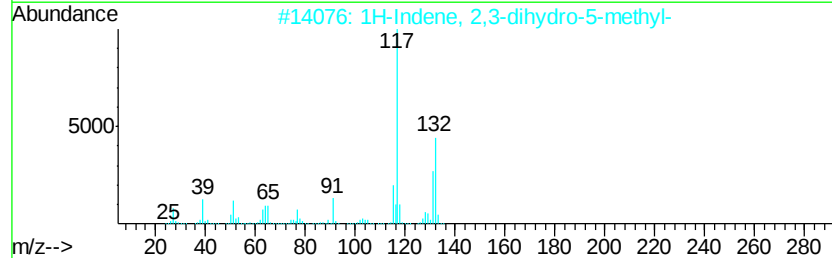
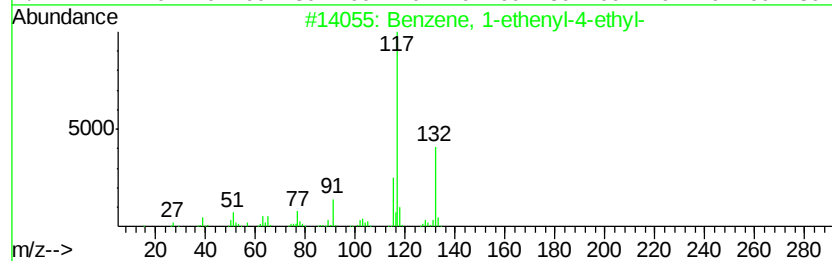
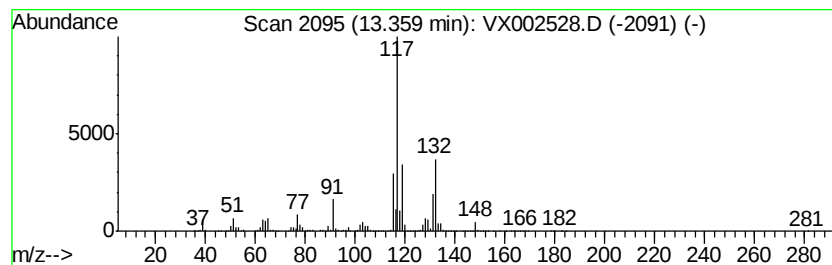
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.36	193.25 ug/l	8301750	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
3		Benzene, 2-butenyl-	132	C10H12	001560-06-1	81
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
5		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX061318\
Data File : VX002528.D
Acq On : 14 Jun 2018 00:55
Operator : JC/MD
Sample : J3413-06
Misc : 5.81g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SW-1RR

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, 3-methyl-	8.51	335.0	ug/l	7605190	3	10.12	1135210	50.0
Octane	8.98	482.3	ug/l	10949700	3	10.12	1135210	50.0
Octane, 4-methyl-	9.96	278.3	ug/l	6317360	3	10.12	1135210	50.0
Octane, 3-methyl-	10.07	227.9	ug/l	5173450	3	10.12	1135210	50.0
Nonane	10.42	332.9	ug/l	7557110	3	10.12	1135210	50.0
Benzene, 1,2,3-tr...	12.15	211.8	ug/l	9099590	4	12.09	2147930	50.0
Benzene, 1-propenyl-	12.30	201.8	ug/l	8671010	4	12.09	2147930	50.0
Benzene, 4-ethyl-...	12.38	296.6	ug/l	12742500	4	12.09	2147930	50.0
Benzene, 1-ethyl-...	12.68	177.9	ug/l	7641120	4	12.09	2147930	50.0
Benzene, 1-etheny...	13.36	193.3	ug/l	8301750	4	12.09	2147930	50.0