

Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
 Data File : VH057091.D
 Acq On : 21 Sep 2015 13:46
 Operator : SY/FY
 Sample : G3727-05
 Misc : 5.72/5mL/100uL/5mL/MSVOA_H/MEOH
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_H
 ClientSampleId :
 MMB7B

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.001

Stop Thrs : 0

Filtering: 9

Min Area: 250 Area counts

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 1

Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.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	1.951	140	145	158	rBV	919574	3087360	4.80%	0.218%
2	2.329	177	181	192	rBV	1529871	3341705	5.20%	0.236%
3	2.876	219	233	250	rVB3	790564	3563988	5.55%	0.252%
4	3.695	301	311	319	rBV	3002825	11867364	18.47%	0.839%
5	3.958	330	336	343	rVV	3753175	12403146	19.30%	0.877%
6	4.233	355	362	366	rBV	1666242	6034026	9.39%	0.426%
7	4.328	366	371	376	rVB2	1324008	4233826	6.59%	0.299%
8	4.433	376	381	383	rBV3	797108	2488972	3.87%	0.176%
9	4.633	394	400	411	rVV	8706017	27207536	42.34%	1.923%
10	4.813	411	417	422	rVV3	778304	3194356	4.97%	0.226%
11	5.404	463	473	490	rBV3	6268400	39491667	61.45%	2.791%
12	5.710	495	502	515	rVB3	2578121	11156915	17.36%	0.789%
13	5.953	515	525	534	rBV3	1967043	10957633	17.05%	0.774%
14	6.184	538	547	554	rBV2	2308858	11015237	17.14%	0.778%
15	6.468	569	574	576	rBV	735758	2288005	3.56%	0.162%
16	6.563	577	583	586	rBV	9998681	29101627	45.28%	2.057%
17	6.774	597	603	616	rBV	6172767	19767710	30.76%	1.397%
18	6.986	616	623	624	rBV	5336339	15268420	23.76%	1.079%
19	7.208	639	644	647	rBV	2269831	5922641	9.22%	0.419%
20	7.272	647	650	653	rVV	2944599	8522452	13.26%	0.602%
21	7.335	653	656	661	rVV	3014799	9874084	15.36%	0.698%
22	7.451	661	667	671	rVV	19719473	55278046	86.01%	3.907%
23	7.546	671	676	680	rVV2	4138179	17337687	26.98%	1.225%
24	7.610	680	682	688	rVV	2515397	7262664	11.30%	0.513%
25	7.727	688	693	697	rVV	2305414	9442151	14.69%	0.667%
26	7.790	697	699	703	rVV2	1545571	3598801	5.60%	0.254%
27	7.864	703	706	711	rVV3	1197799	3714198	5.78%	0.262%
28	8.087	721	727	735	rVV3	1720332	6658413	10.36%	0.471%
29	8.236	735	741	747	rVV	7393292	22191310	34.53%	1.568%
30	8.342	747	751	754	rVV2	3768588	11724679	18.24%	0.829%
31	8.405	754	757	768	rVV2	10614343	33130621	51.55%	2.341%
32	8.583	768	774	788	rVV	8652829	38229843	59.49%	2.702%
33	8.794	789	794	800	rVV	3227588	10741306	16.71%	0.759%
34	8.921	800	806	811	rVV3	6880592	22061057	34.33%	1.559%

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35	9.006	811	814	816	rVV3	2431493	6610587	10.29%	0.467%
36	9.068	816	820	826	rVV2	9690601	28484161	44.32%	2.013%
37	9.173	826	830	832	rVV3	2280138	6393600	9.95%	0.452%
38	9.246	832	837	845	rVV	10617044	30741312	47.83%	2.173%
39	9.372	845	849	853	rVV2	1790623	5234274	8.14%	0.370%
40	9.487	855	860	866	rVV2	2424388	9542552	14.85%	0.674%
41	9.647	867	875	878	rVV2	4538018	18132323	28.21%	1.281%
42	9.710	878	881	885	rVV2	8426332	22977930	35.75%	1.624%
43	9.774	885	887	889	rVV	3738325	8113536	12.62%	0.573%
44	9.859	889	895	899	rVV	23550515	64266169	100.00%	4.542%
45	9.922	899	901	905	rVV2	1871723	4607027	7.17%	0.326%
46	10.016	906	910	915	rVV3	1881048	5559250	8.65%	0.393%
47	10.176	920	925	933	rVV3	5963710	21688813	33.75%	1.533%
48	10.303	933	937	943	rVV2	2584884	9519366	14.81%	0.673%
49	10.461	946	952	959	rVV4	6668825	25138600	39.12%	1.777%
50	10.557	959	961	963	rVV2	1493326	3356379	5.22%	0.237%
51	10.619	963	967	975	rVV2	16764867	49523408	77.06%	3.500%
52	10.732	975	978	980	rVV	2958157	7570534	11.78%	0.535%
53	10.784	980	983	985	rVV	5013920	12263224	19.08%	0.867%
54	10.837	985	988	990	rVV2	2846759	7818906	12.17%	0.553%
55	10.901	990	994	999	rVV4	3306335	11861546	18.46%	0.838%
56	10.995	999	1003	1006	rVV3	5358319	15361157	23.90%	1.086%
57	11.069	1006	1010	1012	rVV2	8959671	21834590	33.98%	1.543%
58	11.111	1012	1014	1018	rVV	8026559	16858086	26.23%	1.191%
59	11.215	1018	1024	1029	rVV2	10253875	30885484	48.06%	2.183%
60	11.375	1033	1039	1043	rVV6	3717303	17391482	27.06%	1.229%
61	11.438	1043	1045	1048	rVV3	3498881	8195113	12.75%	0.579%
62	11.500	1048	1051	1054	rVV	9268734	20522112	31.93%	1.450%
63	11.618	1058	1062	1065	rVV	27811629	55963826	87.08%	3.955%
64	11.671	1065	1067	1069	rVV3	4138210	8975132	13.97%	0.634%
65	11.724	1069	1072	1074	rVV2	3502233	8874436	13.81%	0.627%
66	11.767	1074	1076	1079	rVV3	3803725	8359236	13.01%	0.591%
67	11.830	1079	1082	1084	rVV2	2538421	6310007	9.82%	0.446%
68	11.894	1084	1088	1090	rVV2	3844348	11821744	18.39%	0.835%
69	11.946	1090	1093	1097	rVV	11745013	25083365	39.03%	1.773%
70	12.018	1097	1100	1104	rVV3	2930661	10035721	15.62%	0.709%
71	12.103	1104	1108	1109	rVV2	4292634	10878610	16.93%	0.769%

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Integration Parameters: rteint.p

Integrator: RTE
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72	12.145	1109	1112	1116	rVV2	10059297	25628569	39.88%	1.811%
73	12.229	1116	1120	1122	rVV5	4176716	12014207	18.69%	0.849%
74	12.271	1122	1124	1126	rVV2	2892022	6208596	9.66%	0.439%
75	12.324	1126	1129	1131	rVV	5664169	12786912	19.90%	0.904%
76	12.367	1131	1133	1134	rVV	6066967	9600412	14.94%	0.678%
77	12.398	1134	1136	1141	rVV	6373876	16107424	25.06%	1.138%
78	12.483	1141	1144	1146	rVV3	5040581	10306400	16.04%	0.728%
79	12.535	1146	1149	1153	rVV2	4488694	12898116	20.07%	0.912%
80	12.599	1153	1155	1160	rVV5	2322958	8629107	13.43%	0.610%
81	12.706	1160	1165	1167	rVV4	5796105	18146510	28.24%	1.282%
82	12.759	1167	1170	1177	rVV2	22465474	48392773	75.30%	3.420%
83	12.876	1178	1181	1186	rVV4	3217929	13132854	20.44%	0.928%
84	12.951	1186	1188	1191	rVV3	2979845	6974387	10.85%	0.493%
85	13.054	1191	1198	1199	rVV4	3292635	13342926	20.76%	0.943%
86	13.086	1199	1201	1203	rVV	3209485	5998091	9.33%	0.424%
87	13.128	1203	1205	1209	rVV3	2808086	8739612	13.60%	0.618%
88	13.181	1209	1210	1212	rVV	2738109	4330875	6.74%	0.306%
89	13.223	1212	1214	1217	rVV2	4322440	9414503	14.65%	0.665%
90	13.276	1217	1219	1220	rVV2	2702944	4677957	7.28%	0.331%
91	13.308	1220	1222	1225	rVV2	2829044	6365995	9.91%	0.450%
92	13.351	1225	1226	1231	rVV4	2158274	5706777	8.88%	0.403%
93	13.426	1231	1233	1235	rVV2	1826195	3448817	5.37%	0.244%
94	13.468	1235	1237	1238	rVV2	1540730	2598642	4.04%	0.184%
95	13.511	1238	1241	1243	rVV2	1988499	4873681	7.58%	0.344%
96	13.564	1243	1246	1249	rVV3	1392221	4480748	6.97%	0.317%
97	13.639	1250	1253	1257	rVV	5582002	10801400	16.81%	0.763%
98	13.777	1264	1266	1270	rVV3	2386338	5402226	8.41%	0.382%
99	13.851	1270	1273	1275	rVV2	1168496	2422025	3.77%	0.171%
100	14.084	1291	1295	1299	rVV5	858324	2610462	4.06%	0.184%

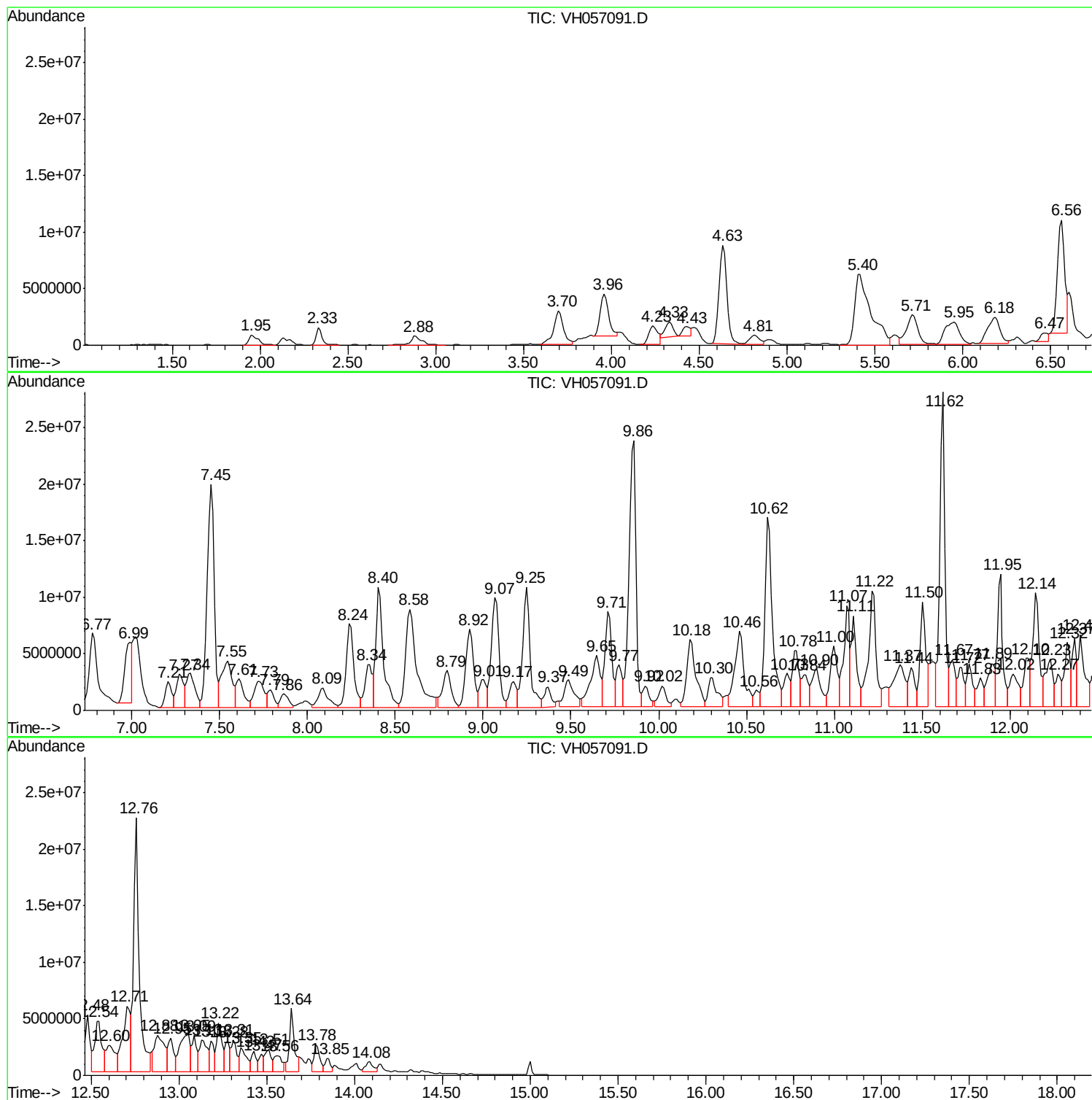
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Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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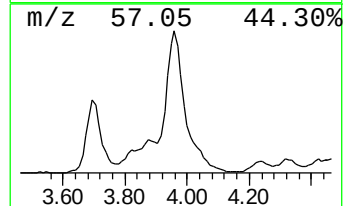
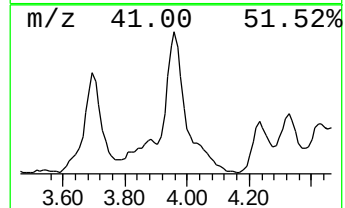
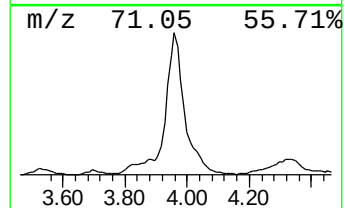
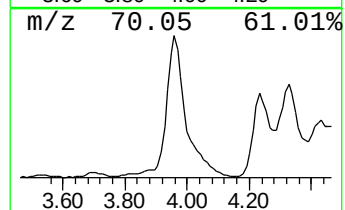
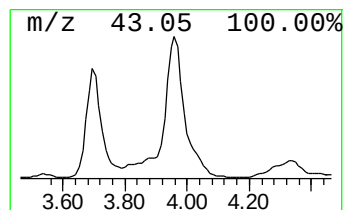
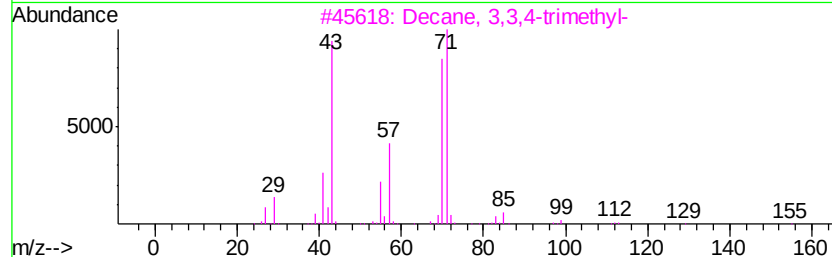
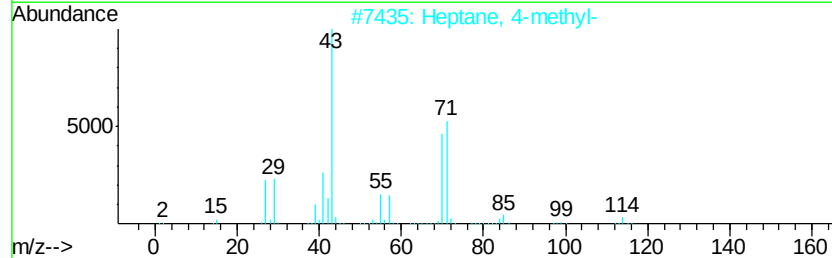
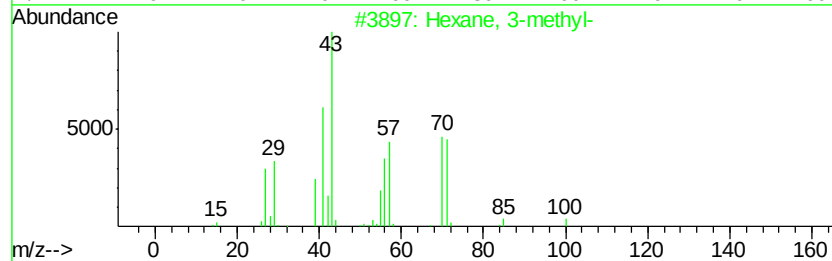
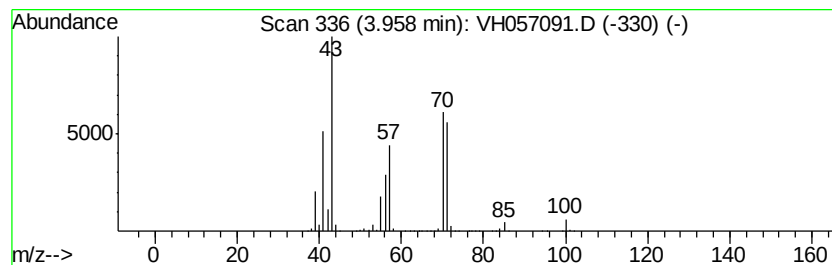
Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Hexane, 3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.96	194.14 ug/l	12403100	Pentafluorobenzene	4.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexane, 3-methyl-	100	C7H16	000589-34-4	76
2		Heptane, 4-methyl-	114	C8H18	000589-53-7	64
3		Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	64
4		Pentane, 3-ethyl-	100	C7H16	000617-78-7	64
5		Octane, 4,5-dimethyl-	142	C10H22	015869-96-2	59



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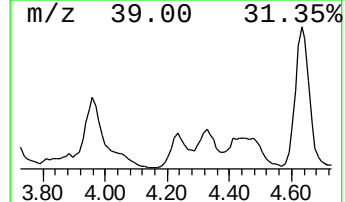
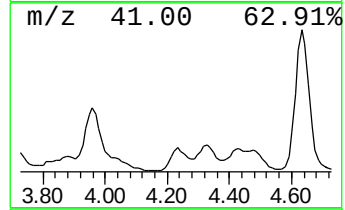
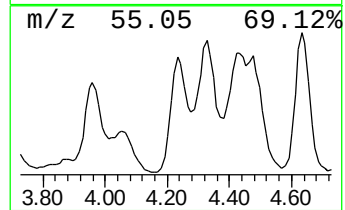
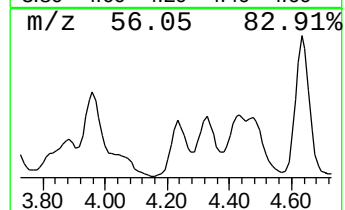
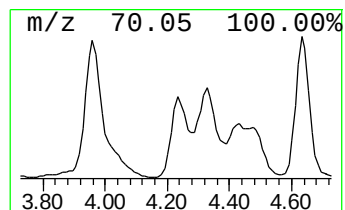
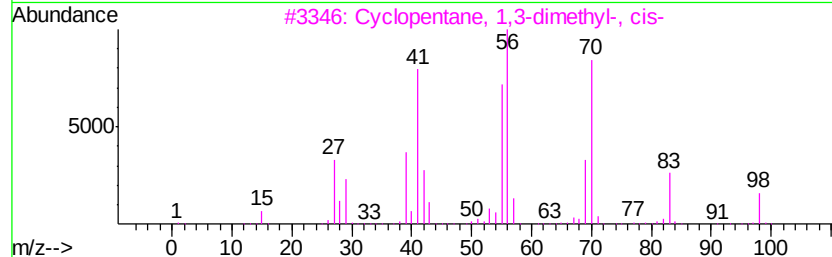
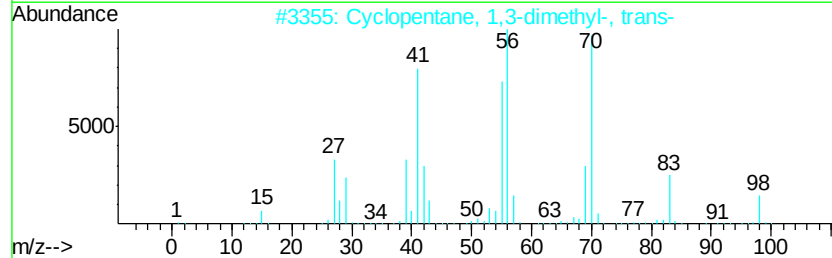
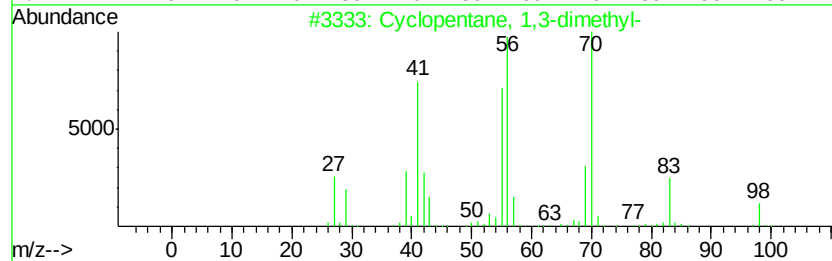
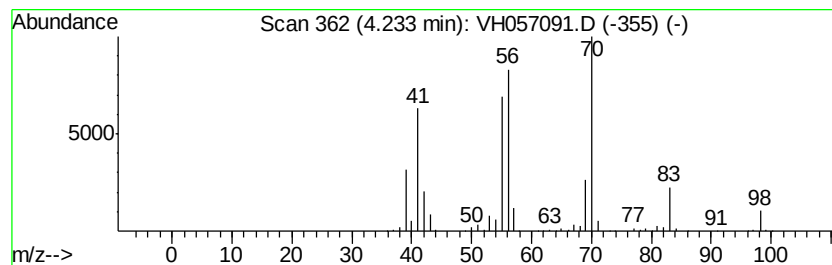
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclopentane, 1,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.23	94.45 ug/l	6034030	Pentafluorobenzene	4.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	97
2		Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	96
3		Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	93
4		Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	93
5		Cyclobutanone, 2,2-dimethyl-	98	C6H10O	001192-14-9	64



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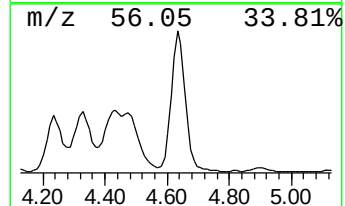
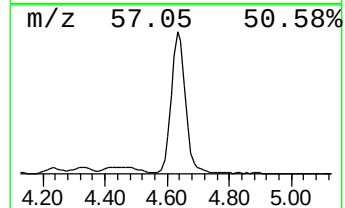
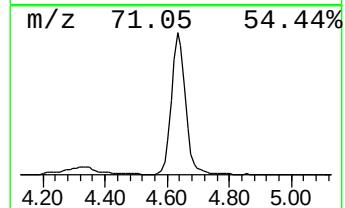
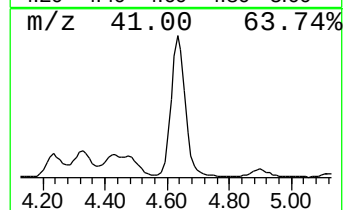
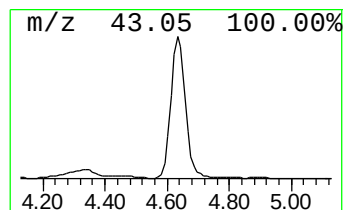
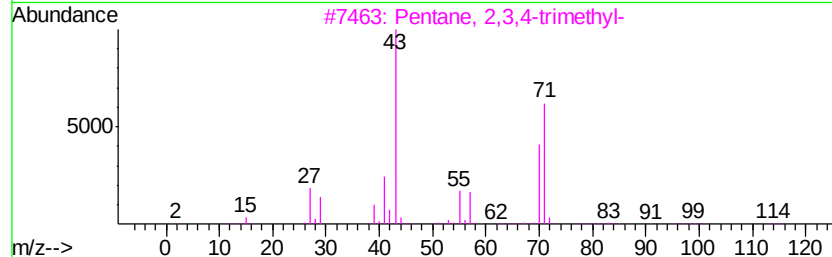
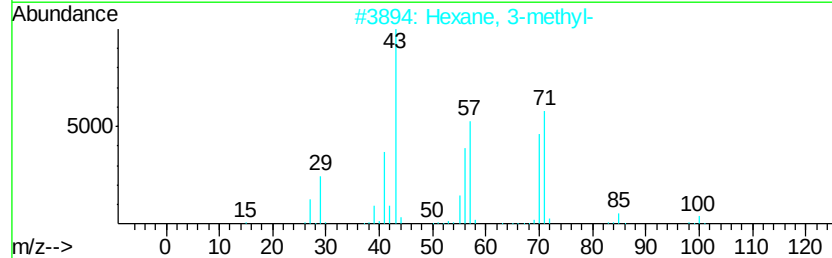
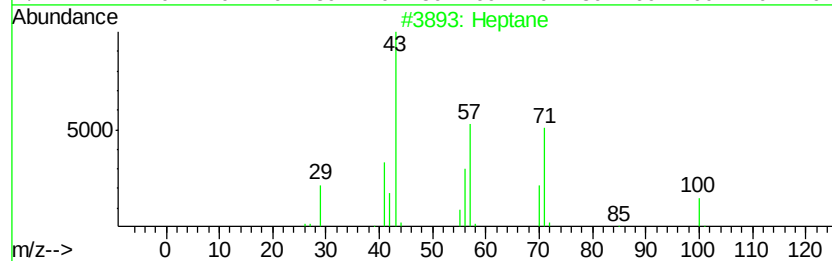
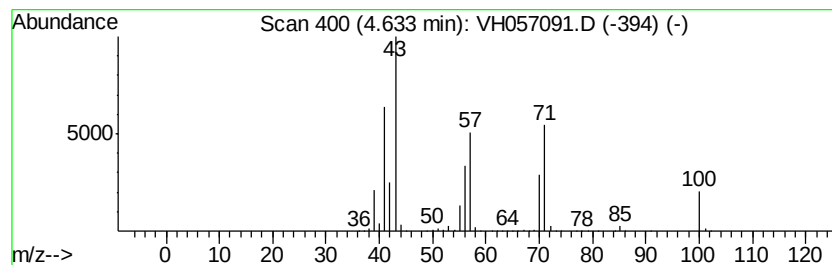
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Heptane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.63	425.87 ug/l	27207500	Pentafluorobenzene	4.81

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane		100	C7H16	000142-82-5	91
2	Hexane, 3-methyl-		100	C7H16	000589-34-4	47
3	Pentane, 2,3,4-trimethyl-		114	C8H18	000565-75-3	38
4	Acetaldehyde, propylhydrazone		100	C5H12N2	007422-88-0	32
5	Pyrrolidine		71	C4H9N	000123-75-1	32



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057091.D
Acq On : 21 Sep 2015 13:46
Operator : SY/FY
Sample : G3727-05
Misc : 5.72/5mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_H
ClientSampled :
MMB7B

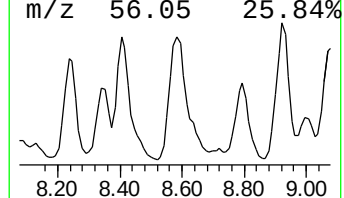
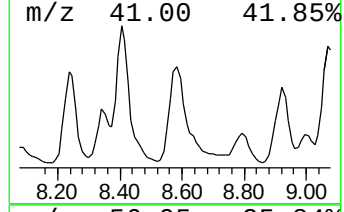
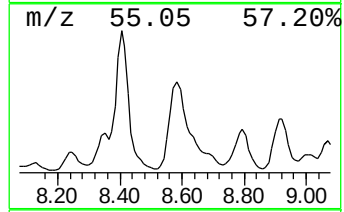
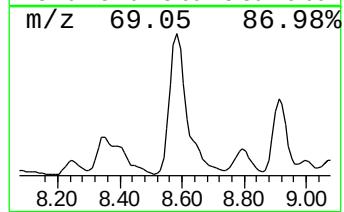
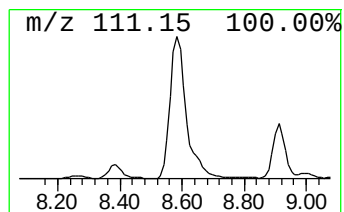
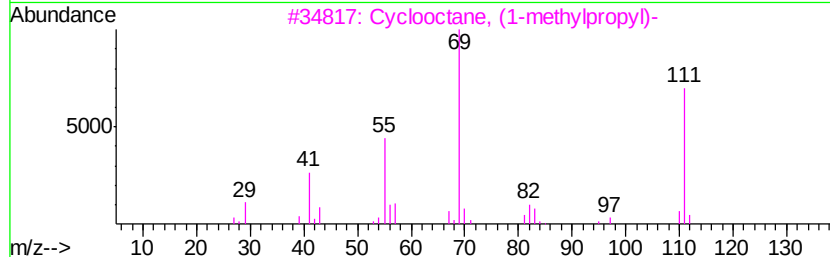
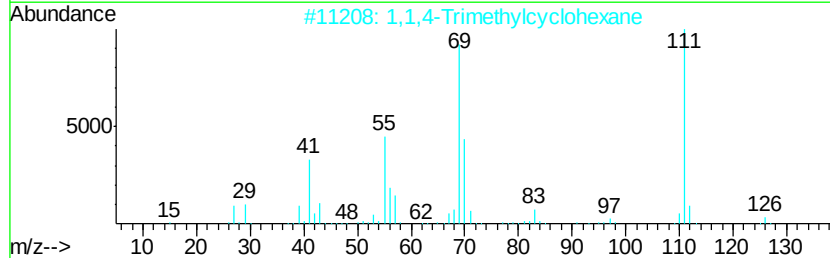
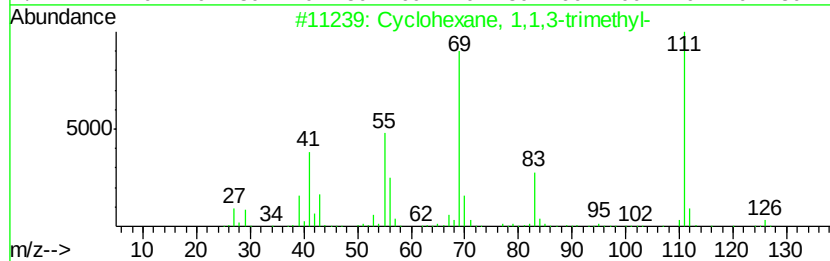
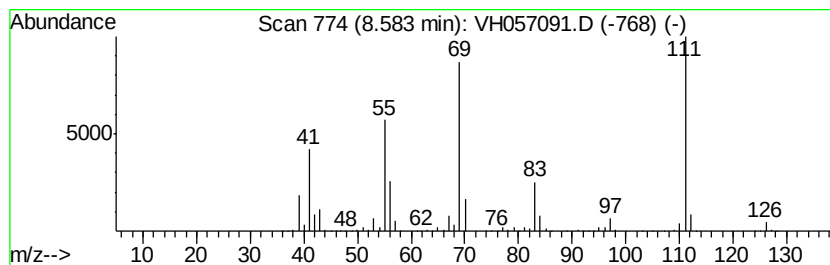
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.58	83.19 ug/l	38229800	Chlorobenzene-d5	9.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	97
2		1,1,4-Trimethylcyclohexane	126	C9H18	007094-27-1	72
3		Cyclooctane, (1-methylpropyl)-	168	C12H24	016538-89-9	64
4		Cyclohexane, 1,2,4-trimethyl-, (...)	126	C9H18	007667-60-9	53
5		Heptane, 4-methylene-	112	C8H16	015918-08-8	43



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057091.D
Acq On : 21 Sep 2015 13:46
Operator : SY/FY
Sample : G3727-05
Misc : 5.72/5mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_H
ClientSampleId :
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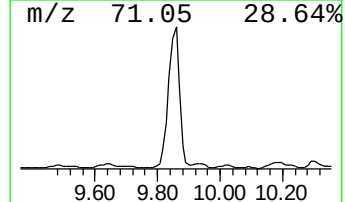
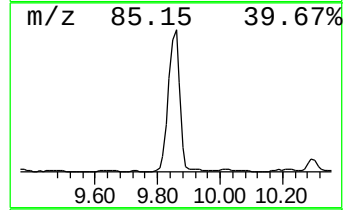
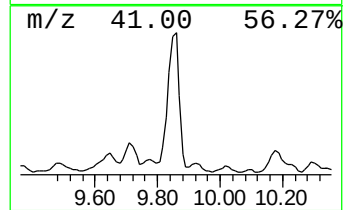
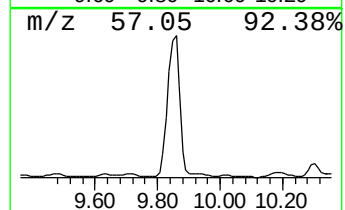
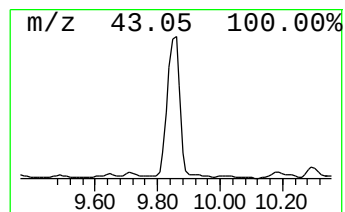
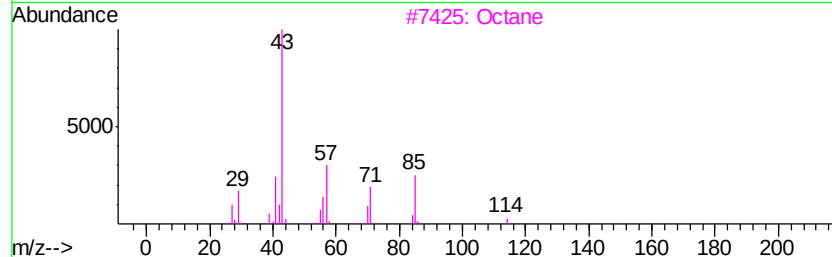
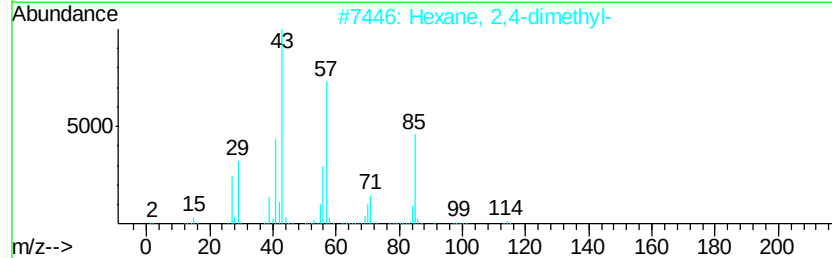
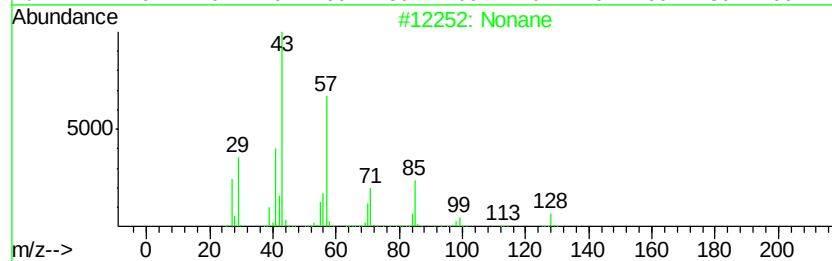
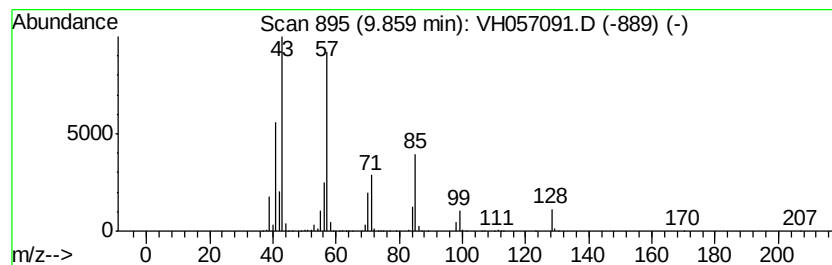
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Nonane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.86	139.84 ug/l	64266200	Chlorobenzene-d5	9.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane		128	C9H20	000111-84-2	94
2	Hexane, 2,4-dimethyl-		114	C8H18	000589-43-5	50
3	Octane		114	C8H18	000111-65-9	50
4	3-Heptanone, 2-methyl-		128	C8H16O	013019-20-0	46
5	Octane, 2,5-dimethyl-		142	C10H22	015869-89-3	38



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057091.D
Acq On : 21 Sep 2015 13:46
Operator : SY/FY
Sample : G3727-05
Misc : 5.72/5mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 6 Sample Multiplier: 1

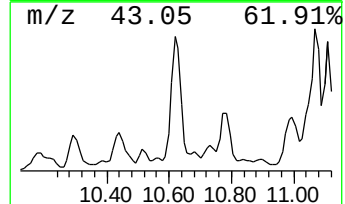
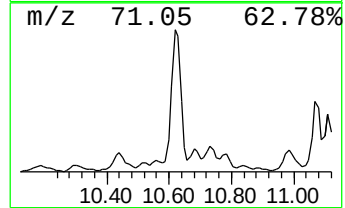
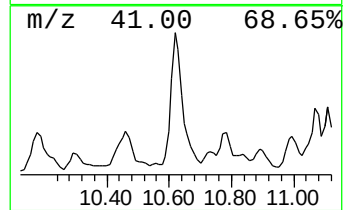
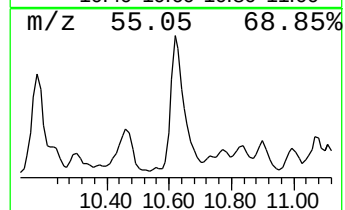
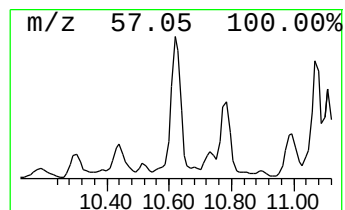
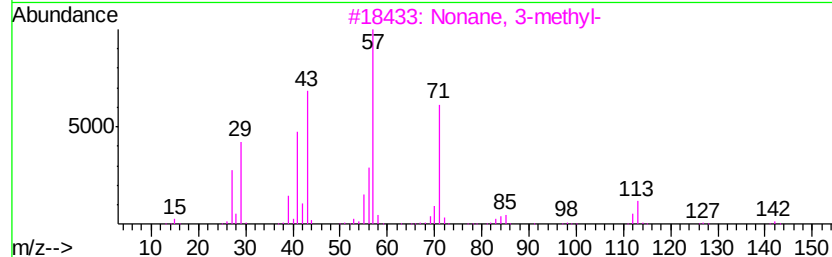
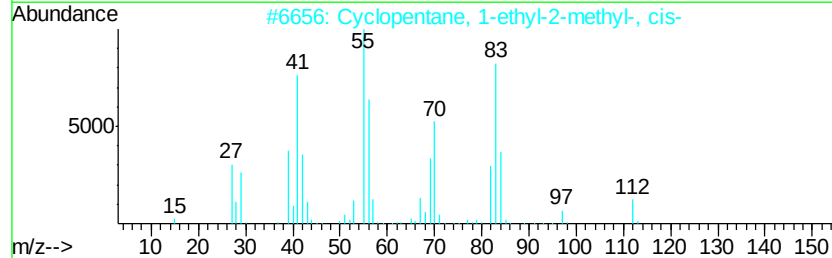
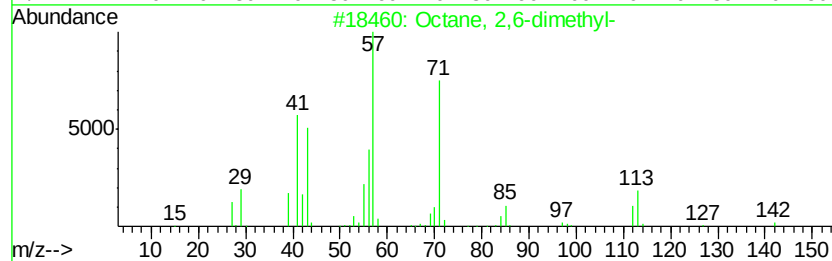
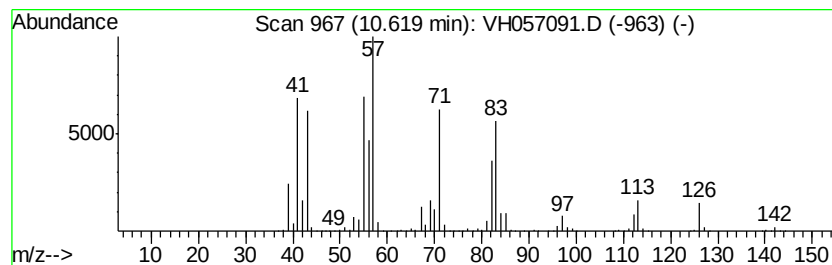
Instrument :
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ClientSampleId :
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Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Octane, 2,6-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.62	107.76 ug/l	49523400	Chlorobenzene-d5	9.71		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	93
2		Cyclopentane, 1-ethyl-2-methyl-, ...	112	C8H16	000930-89-2	46
3		Nonane, 3-methyl-	142	C10H22	005911-04-6	43
4		Ethanone, 1-cyclohexyl-	126	C8H14O	000823-76-7	38
5		Cyclohexanone, 2,3-dimethyl-	126	C8H14O	013395-76-1	38



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057091.D
Acq On : 21 Sep 2015 13:46
Operator : SY/FY
Sample : G3727-05
Misc : 5.72/5mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 6 Sample Multiplier: 1

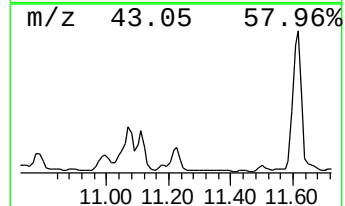
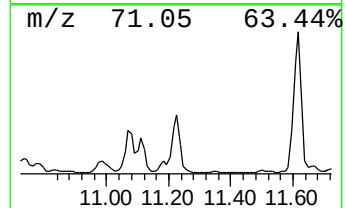
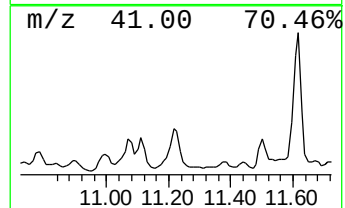
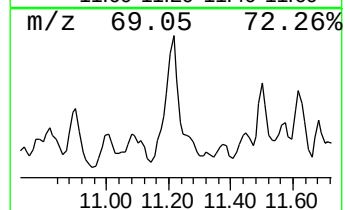
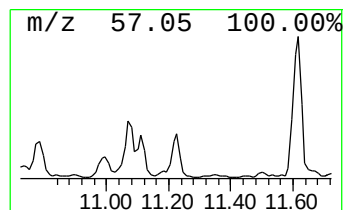
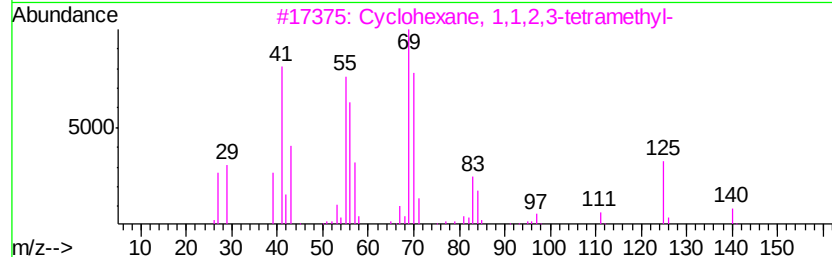
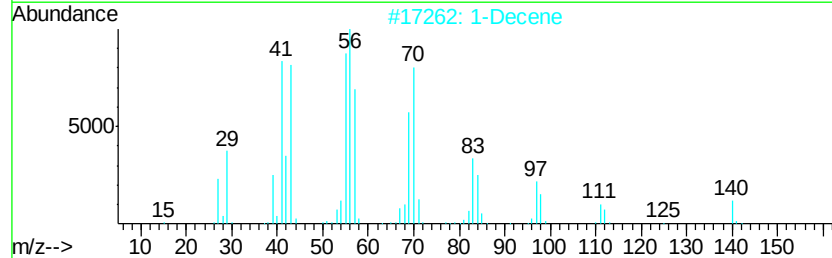
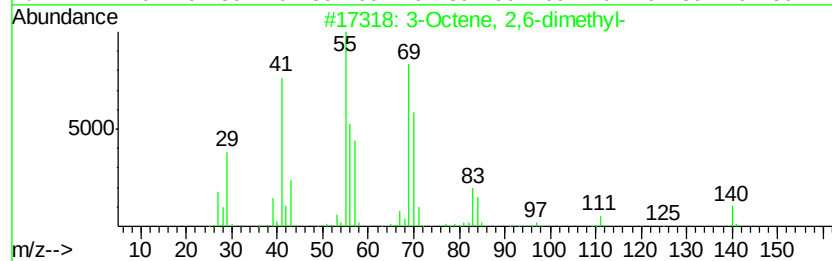
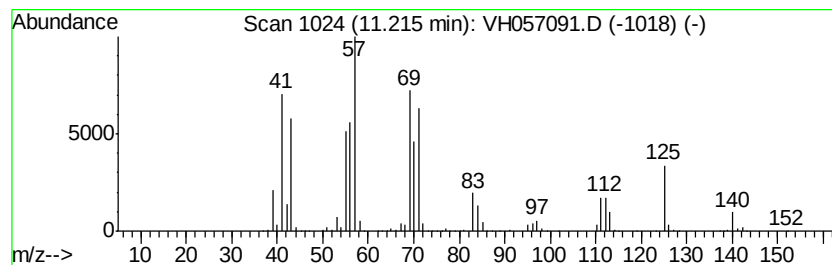
Instrument :
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ClientSampleId :
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Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 3-Octene, 2,6-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.22	149.84 ug/l	30885500	1,4-Dichlorobenzene-d4	12.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Octene, 2,6-dimethyl-	140	C10H20	006874-28-8	60
2	1-Decene	140	C10H20	000872-05-9	49
3	Cyclohexane, 1,1,2,3-tetramethyl-	140	C10H20	006783-92-2	46
4	2-Octene	112	C8H16	000111-67-1	46
5	1R,2c,3t,4t-Tetramethyl-cyclohexane	140	C10H20	1000144-07-3	46



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057091.D
Acq On : 21 Sep 2015 13:46
Operator : SY/FY
Sample : G3727-05
Misc : 5.72/5mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_H
ClientSampleId :
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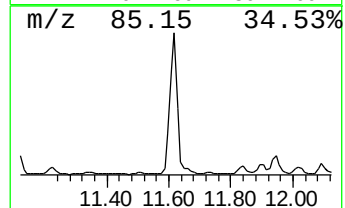
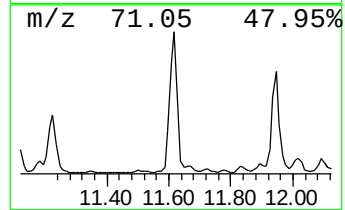
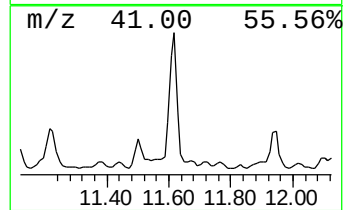
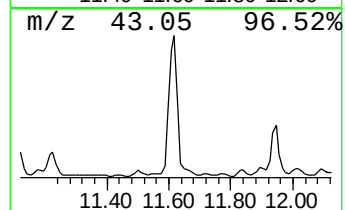
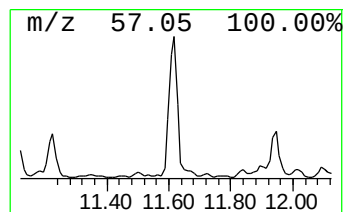
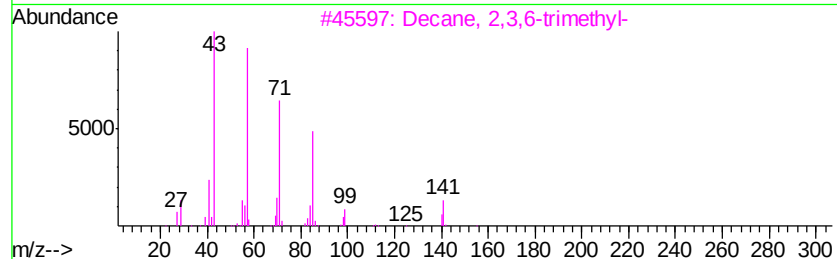
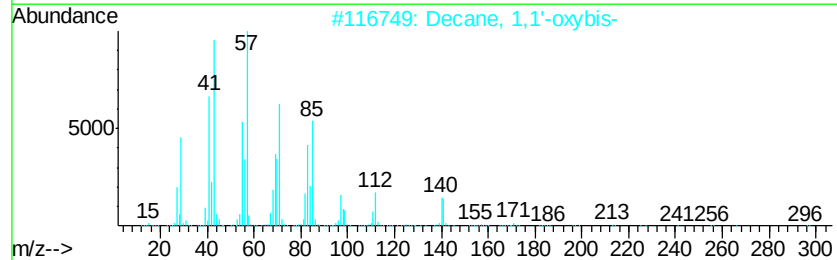
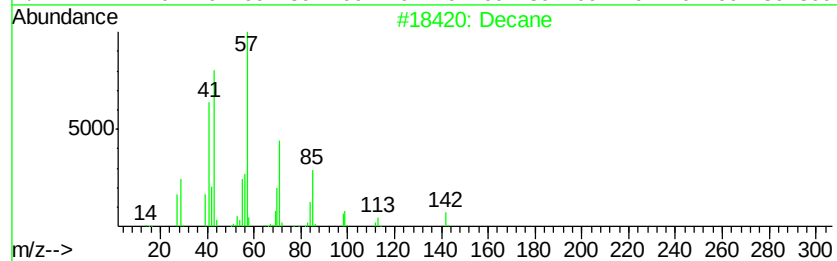
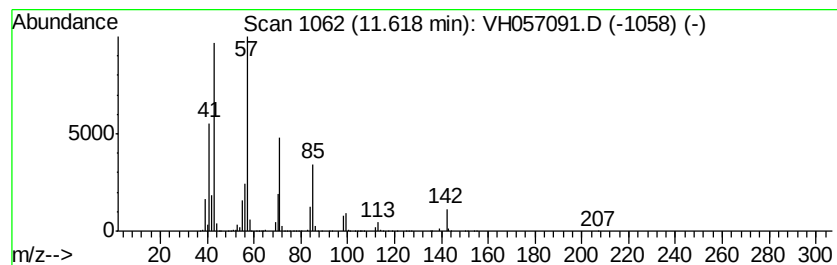
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Decane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.62	271.50 ug/l	55963800	1,4-Dichlorobenzene-d4	12.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane		142	C10H22	000124-18-5	97
2	Decane, 1,1'-oxybis-		298	C20H42O	002456-28-2	64
3	Decane, 2,3,6-trimethyl-		184	C13H28	062238-12-4	59
4	Octane, 2,4,6-trimethyl-		156	C11H24	062016-37-9	59
5	Eicosane		282	C20H42	000112-95-8	59



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057091.D
Acq On : 21 Sep 2015 13:46
Operator : SY/FY
Sample : G3727-05
Misc : 5.72/5mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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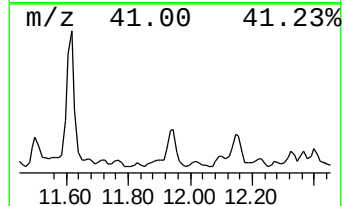
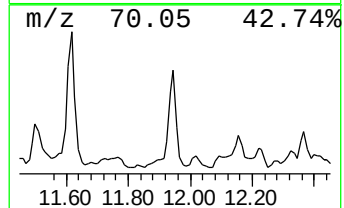
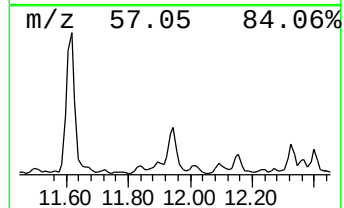
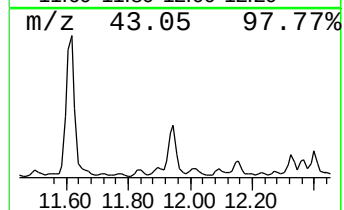
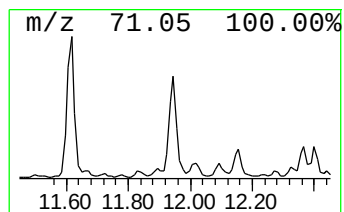
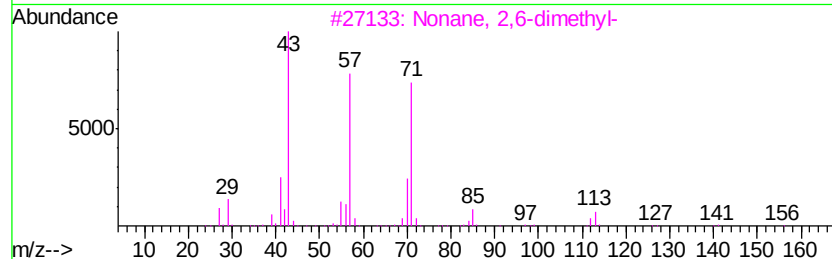
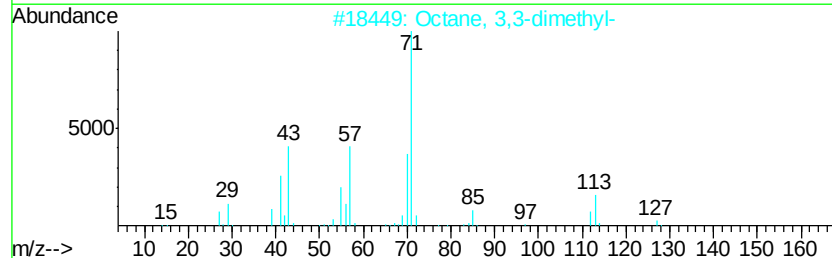
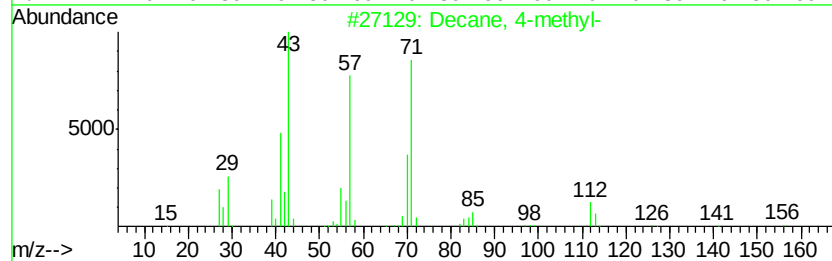
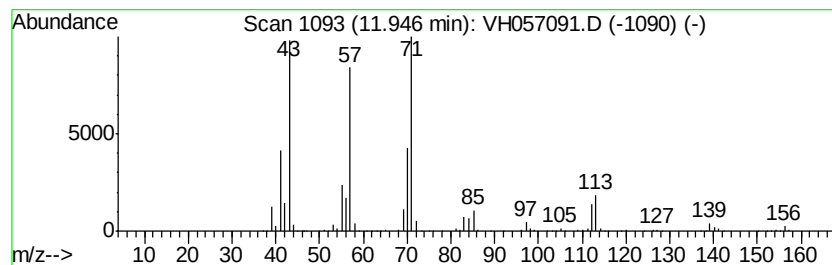
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Decane, 4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.95	121.69 ug/l	25083400	1,4-Dichlorobenzene-d4	12.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 4-methyl-	156	C11H24	002847-72-5	78
2		Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	72
3		Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	64
4		Tridecane, 4-methyl-	198	C14H30	026730-12-1	64
5		Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	59



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057091.D
Acq On : 21 Sep 2015 13:46
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Sample : G3727-05
Misc : 5.72/5mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 6 Sample Multiplier: 1

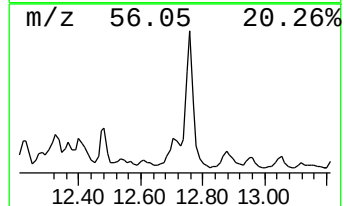
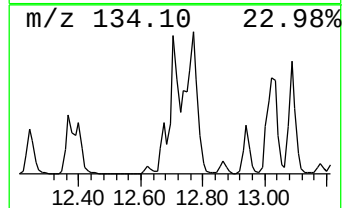
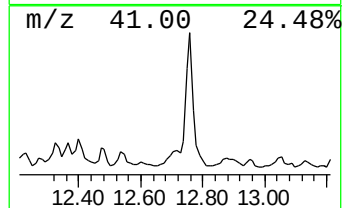
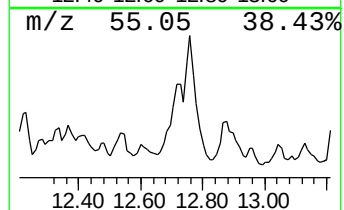
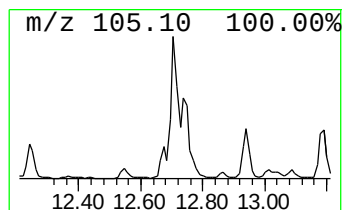
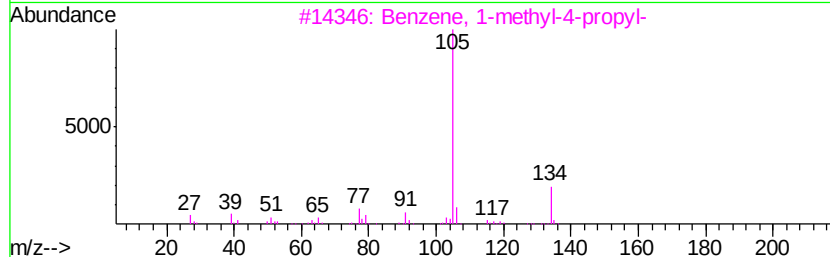
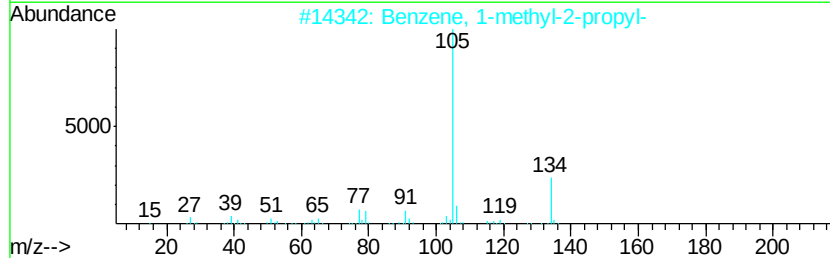
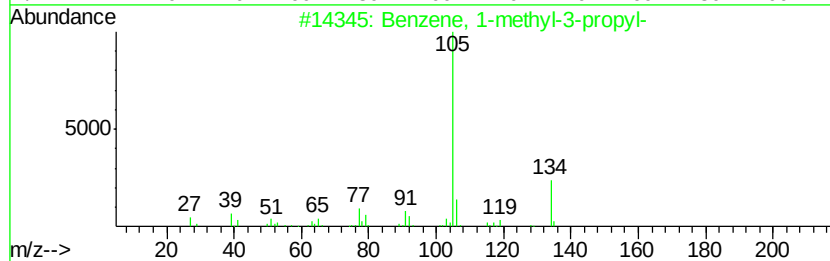
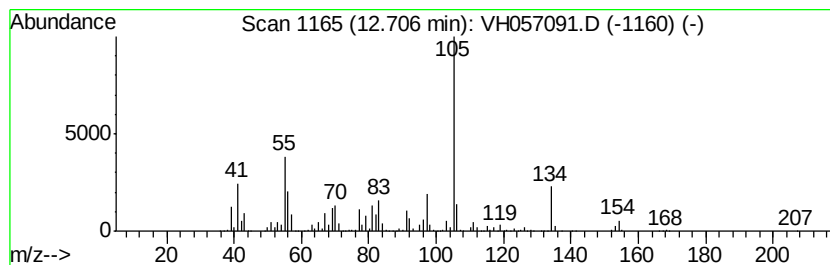
Instrument :
MSVOA_H
ClientSampleId :
MMB7B

Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, 1-methyl-3-propyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.71	88.04 ug/l	18146500	1,4-Dichlorobenzene-d4	12.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-methyl-3-propyl-	134 C10H14	001074-43-7 89
2		Benzene, 1-methyl-2-propyl-	134 C10H14	001074-17-5 50
3		Benzene, 1-methyl-4-propyl-	134 C10H14	001074-55-1 50
4		Benzene, (1-methylpropyl)-	134 C10H14	000135-98-8 50
5		Benzeneacetaldehyde, .alpha.-met...	134 C9H10O	000093-53-8 38



Data Path : W:\HPCHEM1\MSVOA_H\DATA\VH092115\
Data File : VH057091.D
Acq On : 21 Sep 2015 13:46
Operator : SY/FY
Sample : G3727-05
Misc : 5.72/5mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_H
ClientSampleId :
MMB7B

Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Hexane, 3-methyl-	3.96	194.1	ug/l	12403100	1	4.81	3194360	50.0
Cyclopentane, 1,3...	4.23	94.5	ug/l	6034030	1	4.81	3194360	50.0
Heptane	4.63	425.9	ug/l	27207500	1	4.81	3194360	50.0
Cyclohexane, 1,1,...	8.58	83.2	ug/l	38229800	3	9.71	22977900	50.0
Nonane	9.86	139.8	ug/l	64266200	3	9.71	22977900	50.0
Octane, 2,6-dimet...	10.62	107.8	ug/l	49523400	3	9.71	22977900	50.0
3-Octene, 2,6-dim...	11.22	149.8	ug/l	30885500	4	12.49	10306400	50.0
Decane	11.62	271.5	ug/l	55963800	4	12.49	10306400	50.0
Decane, 4-methyl-	11.95	121.7	ug/l	25083400	4	12.49	10306400	50.0
Benzene, 1-methyl...	12.71	88.0	ug/l	18146500	4	12.49	10306400	50.0