

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD111618\
 Data File : VD060395.D
 Acq On : 16 Nov 2018 13:49
 Operator : JC/SY
 Sample : J5911-06
 Misc : 5.00g/5ml/MSVOA D/SOIL
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 FDSB-07(20-22)

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_D\METHOD\82D110618S.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	3.756	217	227	254	rVB2	633911	2497842	9.47%	0.345%
2	4.130	257	265	278	rBV	330176	1082184	4.10%	0.150%
3	5.429	392	397	409	rVB	624195	1882007	7.14%	0.260%
4	6.383	482	494	496	rBV2	2197783	7303165	27.70%	1.010%
5	6.491	503	505	514	rVB	613803	1600598	6.07%	0.221%
6	6.639	514	520	532	rBV	1921251	6023284	22.85%	0.833%
7	6.806	532	537	545	rVB	511446	1209109	4.59%	0.167%
8	6.944	545	551	555	rBV	993380	2701989	10.25%	0.374%
9	7.033	555	560	564	rVV2	883397	2659940	10.09%	0.368%
10	7.111	564	568	579	rVB	1233448	3489143	13.23%	0.482%
11	7.289	579	586	597	rVB	449986	1211318	4.59%	0.167%
12	7.466	597	604	614	rBV	2308174	6061601	22.99%	0.838%
13	8.066	654	665	673	rBV2	5551413	19116861	72.51%	2.643%
14	8.174	673	676	682	rVB	810938	1913273	7.26%	0.265%
15	8.322	682	691	697	rVB	797559	1900891	7.21%	0.263%
16	8.460	697	705	711	rBV2	1305001	3780453	14.34%	0.523%
17	8.656	720	725	732	rBV2	1576556	4267193	16.19%	0.590%
18	8.893	744	749	752	rBV	749273	1631523	6.19%	0.226%
19	8.991	754	759	768	rVB2	5640108	17099113	64.86%	2.364%
20	9.129	768	773	775	rBV3	410451	1270980	4.82%	0.176%
21	9.188	775	779	787	rVB	3190888	8796179	33.36%	1.216%
22	9.316	787	792	799	rVB4	359958	1663214	6.31%	0.230%
23	9.473	799	808	811	rBV2	5819815	16314218	61.88%	2.256%
24	9.522	811	813	818	rVB	2039812	3921971	14.88%	0.542%
25	9.670	823	828	830	rBV	1015969	2517281	9.55%	0.348%
26	9.759	835	837	840	rVB	623142	864118	3.28%	0.119%
27	9.818	840	843	847	rVB	3373581	6660520	25.26%	0.921%
28	9.975	850	859	868	rVB3	2779065	9322603	35.36%	1.289%
29	10.133	868	875	879	rBV3	2182241	5830884	22.12%	0.806%
30	10.211	879	883	885	rBV3	400918	822604	3.12%	0.114%
31	10.261	885	888	892	rVB2	711914	1418326	5.38%	0.196%
32	10.329	892	895	901	rVB2	1399263	3022316	11.46%	0.418%
33	10.467	901	909	913	rBV	4873758	10626948	40.31%	1.469%
34	10.605	918	923	925	rBV	2641466	5874705	22.28%	0.812%

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35	10.654	925	928	932	rVB3	1165481	2381480	9.03%	0.329%
36	10.723	932	935	939	rVB	5341594	10586246	40.15%	1.464%
37	10.802	939	943	946	rBV	6350703	13437526	50.97%	1.858%
38	10.861	946	949	953	rVB4	1250530	3046573	11.56%	0.421%
39	10.949	953	958	961	rBV4	1587222	4096855	15.54%	0.567%
40	11.008	961	964	966	rBV	3542476	5997335	22.75%	0.829%
41	11.067	966	970	972	rBV2	3775622	8386856	31.81%	1.160%
42	11.117	972	975	981	rVB2	6487816	14955997	56.73%	2.068%
43	11.235	981	987	991	rBV	7589868	15894124	60.29%	2.198%
44	11.304	991	994	997	rVB	2631890	3913480	14.84%	0.541%
45	11.392	997	1003	1007	rVB4	1530273	5096196	19.33%	0.705%
46	11.471	1007	1011	1015	rBV2	2363021	5866863	22.25%	0.811%
47	11.550	1015	1019	1023	rBV2	3282548	8812617	33.43%	1.219%
48	11.619	1023	1026	1028	rBV	4982062	7753531	29.41%	1.072%
49	11.668	1028	1031	1037	rVB	5469764	10384154	39.39%	1.436%
50	11.766	1037	1041	1045	rBV3	1926178	4917644	18.65%	0.680%
51	11.835	1045	1048	1051	rVB2	1056595	1990000	7.55%	0.275%
52	11.904	1051	1055	1059	rBV3	6270846	17970588	68.16%	2.485%
53	11.992	1061	1064	1067	rVB2	2797294	5994925	22.74%	0.829%
54	12.042	1067	1069	1072	rBV3	1379555	1942535	7.37%	0.269%
55	12.111	1072	1076	1080	rBV3	10498533	22732964	86.23%	3.143%
56	12.209	1080	1086	1090	rBV3	8322820	26364486	100.00%	3.646%
57	12.298	1093	1095	1098	rVB	3485406	5066703	19.22%	0.701%
58	12.366	1098	1102	1105	rBV2	6188270	15515522	58.85%	2.145%
59	12.435	1105	1109	1110	rBV2	5248136	9204233	34.91%	1.273%
60	12.504	1114	1116	1118	rBV2	2576192	3487110	13.23%	0.482%
61	12.544	1118	1120	1123	rBV	5347404	8301113	31.49%	1.148%
62	12.603	1123	1126	1129	rVB2	6704793	11088788	42.06%	1.533%
63	12.652	1129	1131	1133	rVB	1992857	2376088	9.01%	0.329%
64	12.711	1133	1137	1140	rBV4	3863720	8974654	34.04%	1.241%
65	12.799	1140	1146	1149	rBV5	6832927	21838377	82.83%	3.020%
66	12.849	1149	1151	1154	rVB	7290700	11549034	43.81%	1.597%
67	12.908	1154	1157	1159	rBV4	1963342	4320309	16.39%	0.597%
68	12.937	1159	1160	1162	rVB	2385050	2410101	9.14%	0.333%
69	12.986	1162	1165	1168	rBV3	4699567	12074917	45.80%	1.670%
70	13.036	1168	1170	1172	rBV2	2700307	3961692	15.03%	0.548%
71	13.075	1172	1174	1177	rVB2	6212156	9162581	34.75%	1.267%

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72	13.134	1177	1180	1184	rVB4	4725161	9096672	34.50%	1.258%
73	13.193	1184	1186	1188	rBV	3993610	6302648	23.91%	0.872%
74	13.242	1188	1191	1195	rBV4	5688506	10189429	38.65%	1.409%
75	13.311	1195	1198	1200	rBV2	7090070	12139064	46.04%	1.679%
76	13.390	1204	1206	1210	rBV4	4361145	7693224	29.18%	1.064%
77	13.449	1210	1212	1215	rVB2	2324339	2312784	8.77%	0.320%
78	13.518	1217	1219	1222	rVB2	4049923	5382156	20.41%	0.744%
79	13.577	1222	1225	1230	rVB5	5249975	11877657	45.05%	1.642%
80	13.675	1230	1235	1238	rBV4	7697400	24488698	92.89%	3.386%
81	13.764	1242	1244	1246	rBV	2895750	4575617	17.36%	0.633%
82	13.892	1254	1257	1259	rBV2	4575906	9375152	35.56%	1.296%
83	13.921	1259	1260	1263	rVB	4576647	4859319	18.43%	0.672%
84	14.010	1265	1269	1273	rBV4	4687854	11589937	43.96%	1.603%
85	14.089	1274	1277	1280	rVB4	3996937	8624445	32.71%	1.193%
86	14.148	1280	1283	1285	rBV4	5446124	10637426	40.35%	1.471%
87	14.187	1285	1287	1292	rVB3	5786943	13095757	49.67%	1.811%
88	14.295	1295	1298	1301	rBV4	4284526	7898150	29.96%	1.092%
89	14.443	1310	1313	1317	rBV3	2263901	5612908	21.29%	0.776%
90	14.541	1320	1323	1327	rVV3	5563659	15767863	59.81%	2.180%
91	14.600	1327	1329	1332	rVB2	4404252	6105707	23.16%	0.844%
92	14.669	1334	1336	1340	rVB3	2827798	5072823	19.24%	0.701%
93	14.856	1352	1355	1358	rVB2	4083683	7457185	28.28%	1.031%
94	14.905	1358	1360	1361	rVV2	2724449	3300717	12.52%	0.456%
95	14.964	1364	1366	1368	rVB2	1080248	1016900	3.86%	0.141%
96	15.004	1368	1370	1372	rBV3	1907224	3182155	12.07%	0.440%
97	15.083	1375	1378	1384	rVB6	4094972	11054020	41.93%	1.529%
98	15.151	1384	1385	1389	rVB2	707233	979820	3.72%	0.135%
99	15.269	1395	1397	1400	rVB3	1163382	2005325	7.61%	0.277%
100	15.368	1400	1407	1410	rVB4	1229211	3275153	12.42%	0.453%

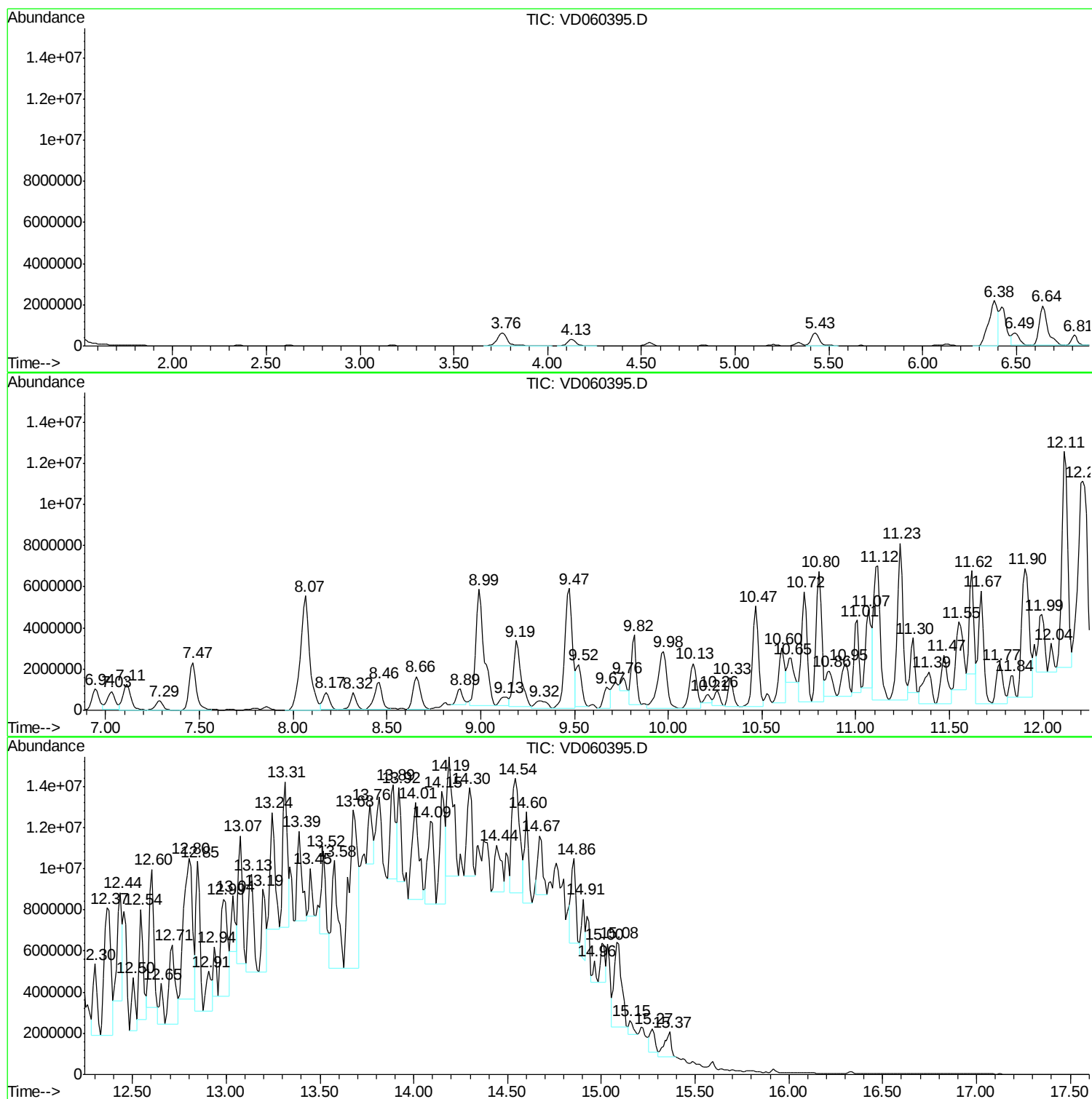
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Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110618S.M
Quant Title : SW846 8260

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



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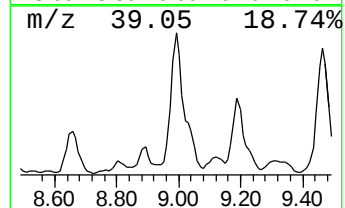
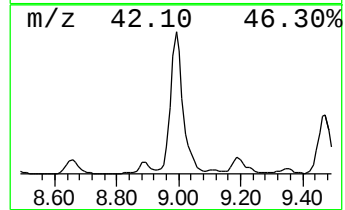
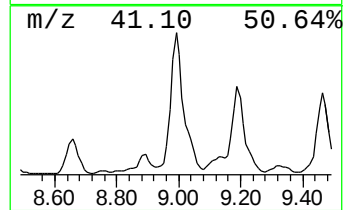
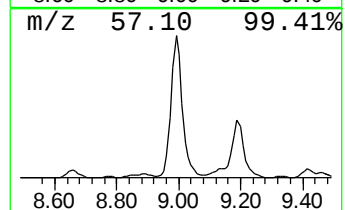
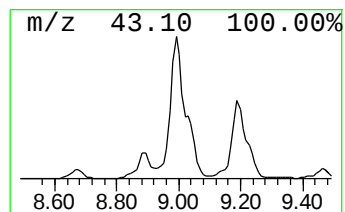
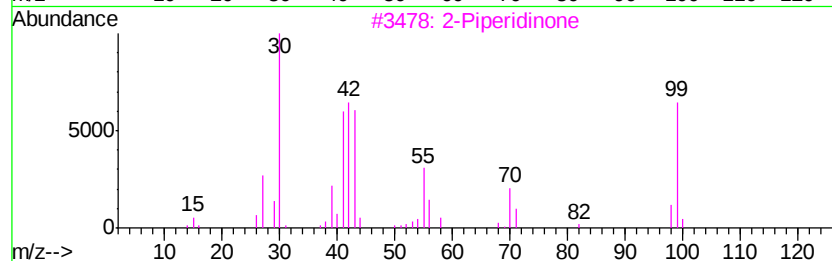
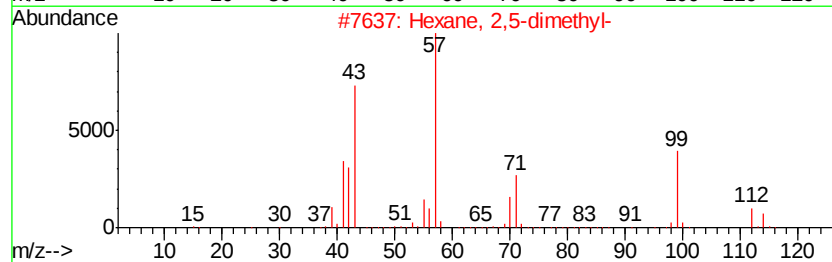
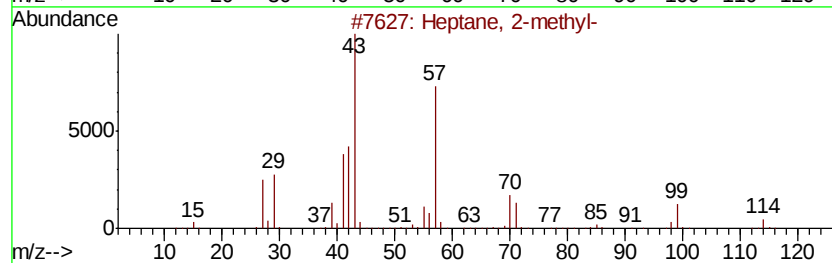
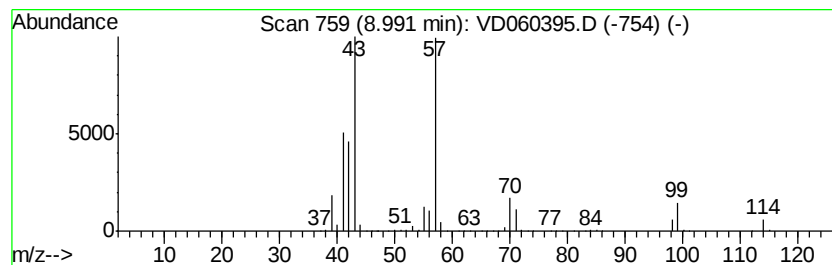
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110618S.M
Quant Title : SW846 8260

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

Peak Number 1 Heptane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.99	141.05 ug/l	17099100	1,4-Difluorobenzene	7.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane, 2-methyl-	114	C8H18	000592-27-8	90
2		Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	78
3		2-Piperidinone	99	C5H9NO	000675-20-7	40
4		Heptane, 3,5-dimethyl-	128	C9H20	000926-82-9	38
5		Isobutane	58	C4H10	000075-28-5	38



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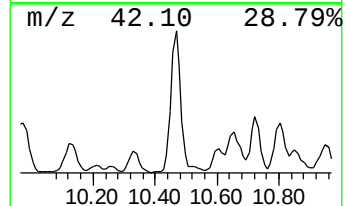
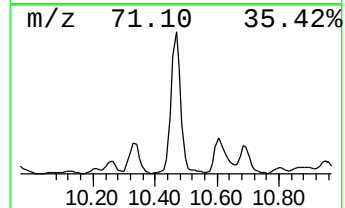
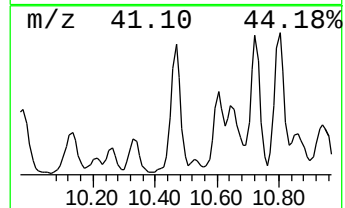
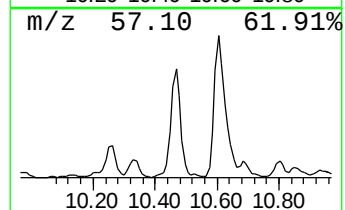
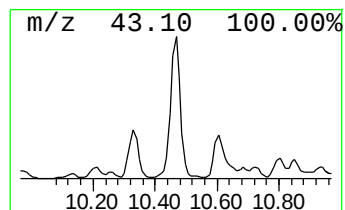
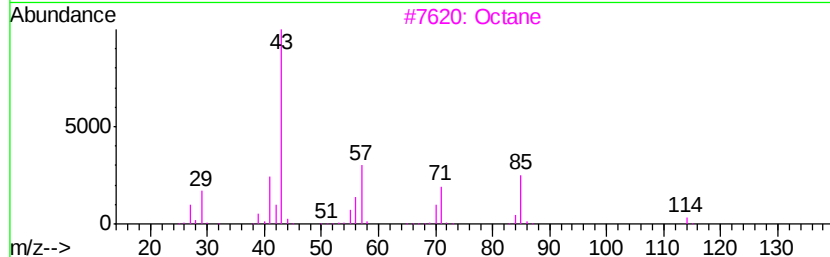
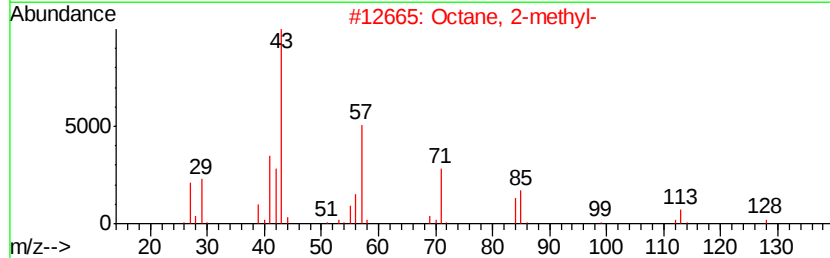
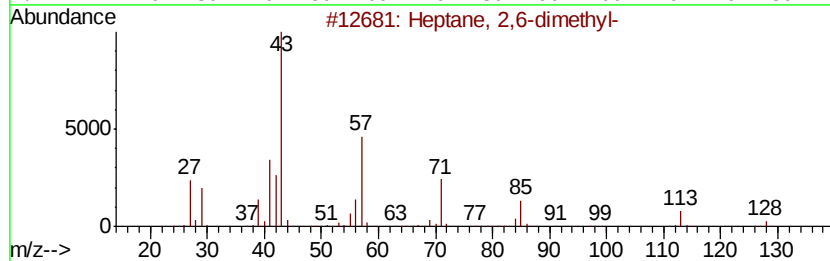
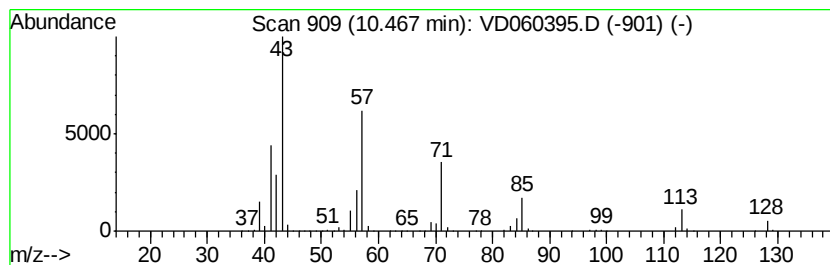
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Heptane, 2,6-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.47	135.77 ug/l	10626900	Chlorobenzene-d5	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	90
2		Octane, 2-methyl-	128	C9H20	003221-61-2	78
3		Octane	114	C8H18	000111-65-9	53
4		Octane, 3-ethyl-2,7-dimethyl-	170	C12H26	062183-55-5	50
5		Decane, 2,4-dimethyl-	170	C12H26	002801-84-5	50



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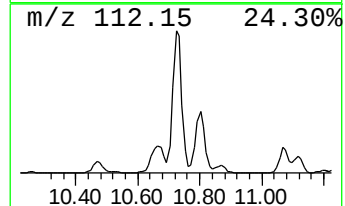
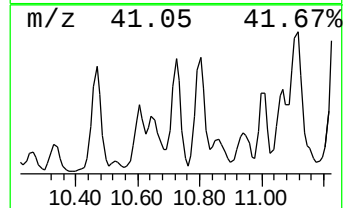
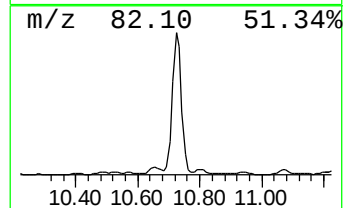
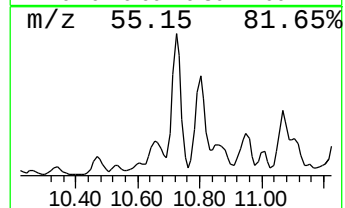
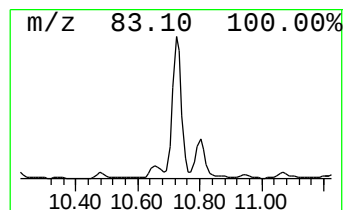
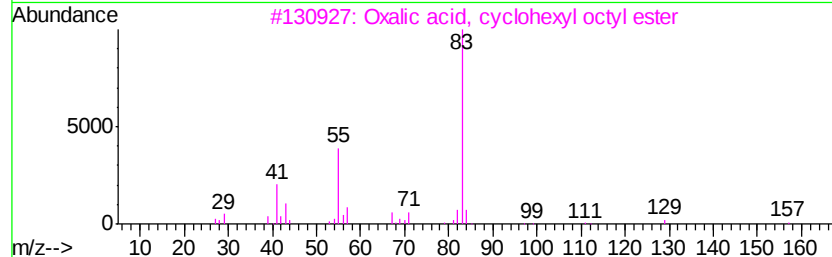
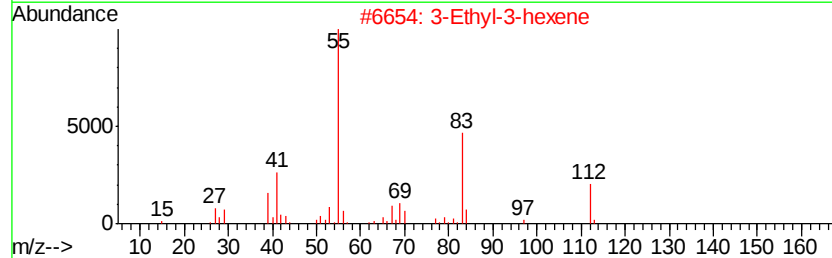
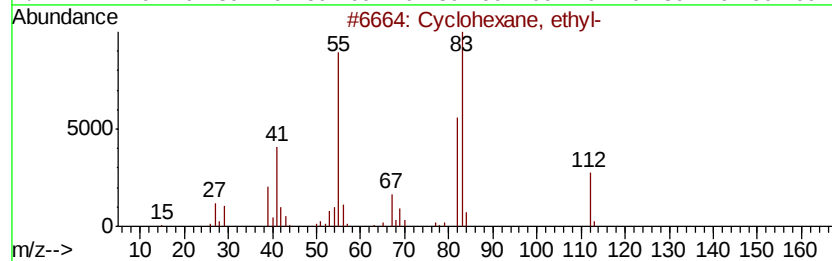
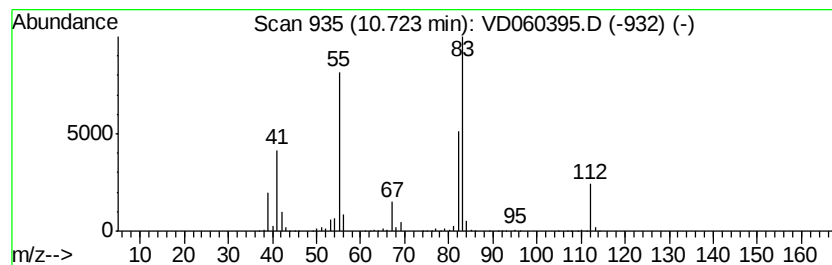
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Quant Title : SW846 8260

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

Peak Number 3 Cyclohexane, ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.72	135.25 ug/l	10586200	Chlorobenzene-d5	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, ethyl-	112	C8H16	001678-91-7	90
2		3-Ethyl-3-hexene	112	C8H16	016789-51-8	53
3		Oxalic acid, cyclohexyl octyl ester	284	C16H28O4	1000309-31-0	53
4		2-Hexene, 2,4-dimethyl-	112	C8H16	014255-23-3	52
5		3,4-Dimethyl-2-hexene	112	C8H16	002213-37-8	52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111618\
Data File : VD060395.D
Acq On : 16 Nov 2018 13:49
Operator : JC/SY
Sample : J5911-06
Misc : 5.00µl/5ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
FDSB-07(20-22)

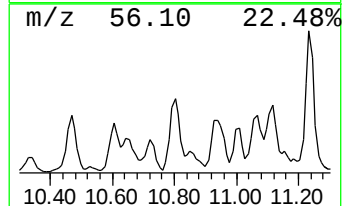
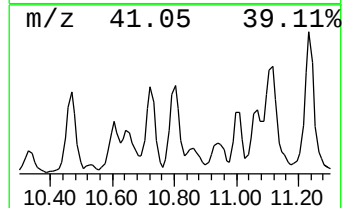
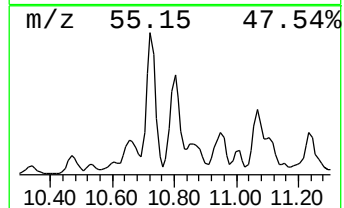
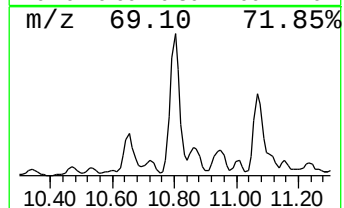
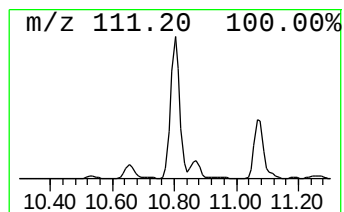
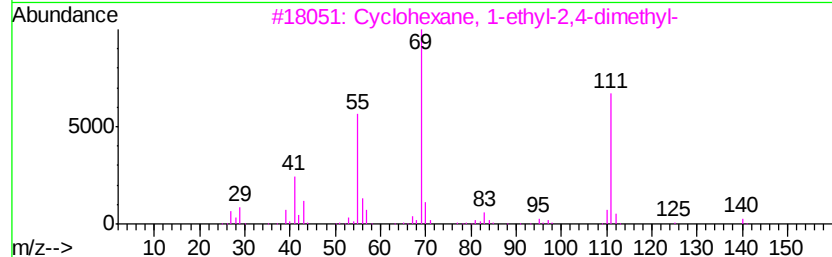
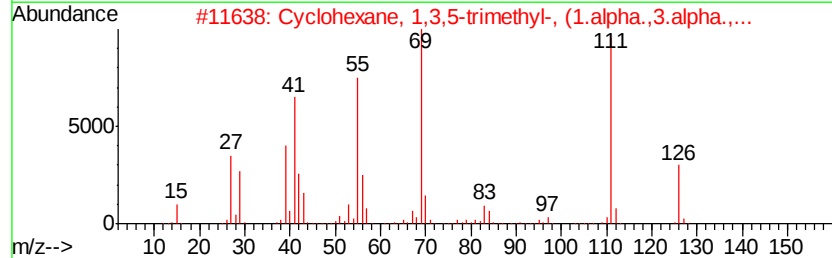
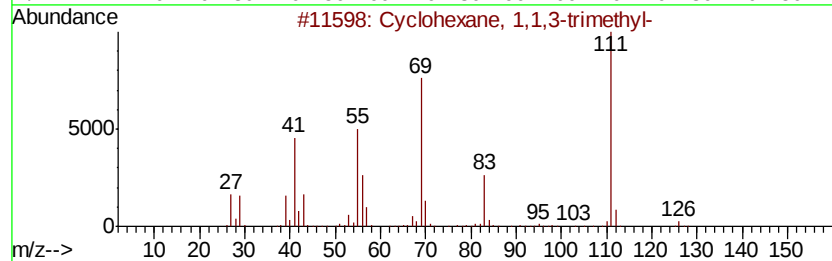
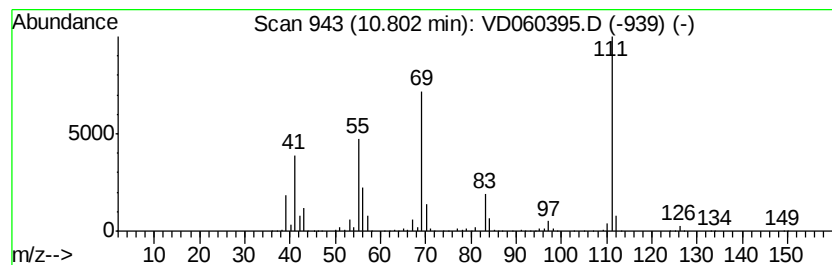
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Quant Title : SW846 8260

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

Peak Number 4 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.80	171.68 ug/l	13437500	Chlorobenzene-d5	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	95
2		Cyclohexane, 1,3,5-trimethyl-, (...)	126	C9H18	001795-26-2	64
3		Cyclohexane, 1-ethyl-2,4-dimethyl-	140	C10H20	061142-69-6	59
4		Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	53
5		Cyclooctane, butyl-	168	C12H24	016538-93-5	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111618\
Data File : VD060395.D
Acq On : 16 Nov 2018 13:49
Operator : JC/SY
Sample : J5911-06
Misc : 5.00µl/5ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
FDSB-07(20-22)

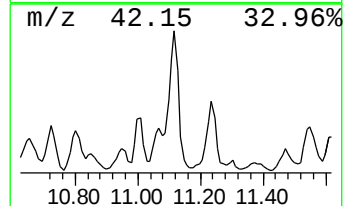
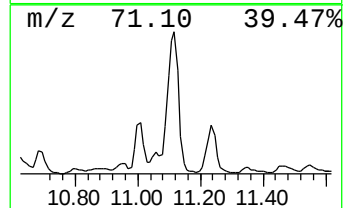
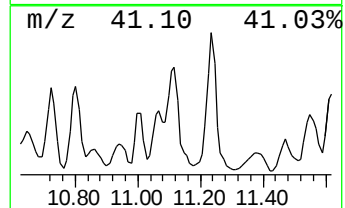
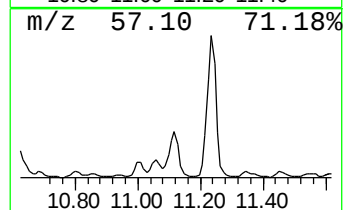
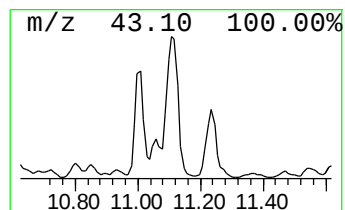
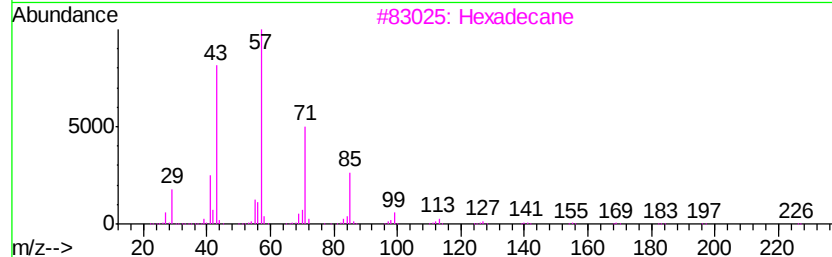
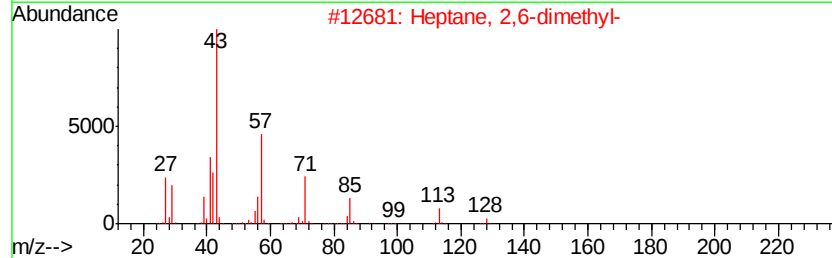
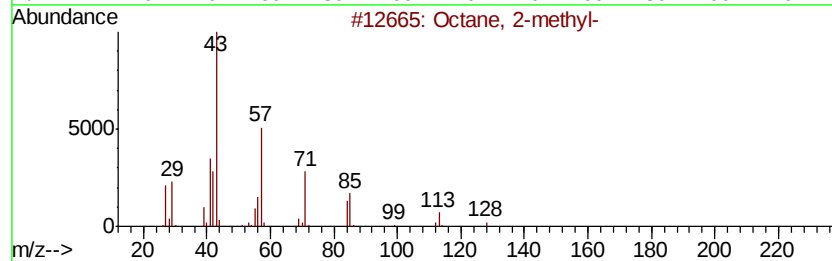
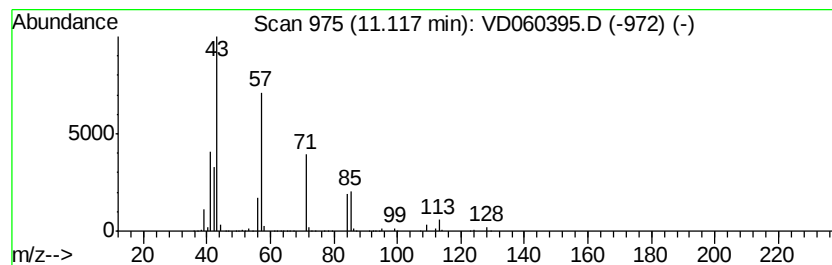
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.12	191.08 ug/l	14956000	Chlorobenzene-d5	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 2-methyl-	128	C9H20	003221-61-2	86
2		Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	78
3		Hexadecane	226	C16H34	000544-76-3	59
4		Decane, 2,4-dimethyl-	170	C12H26	002801-84-5	59
5		Octane, 3-ethyl-2,7-dimethyl-	170	C12H26	062183-55-5	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111618\
Data File : VD060395.D
Acq On : 16 Nov 2018 13:49
Operator : JC/SY
Sample : J5911-06
Misc : 5.00µl/5ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
FDSB-07(20-22)

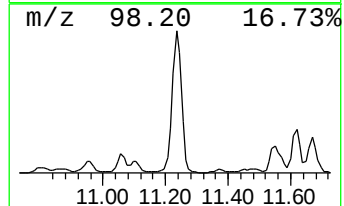
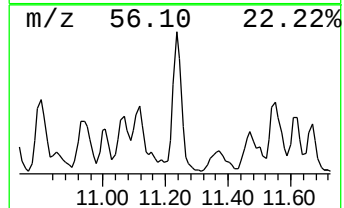
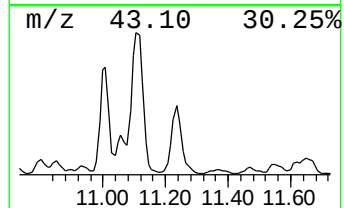
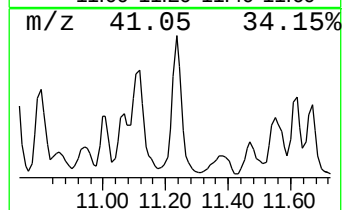
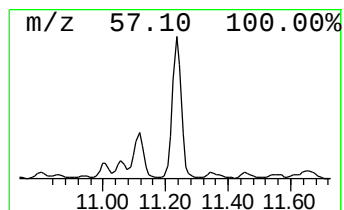
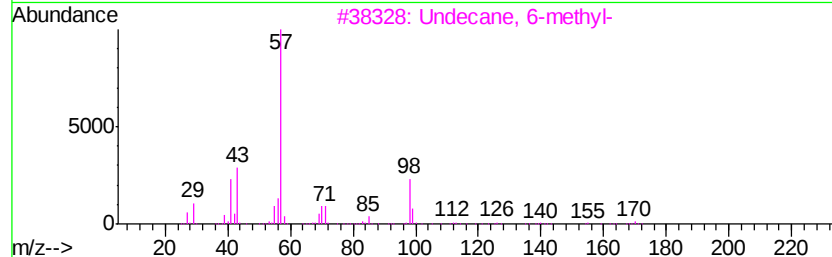
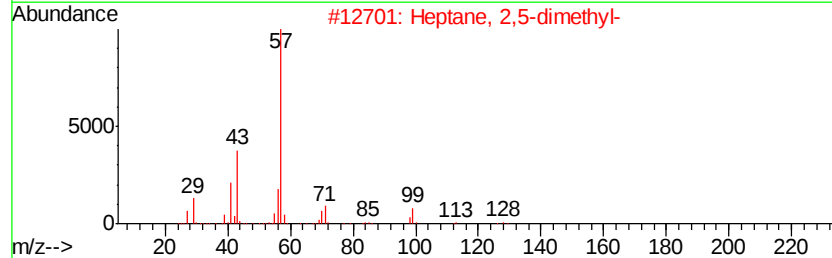
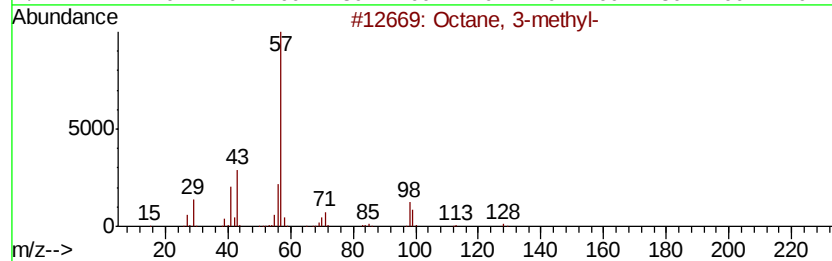
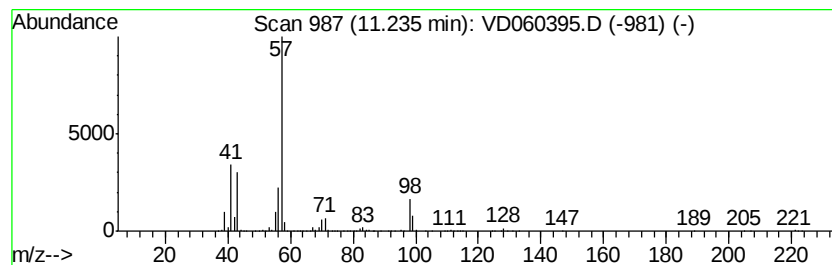
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.23	203.07 ug/l	15894100	Chlorobenzene-d5	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 3-methyl-	128	C9H20	002216-33-3	91
2		Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	64
3		Undecane, 6-methyl-	170	C12H26	017302-33-9	59
4		Heptane, 4-(1-methylethyl)-	142	C10H22	052896-87-4	59
5		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111618\
Data File : VD060395.D
Acq On : 16 Nov 2018 13:49
Operator : JC/SY
Sample : J5911-06
Misc : 5.00g/5ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

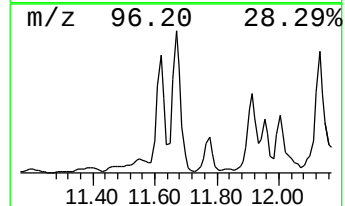
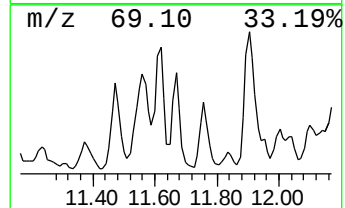
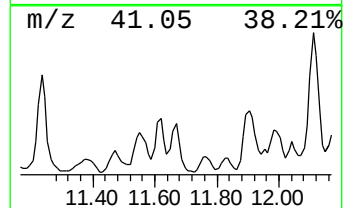
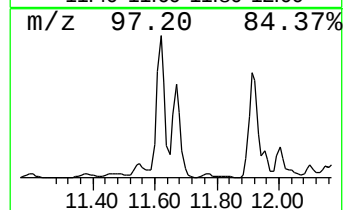
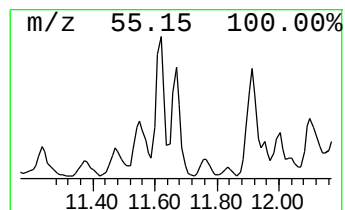
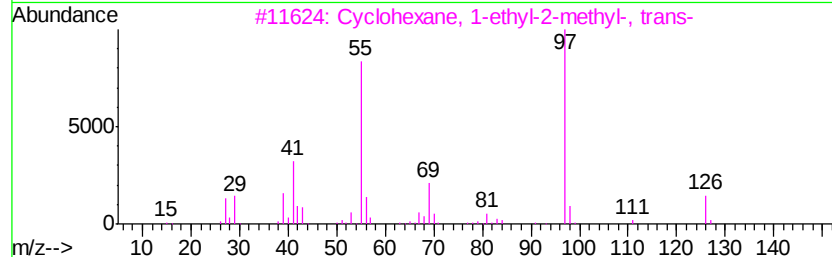
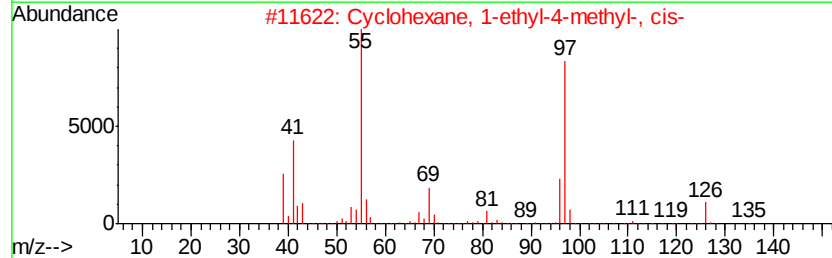
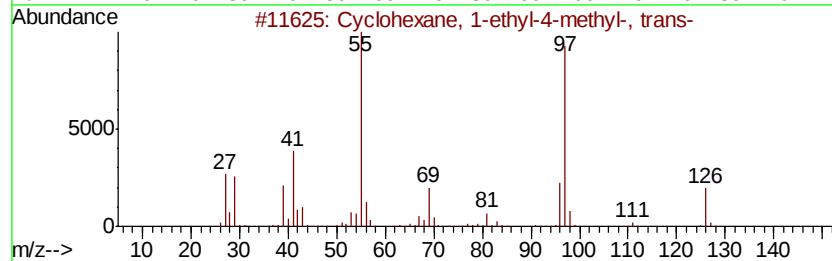
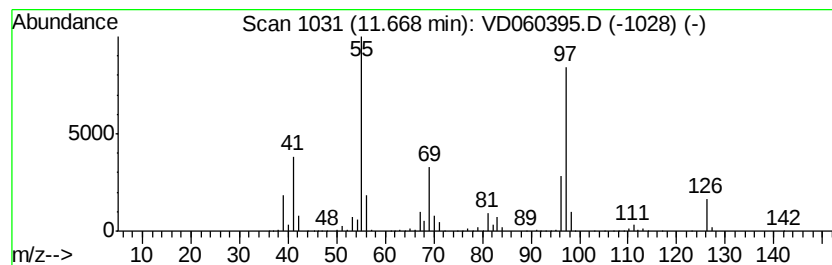
Instrument :
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ClientSampleId :
FDSB-07(20-22)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110618S.M
Quant Title : SW846 8260

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

Peak Number 7 Cyclohexane, 1-ethyl-4-meth... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.67	132.67 ug/l	10384200	Chlorobenzene-d5	11.30		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cvclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	91
2		Cvclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	91
3		Cvclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	87
4		1-Ethyl-4-methylcvclohexane	126	C9H18	003728-56-1	80
5		cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	80



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111618\
Data File : VD060395.D
Acq On : 16 Nov 2018 13:49
Operator : JC/SY
Sample : J5911-06
Misc : 5.00µl/5ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
FDSB-07(20-22)

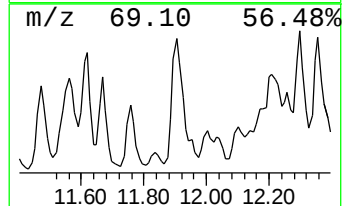
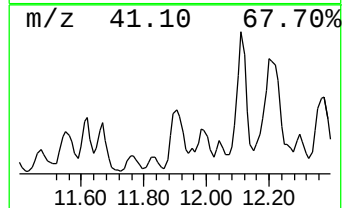
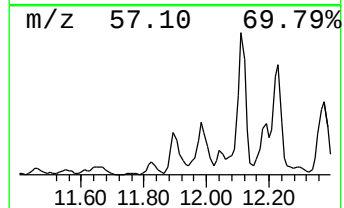
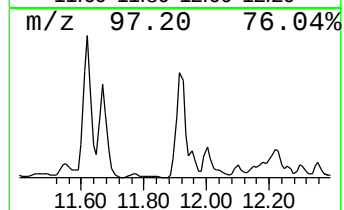
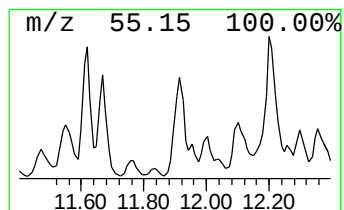
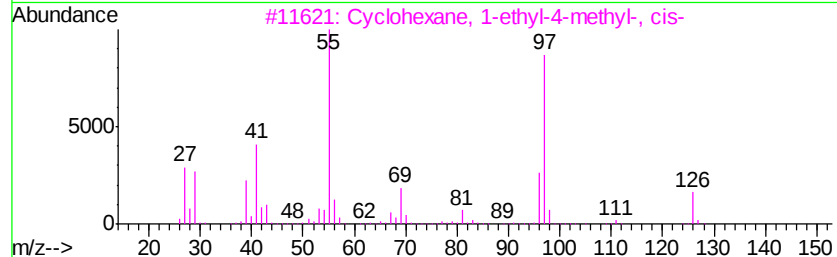
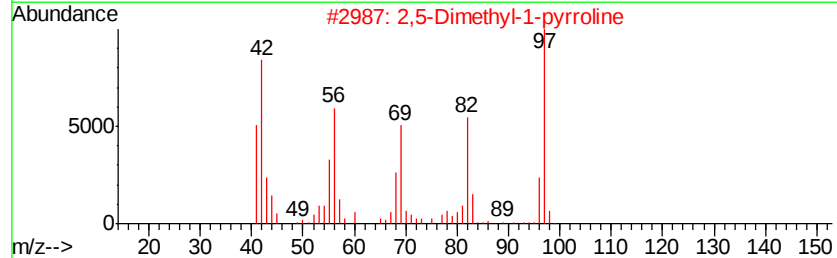
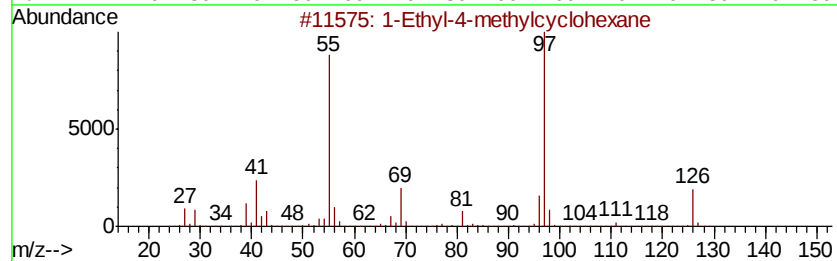
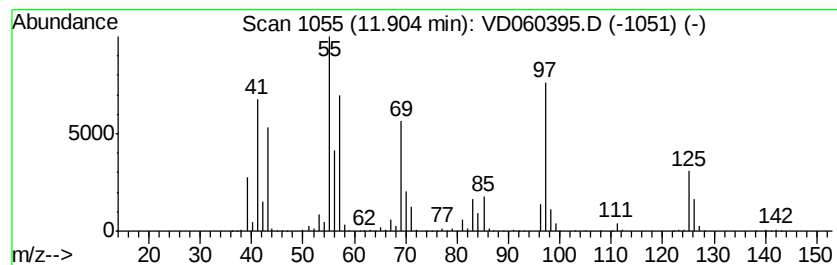
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Quant Title : SW846 8260

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

Peak Number 8 unknown11.90 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	229.60 ug/l	17970600	Chlorobenzene-d5	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	49
2		2,5-Dimethyl-1-pyrroline	97	C6H11N	000822-64-0	46
3		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	46
4		3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	46
5		1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111618\
Data File : VD060395.D
Acq On : 16 Nov 2018 13:49
Operator : JC/SY
Sample : J5911-06
Misc : 5.00µl/5ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

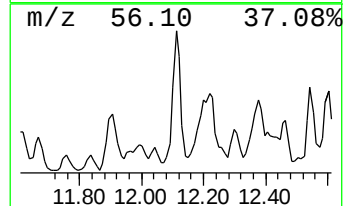
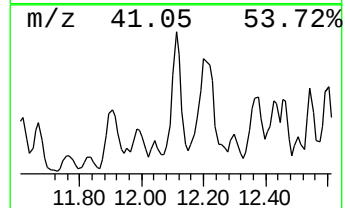
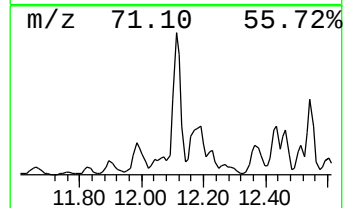
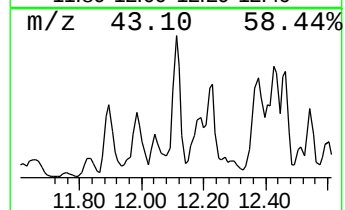
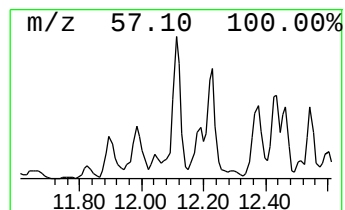
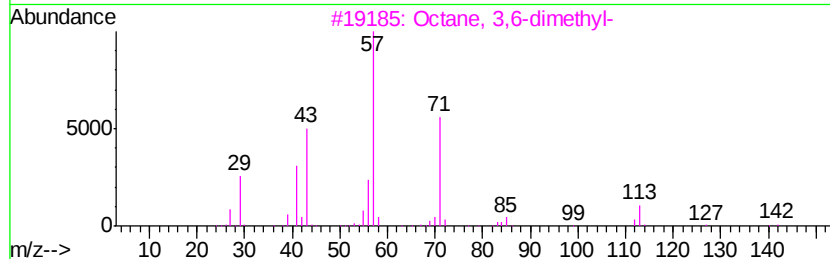
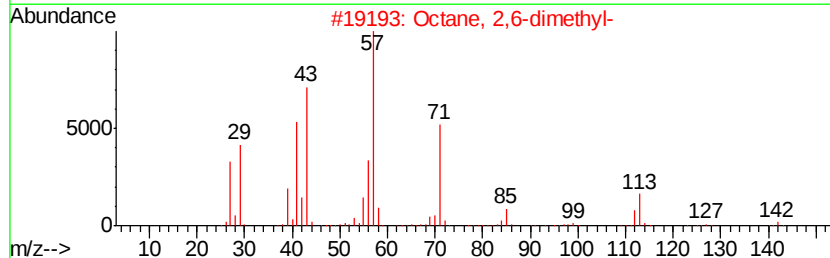
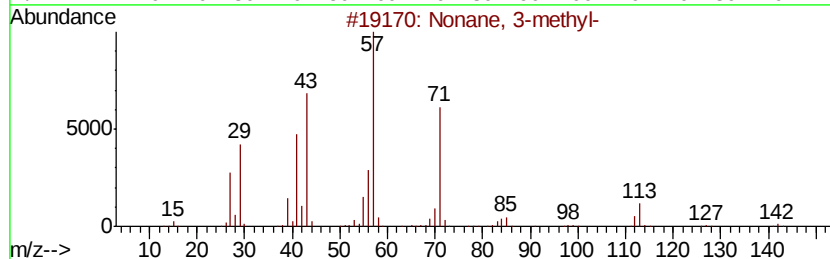
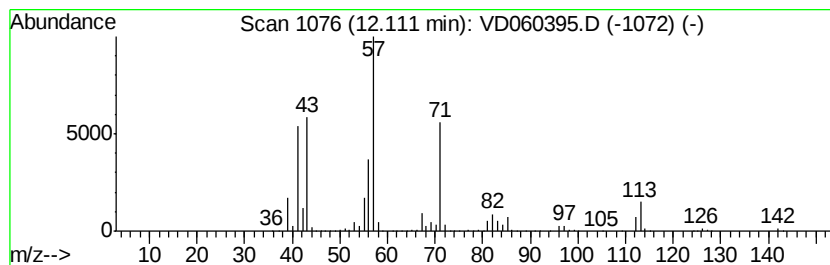
Instrument :
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ClientSampleId :
FDSB-07(20-22)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110618S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.11	290.44 ug/l	22733000	Chlorobenzene-d5	11.30		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 3-methyl-	142	C10H22	005911-04-6	94
2		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	91
3		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	86
4		Sulfurous acid, butyl 2-ethylhex...	250	C12H26O3S	1000309-18-8	64
5		Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111618\
Data File : VD060395.D
Acq On : 16 Nov 2018 13:49
Operator : JC/SY
Sample : J5911-06
Misc : 5.00g/5ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
FDSB-07(20-22)

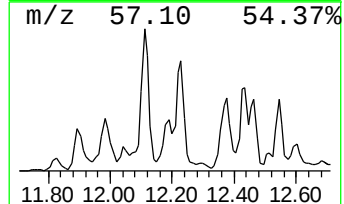
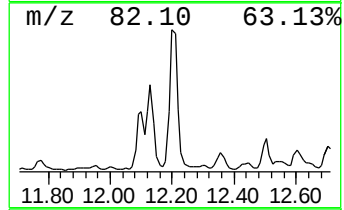
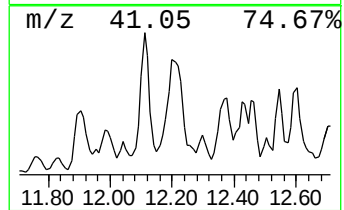
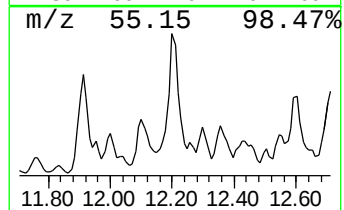
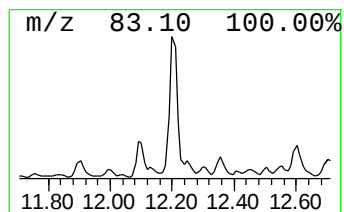
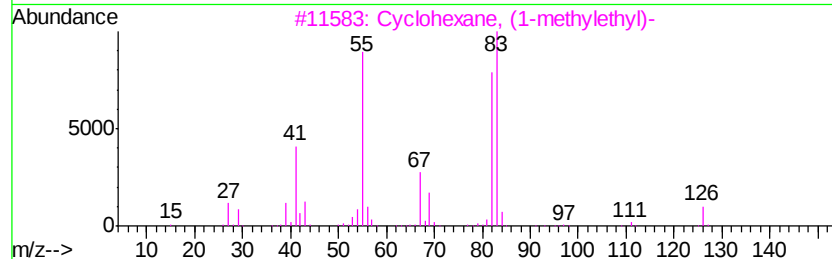
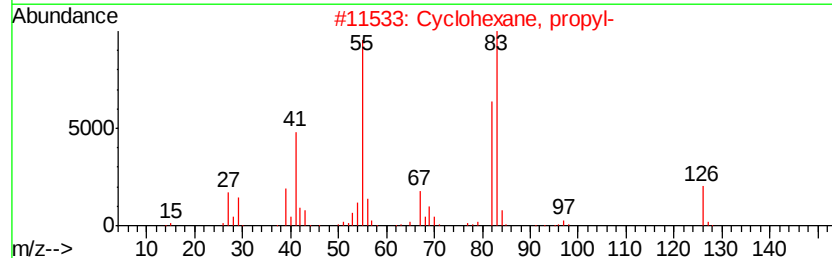
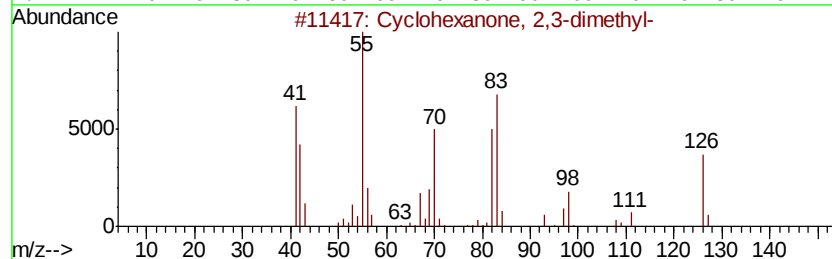
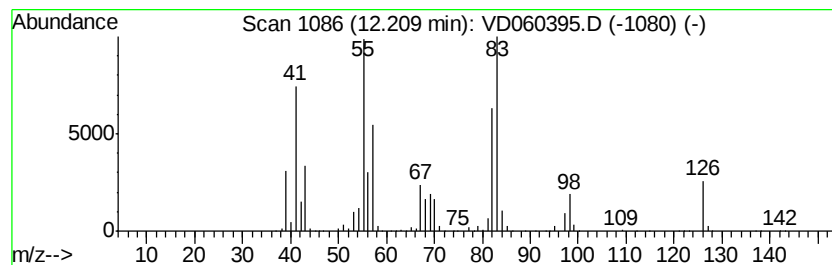
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Quant Title : SW846 8260

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

Peak Number 10 Cyclohexanone, 2,3-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.21	336.84 ug/l	26364500	Chlorobenzene-d5	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanone, 2,3-dimethyl-	126	C8H14O	013395-76-1	90
2		Cyclohexane, propyl-	126	C9H18	001678-92-8	83
3		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	81
4		Cyclohexane, methyl-	98	C7H14	000108-87-2	58
5		2,3-Dimethyl-2-heptene	126	C9H18	003074-64-4	58



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD111618\
Data File : VD060395.D
Acq On : 16 Nov 2018 13:49
Operator : JC/SY
Sample : J5911-06
Misc : 5.00g/5ml/MSVOA_D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
FDSB-07(20-22)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D110618S.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, 2-methyl-	8.99	141.1	ug/l	17099100	2	7.47	6061600	50.0
Heptane, 2,6-dime...	10.47	135.8	ug/l	10626900	3	11.30	3913480	50.0
Cyclohexane, ethyl-	10.72	135.3	ug/l	10586200	3	11.30	3913480	50.0
Cyclohexane, 1,1,...	10.80	171.7	ug/l	13437500	3	11.30	3913480	50.0
Octane, 2-methyl-	11.12	191.1	ug/l	14956000	3	11.30	3913480	50.0
Octane, 3-methyl-	11.23	203.1	ug/l	15894100	3	11.30	3913480	50.0
Cyclohexane, 1-et...	11.67	132.7	ug/l	10384200	3	11.30	3913480	50.0
unknown11.90	11.90	229.6	ug/l	17970600	3	11.30	3913480	50.0
Nonane, 3-methyl-	12.11	290.4	ug/l	22733000	3	11.30	3913480	50.0
Cyclohexanone, 2,...	12.21	336.8	ug/l	26364500	3	11.30	3913480	50.0