

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_F\DATA\VF111518\
 Data File : VF060755.D
 Acq On : 15 Nov 2018 19:14
 Operator : VA/AP
 Sample : J5911-01
 Misc : 5.00g/5mL/MSVOA-F/SOIL
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 FDSB-09(10-12)

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_F\METHODS\82F111318S.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	2.042	129	143	153	rBV	986580	5272626	5.92%	1.277%
2	2.220	153	161	173	rVB	575522	2725512	3.06%	0.660%
3	2.387	173	178	179	rBV	725597	1802849	2.02%	0.437%
4	2.989	222	239	246	rBV	1183913	8216571	9.22%	1.990%
5	3.068	246	247	281	rVB	803344	4149755	4.66%	1.005%
6	3.758	284	317	319	rBV2	2106719	10926486	12.26%	2.646%
7	4.715	408	414	435	rVB3	2843350	18708813	21.00%	4.531%
8	5.583	483	502	503	rBV	7718829	48051076	53.94%	11.637%
9	6.698	604	615	616	rBV	3296511	15778802	17.71%	3.821%
10	10.486	983	999	1006	rBV7	8640270	89089679	100.00%	21.576%
11	11.679	1117	1120	1122	rBV3	4060825	9361984	10.51%	2.267%
12	12.104	1154	1163	1167	rVB6	12840140	47861797	53.72%	11.591%
13	12.163	1167	1169	1172	rBV2	1364788	2745083	3.08%	0.665%
14	12.291	1178	1182	1186	rVB2	10270100	27531834	30.90%	6.668%
15	12.370	1186	1190	1194	rVV3	6048012	15875687	17.82%	3.845%
16	12.439	1194	1197	1199	rVV3	2507746	4047956	4.54%	0.980%
17	12.478	1199	1201	1202	rVV2	2553990	3819928	4.29%	0.925%
18	12.508	1202	1204	1208	rVB2	3732209	6915986	7.76%	1.675%
19	12.676	1216	1221	1224	rBV2	5651067	14674663	16.47%	3.554%
20	12.774	1228	1231	1235	rVV2	8356710	17956209	20.16%	4.349%
21	12.834	1235	1237	1240	rVV2	3403790	7328687	8.23%	1.775%
22	12.883	1240	1242	1246	rVV2	4148291	6960307	7.81%	1.686%
23	12.962	1247	1250	1255	rVB3	1606688	3349800	3.76%	0.811%
24	13.041	1255	1258	1261	rVB	2441081	4622670	5.19%	1.120%
25	13.100	1261	1264	1266	rBV	1830157	2856413	3.21%	0.692%
26	13.179	1270	1272	1275	rBV	4372346	7689980	8.63%	1.862%
27	13.258	1277	1280	1283	rVB2	5277704	9373832	10.52%	2.270%
28	13.307	1283	1285	1289	rVB2	1163180	1822518	2.05%	0.441%
29	13.406	1293	1295	1297	rVB	1256007	1442190	1.62%	0.349%
30	13.435	1297	1298	1303	rVB4	654477	1093671	1.23%	0.265%
31	13.514	1303	1306	1308	rVB	1342228	1898426	2.13%	0.460%
32	13.593	1311	1314	1318	rVB2	838166	1657435	1.86%	0.401%
33	13.721	1324	1327	1329	rBV	1848907	2822495	3.17%	0.684%
34	13.859	1339	1341	1346	rBV2	1821248	3467460	3.89%	0.840%

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Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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35	14.086	1358	1364	1367	rVB2	483450	1014666	1.14%	0.246%
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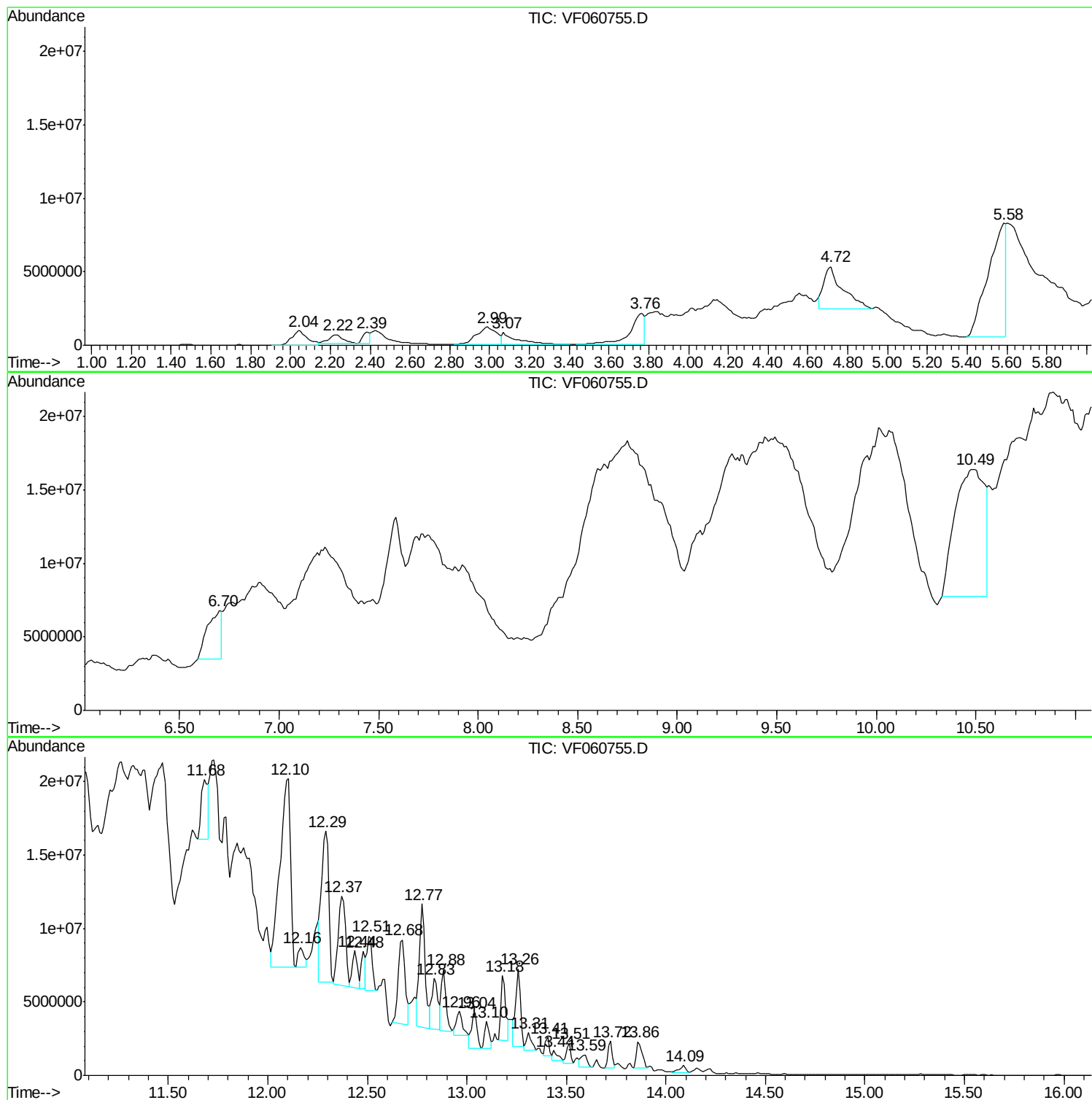
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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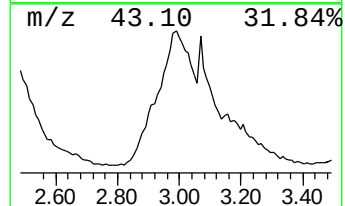
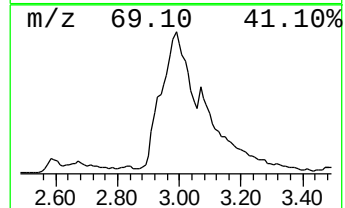
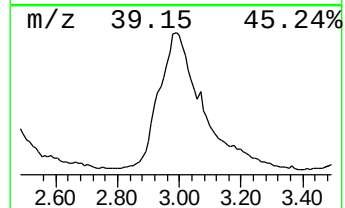
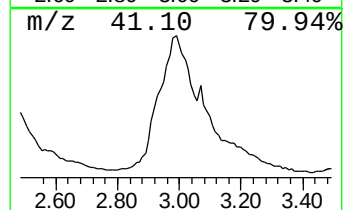
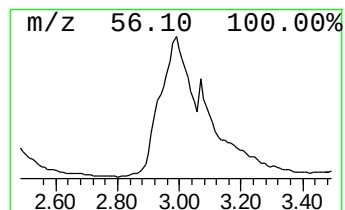
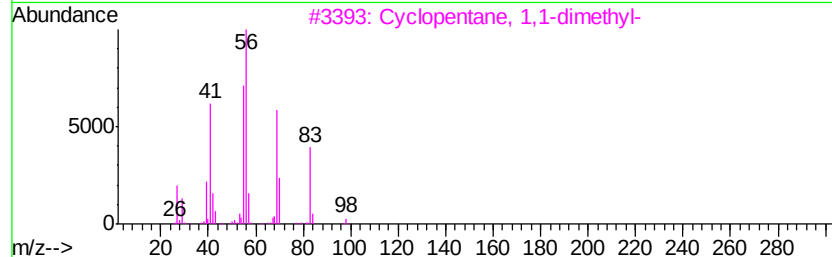
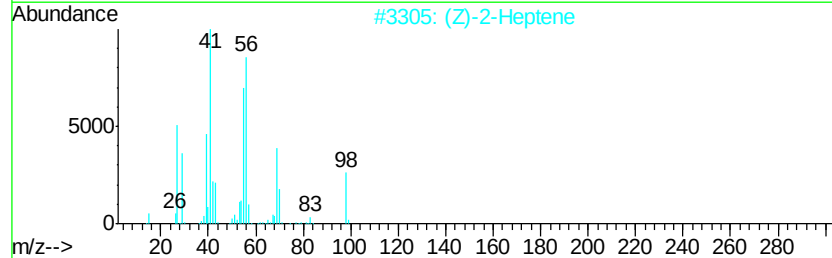
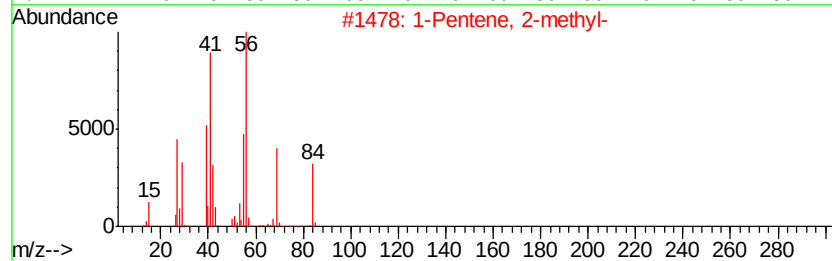
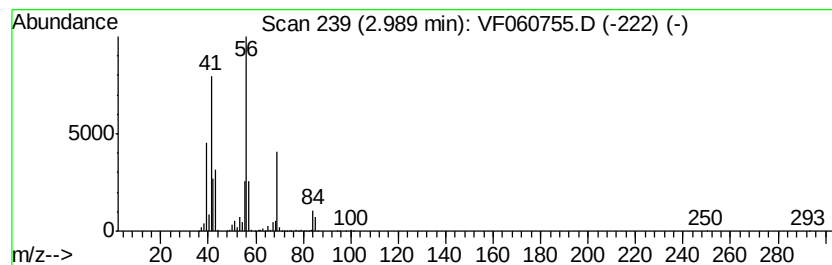
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Quant Title : SW846 8260

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

Peak Number 1 1-Pentene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
2.99	21.96 ug/l	8216570	Pentafluorobenzene	4.88		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentene, 2-methyl-		84	C6H12	000763-29-1	90
2	(Z)-2-Heptene		98	C7H14	006443-92-1	64
3	Cyclopentane, 1,1-dimethyl-		98	C7H14	001638-26-2	59
4	Cyclopentane, methyl-		84	C6H12	000096-37-7	59
5	Cyclohexane		84	C6H12	000110-82-7	53



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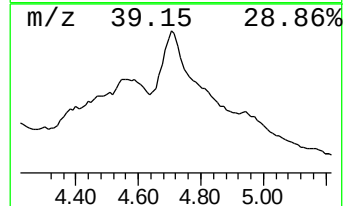
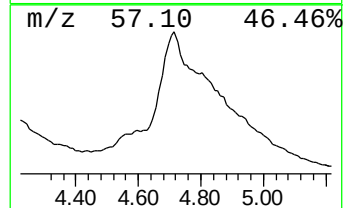
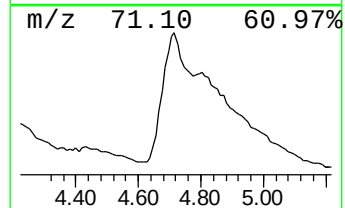
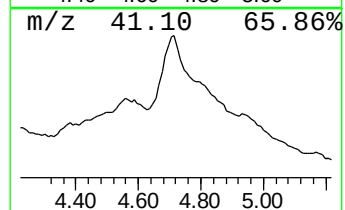
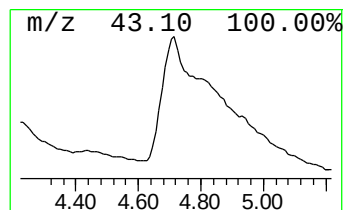
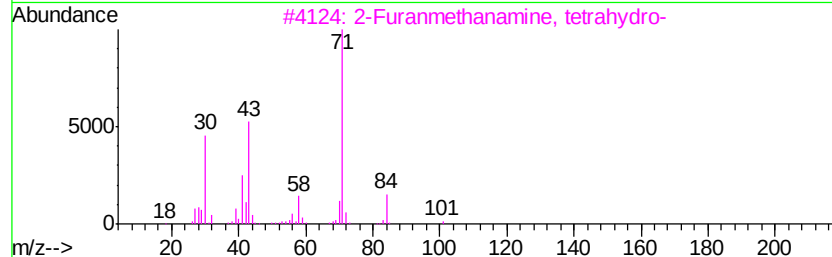
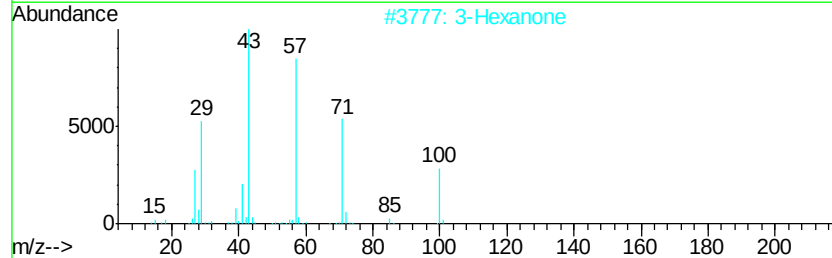
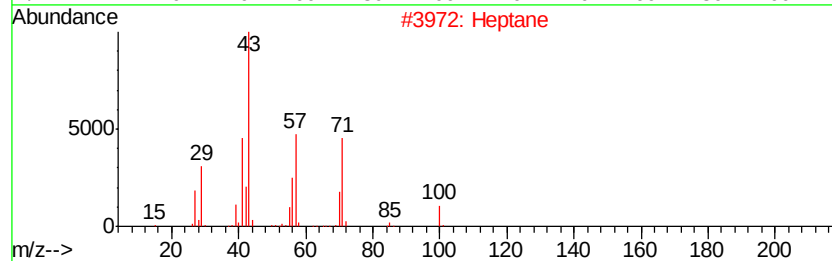
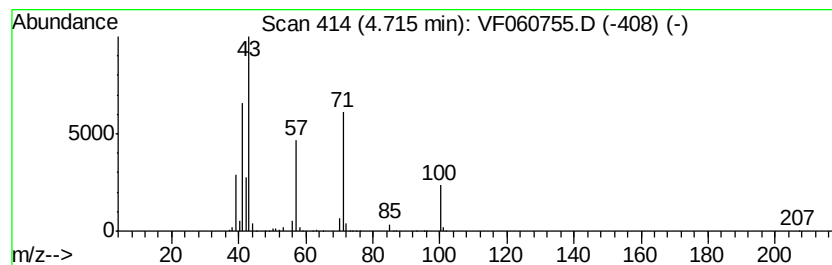
Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_F\METHODS\82F111318S.M
Quant Title : SW846 8260

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

Peak Number 2 Heptane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.72	50.00 ug/l	18708800	Pentafluorobenzene	4.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane	100	C7H16	000142-82-5	86
2		3-Hexanone	100	C6H12O	000589-38-8	42
3		2-Furanmethanamine, tetrahydro-	101	C5H11NO	004795-29-3	38
4		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	38
5		Pentane, 1-bromo-	150	C5H11Br	000110-53-2	36



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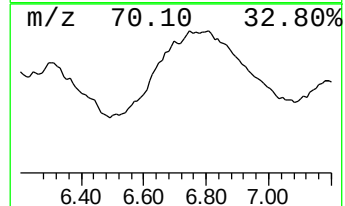
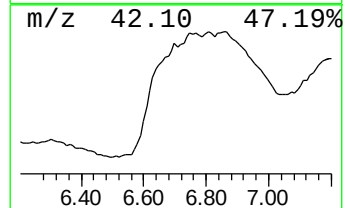
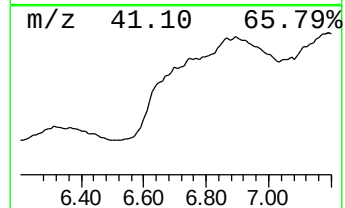
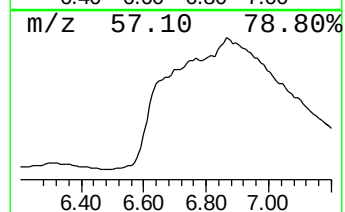
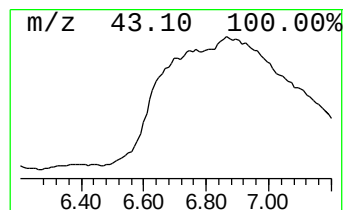
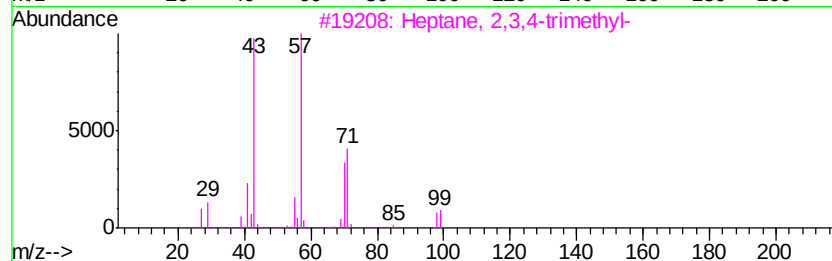
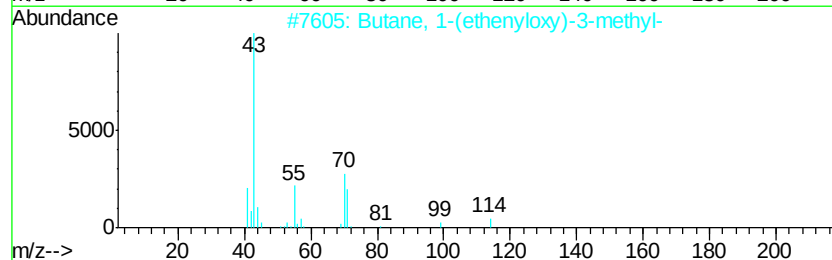
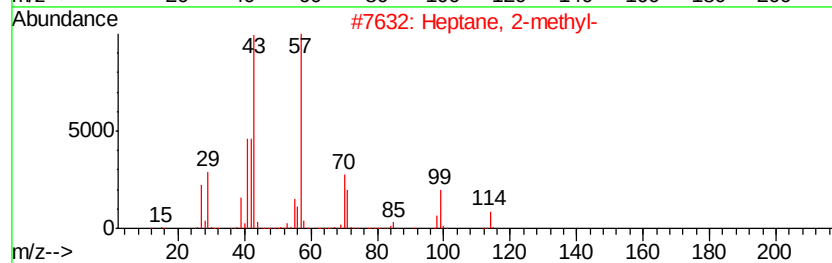
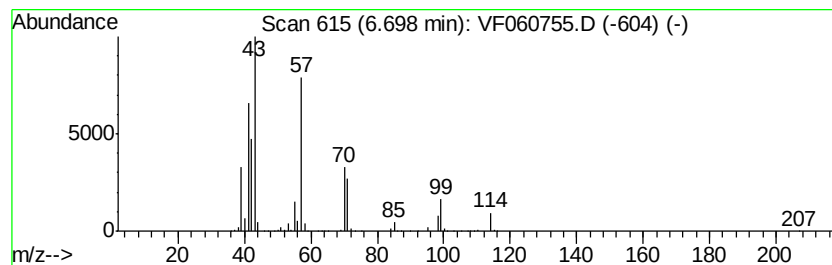
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Peak Number 3 Heptane, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.70	16.42 ug/l	15778800	1,4-Difluorobenzene	5.61
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptane, 2-methyl-	114 C8H18	000592-27-8	87
2	Butane, 1-(ethenyloxy)-3-methyl-	114 C7H14O	039782-38-2	43
3	Heptane, 2,3,4-trimethyl-	142 C10H22	052896-95-4	43
4	Isopentyloxyethyl acetate	174 C9H18O3	204652-53-9	43
5	Pentane, 2-bromo-	150 C5H11Br	000107-81-3	43



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ALS Vial : 20 Sample Multiplier: 1

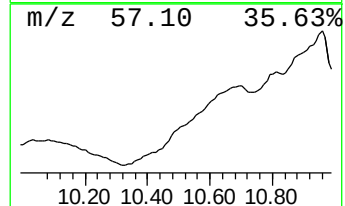
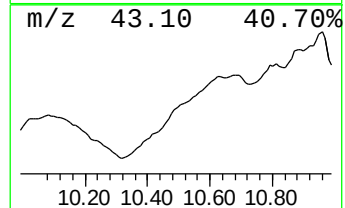
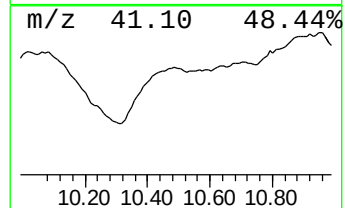
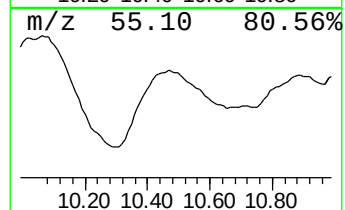
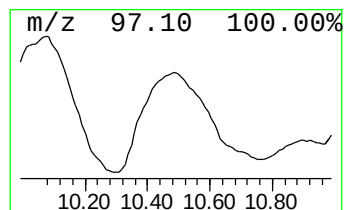
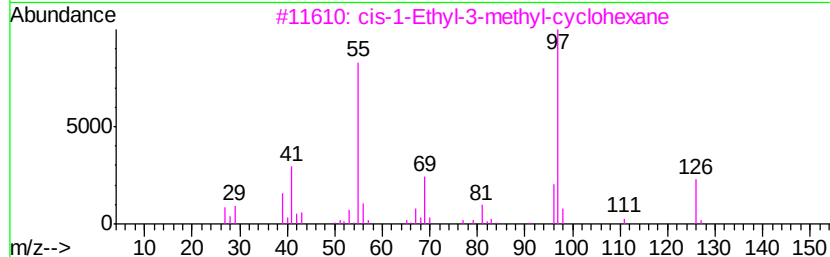
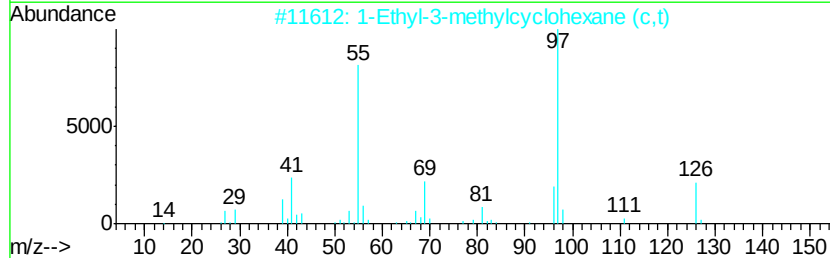
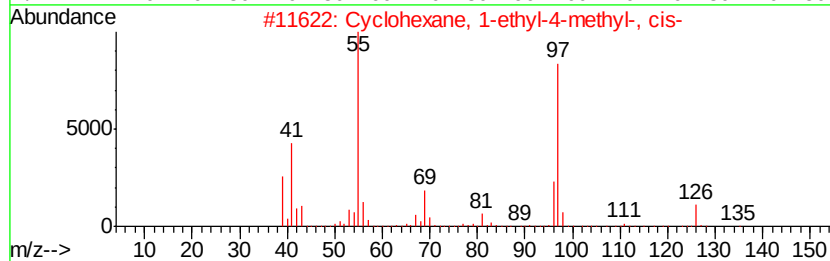
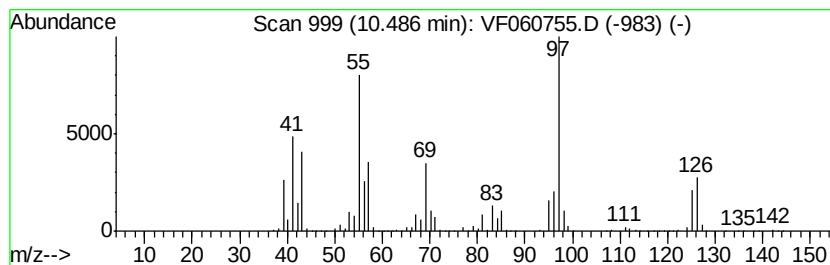
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Cyclohexane, 1-ethyl-4-meth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.49	50.00 ug/l	89089700	Chlorobenzene-d5	9.84		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	70
2		1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	68
3		cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	68
4		trans-1,2-Diethyl cyclopentane	126	C9H18	000932-40-1	59
5		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	53



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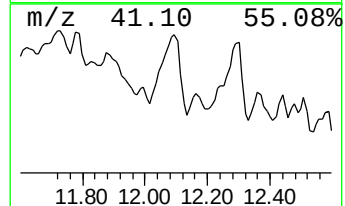
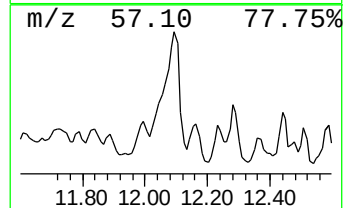
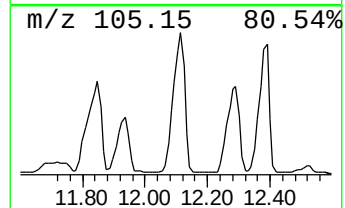
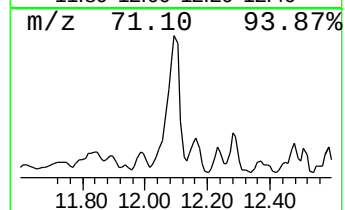
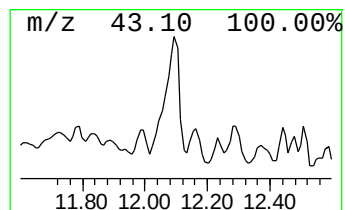
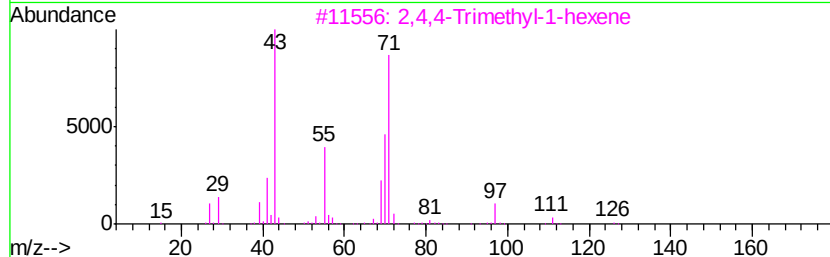
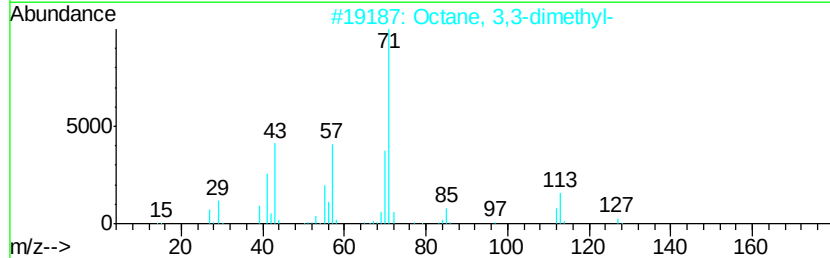
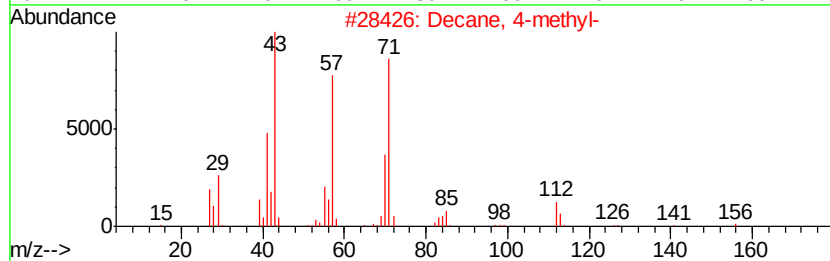
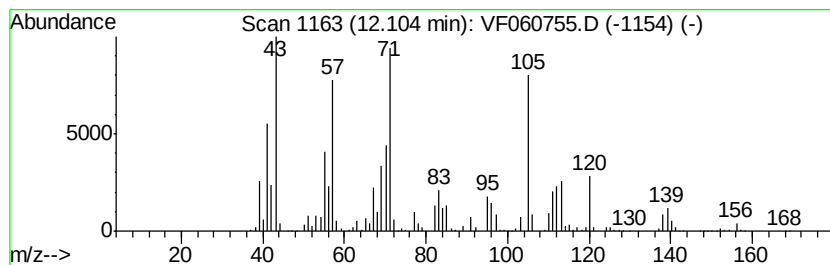
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Peak Number 5 Decane, 4-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.10	163.08 ug/l	47861800	1,4-Dichlorobenzene-d4	12.60
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Decane, 4-methyl-	156 C11H24	002847-72-5	60
2	Octane, 3,3-dimethyl-	142 C10H22	004110-44-5	49
3	2,4,4-Trimethyl-1-hexene	126 C9H18	051174-12-0	43
4	Dodecane, 4-methyl-	184 C13H28	006117-97-1	38
5	Octane, 3,6-dimethyl-	142 C10H22	015869-94-0	35



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF111518\
Data File : VF060755.D
Acq On : 15 Nov 2018 19:14
Operator : VA/AP
Sample : J5911-01
Misc : 5.00g/5mL/MSVOA-F/SOIL
ALS Vial : 20 Sample Multiplier: 1

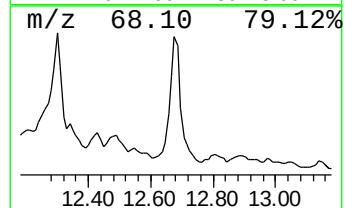
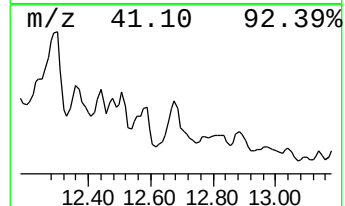
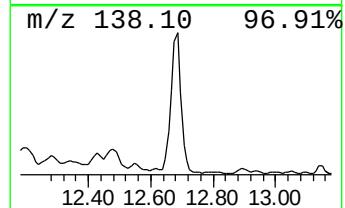
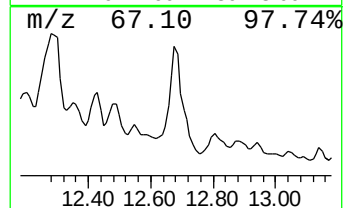
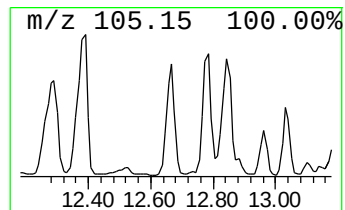
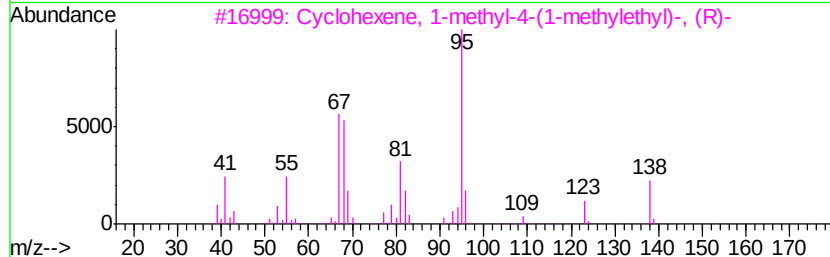
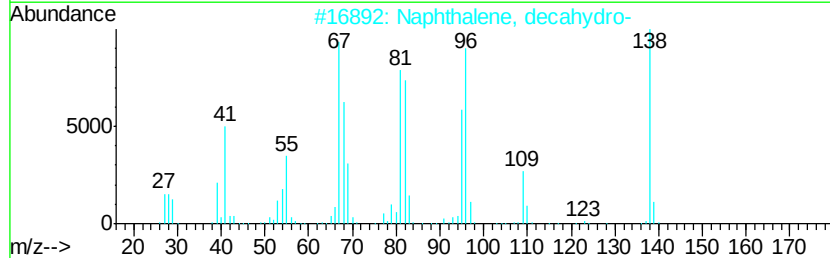
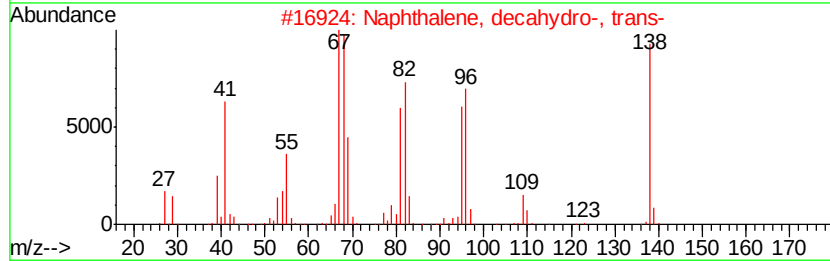
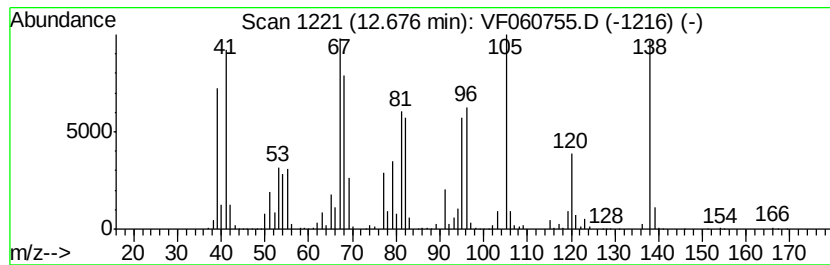
Instrument :
MSVOA_F
ClientSampleId :
FDSB-09(10-12)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F111318S.M
Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, decahydro-, tr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.68	50.00 ug/l	14674700	1,4-Dichlorobenzene-d4	12.60	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	91
2	Naphthalene, decahydro-	138	C10H18	000091-17-8	76
3	Cyclohexene, 1-methyl-4-(1-methy...	138	C10H18	001195-31-9	70
4	Spiro[4.4]nonan-2-one	138	C9H14O	034177-18-9	58
5	Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF111518\
Data File : VF060755.D
Acq On : 15 Nov 2018 19:14
Operator : VA/AP
Sample : J5911-01
Misc : 5.00g/5mL/MSVOA-F/SOIL
ALS Vial : 20 Sample Multiplier: 1

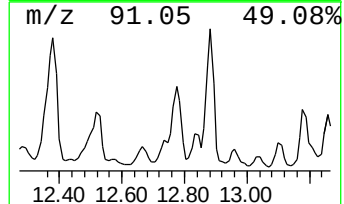
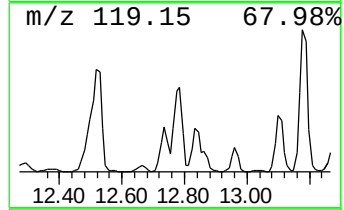
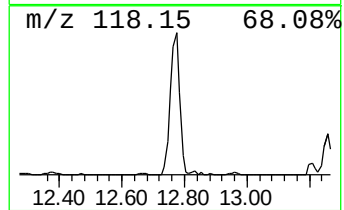
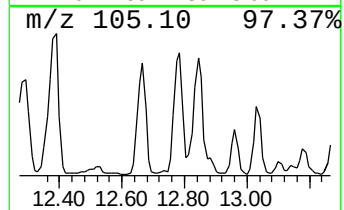
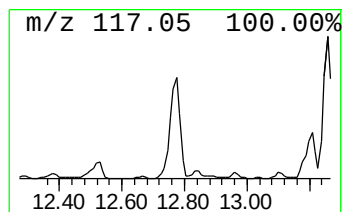
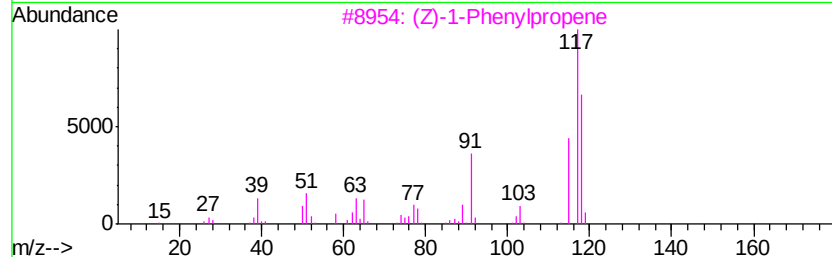
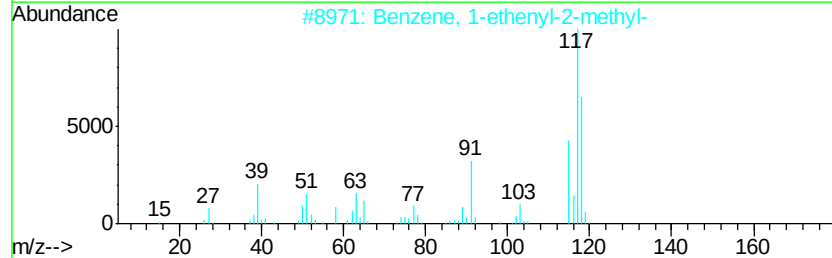
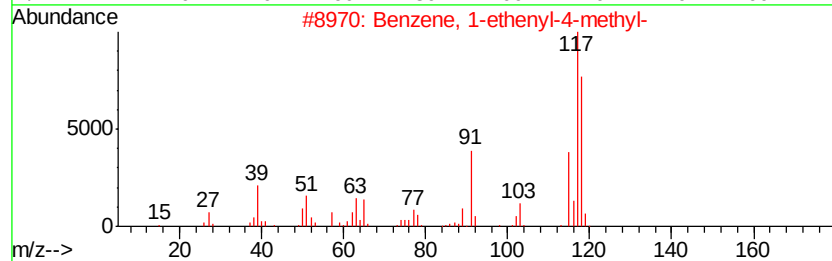
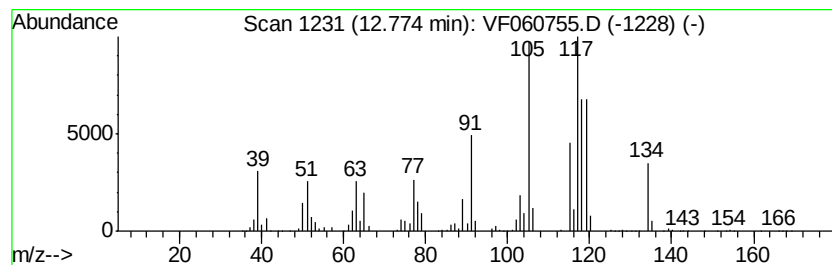
Instrument :
MSVOA_F
ClientSampleId :
FDSB-09(10-12)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F111318S.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethenyl-4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	61.18 ug/l	17956200	1,4-Dichlorobenzene-d4	12.60
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-4-methyl-	118 C9H10	000622-97-9 90
2		Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 55
3		(Z)-1-Phenylpropene	118 C9H10	000766-90-5 50
4		Benzene, 1-propenyl-	118 C9H10	000637-50-3 50
5		Indane	118 C9H10	000496-11-7 49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF111518\
Data File : VF060755.D
Acq On : 15 Nov 2018 19:14
Operator : VA/AP
Sample : J5911-01
Misc : 5.00g/5mL/MSVOA-F/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
FDSB-09(10-12)

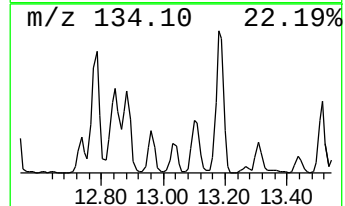
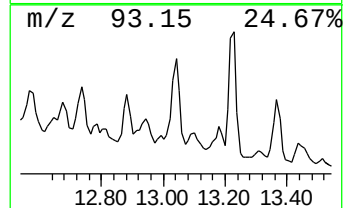
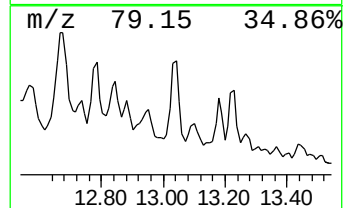
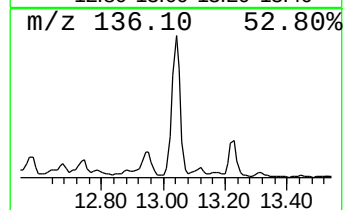
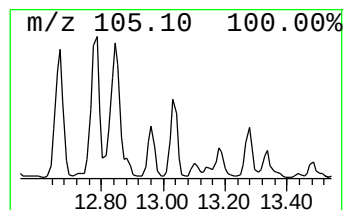
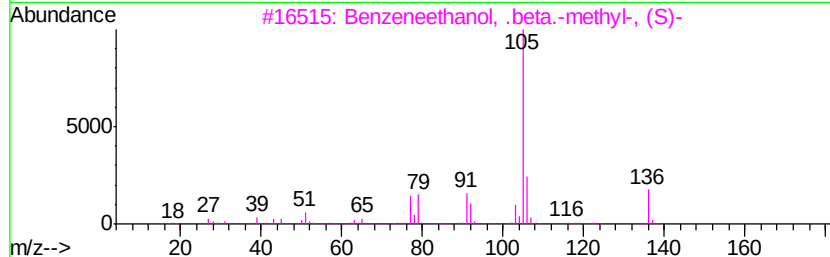
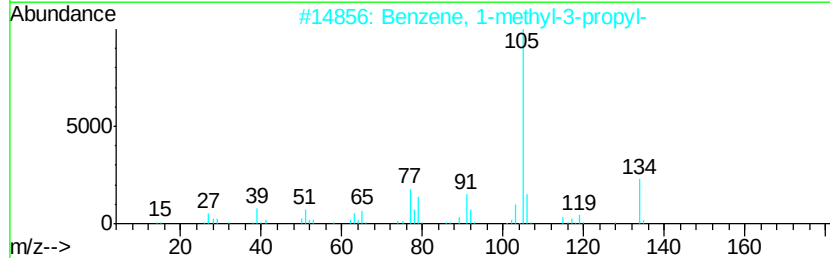
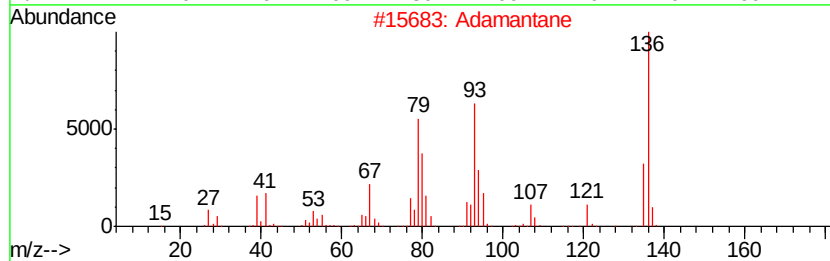
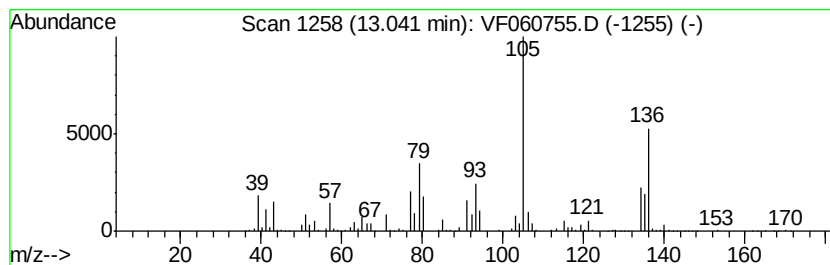
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F111318S.M
Quant Title : SW846 8260

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

Peak Number 8 unknown13.04 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.04	15.75 ug/l	4622670	1,4-Dichlorobenzene-d4	12.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Adamantane	136	C10H16	000281-23-2	46
2		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	38
3		Benzeneethanol, .beta.-methyl-, ...	136	C9H12O	037778-99-7	38
4		Benzeneethanol, 3-methyl-	136	C9H12O	001875-89-4	30
5		Benzeneethanol, 2-methyl-	136	C9H12O	019819-98-8	27



Data Path : Z:\V0ASRV\HPCHEM1\MSV0A_F\DATA\VF111518\
Data File : VF060755.D
Acq On : 15 Nov 2018 19:14
Operator : VA/AP
Sample : J5911-01
Misc : 5.00g/5mL/MSV0A-F/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSV0A_F
ClientSampleId :
FDSB-09(10-12)

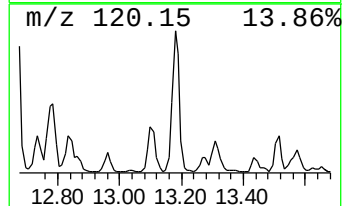
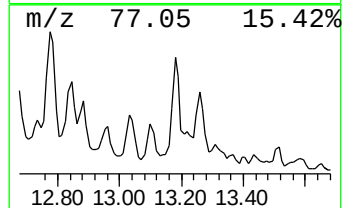
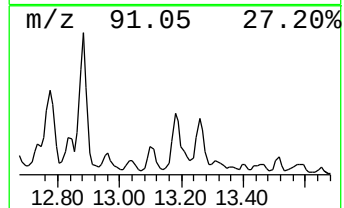
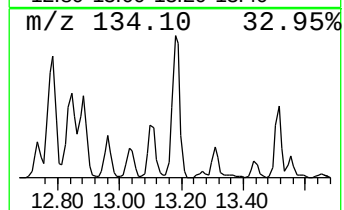
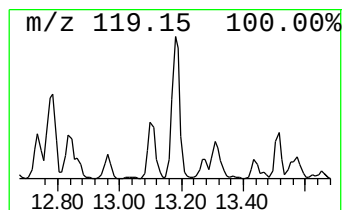
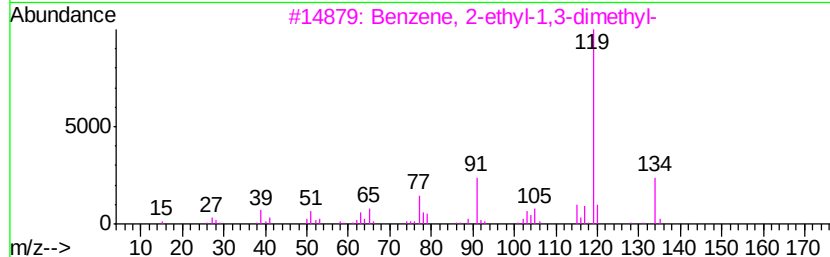
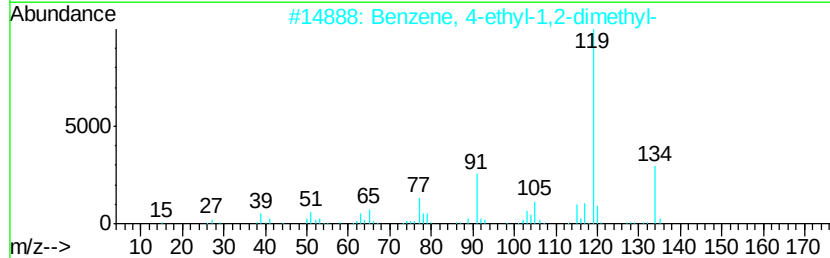
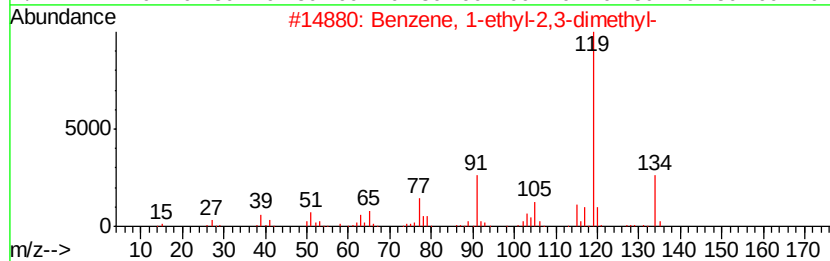
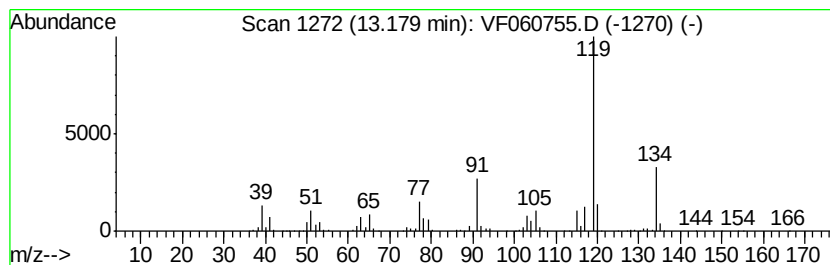
Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_F\METHODS\82F111318S.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.18	26.20 ug/l	7689980	1,4-Dichlorobenzene-d4	12.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	96
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	96
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95



Data Path : Z:\V0ASRV\HPCHEM1\MSV0A_F\DATA\VF111518\
Data File : VF060755.D
Acq On : 15 Nov 2018 19:14
Operator : VA/AP
Sample : J5911-01
Misc : 5.00g/5mL/MSV0A-F/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSV0A_F
ClientSampleId :
FDSB-09(10-12)

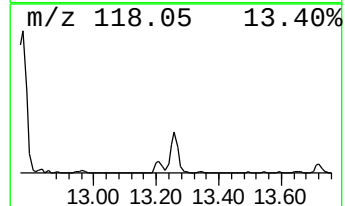
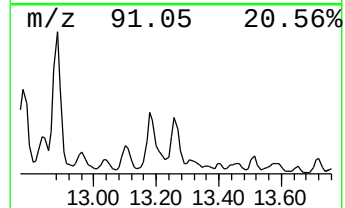
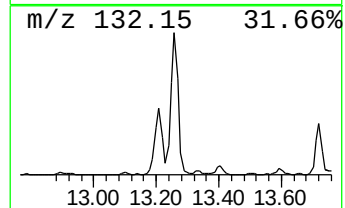
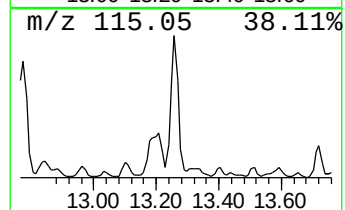
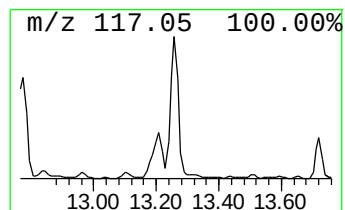
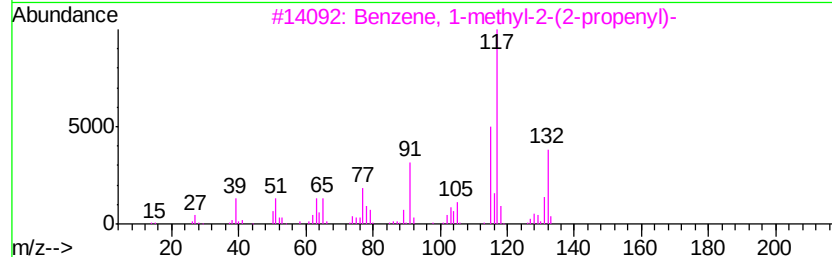
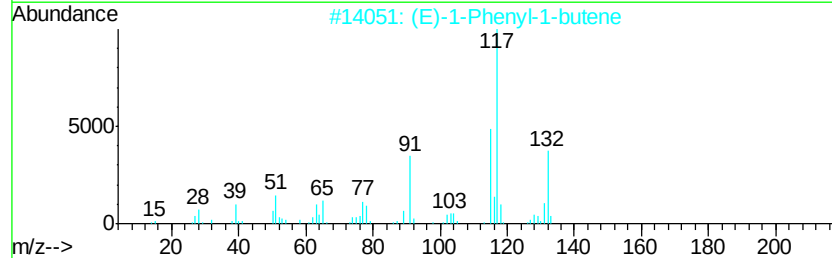
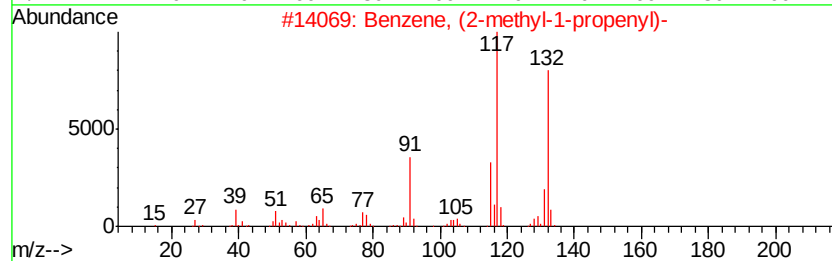
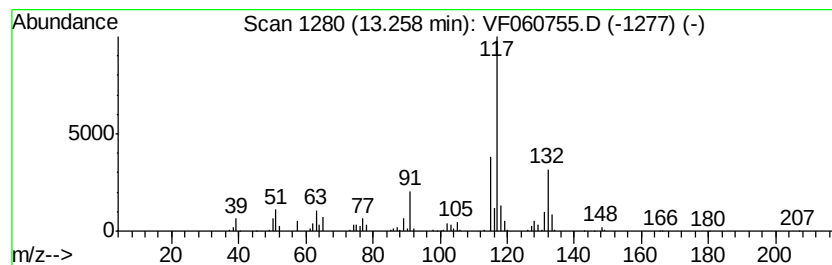
Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_F\METHODS\82F111318S.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, (2-methyl-1-propen... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.26	31.94 ug/l	9373830	1,4-Dichlorobenzene-d4	12.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	91
2		(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	90
3		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	86
4		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	80
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	72



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_F\DATA\VF111518\
 Data File : VF060755.D
 Acq On : 15 Nov 2018 19:14
 Operator : VA/AP
 Sample : J5911-01
 Misc : 5.00g/5mL/MSVOA-F/SOIL
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 FDSB-09(10-12)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F111318S.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
1-Pentene, 2-methyl-	2.99	22.0	ug/l	8216570	1	4.88	18708800	50.0
Heptane	4.72	50.0	ug/l	18708800	1	4.88	18708800	50.0
Heptane, 2-methyl-	6.70	16.4	ug/l	15778800	2	5.61	48051100	50.0
Cyclohexane, 1-et...	10.49	50.0	ug/l	89089700	3	9.84	89089700	50.0
Decane, 4-methyl-	12.10	163.1	ug/l	47861800	4	12.60	14674700	50.0
Naphthalene, deca...	12.68	50.0	ug/l	14674700	4	12.60	14674700	50.0
Benzene, 1-etheny...	12.77	61.2	ug/l	17956200	4	12.60	14674700	50.0
unknown13.04	13.04	15.8	ug/l	4622670	4	12.60	14674700	50.0
Benzene, 1-ethyl-...	13.18	26.2	ug/l	7689980	4	12.60	14674700	50.0
Benzene, (2-methy...	13.26	31.9	ug/l	9373830	4	12.60	14674700	50.0