

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF080718\
 Data File : VF059911.D
 Acq On : 7 Aug 2018 9:48
 Operator : VA/AP
 Sample : J4297-02
 Misc : 6.00µ/5mL/MSVOA F/SOIL
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 TP-2

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\82F080418S.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	5.040	449	457	468	rBV2	287425	960950	1.44%	0.141%
2	5.770	527	531	541	rVB	288075	790012	1.19%	0.116%
3	6.765	625	632	643	rVB2	346084	1335095	2.00%	0.196%
4	6.982	648	654	670	rVV	256525	859437	1.29%	0.126%
5	7.238	678	680	694	rVB3	215390	779911	1.17%	0.114%
6	7.642	716	721	726	rBV	1302895	3353977	5.03%	0.492%
7	8.441	796	802	809	rBV	607491	1851506	2.78%	0.271%
8	8.618	815	820	832	rVB	1063302	3544047	5.32%	0.519%
9	8.806	832	839	855	rBV2	694193	3980285	5.97%	0.583%
10	9.131	866	872	877	rBV2	843060	2693422	4.04%	0.395%
11	9.269	881	886	894	rVB2	1673720	5215453	7.83%	0.764%
12	9.446	894	904	910	rBV	1607389	4726606	7.09%	0.693%
13	9.703	923	930	937	rBV4	222940	920029	1.38%	0.135%
14	9.860	937	946	949	rBV	545090	2105942	3.16%	0.309%
15	9.929	949	953	957	rBV2	1372821	3728395	5.60%	0.546%
16	10.057	960	966	972	rVB	6504069	17596299	26.41%	2.579%
17	10.284	985	989	994	rVB	1652839	3581490	5.38%	0.525%
18	10.393	994	1000	1007	rBV3	1312287	5390333	8.09%	0.790%
19	10.501	1007	1011	1017	rVB3	689834	1868887	2.81%	0.274%
20	10.659	1019	1027	1035	rBV6	1604820	6922454	10.39%	1.015%
21	10.807	1037	1042	1050	rBV3	4749437	18891722	28.36%	2.769%
22	10.915	1050	1053	1054	rBV	746868	1353721	2.03%	0.198%
23	10.954	1055	1057	1061	rVB	1358849	2628178	3.94%	0.385%
24	11.083	1066	1070	1074	rVB2	982050	2203759	3.31%	0.323%
25	11.181	1074	1080	1082	rBV3	1665702	5524545	8.29%	0.810%
26	11.240	1082	1086	1088	rBV2	2503152	5214729	7.83%	0.764%
27	11.280	1088	1090	1095	rVB	3436157	6825194	10.24%	1.000%
28	11.398	1095	1102	1107	rVB3	5740631	15110271	22.68%	2.215%
29	11.516	1110	1114	1115	rBV3	809347	1614053	2.42%	0.237%
30	11.566	1116	1119	1121	rBV3	511381	881849	1.32%	0.129%
31	11.615	1121	1124	1127	rVB4	619342	981465	1.47%	0.144%
32	11.674	1127	1130	1135	rBV2	3115746	10586283	15.89%	1.552%
33	11.792	1136	1142	1145	rBV	15527938	37197865	55.83%	5.452%
34	11.891	1150	1152	1154	rVB	870868	1073872	1.61%	0.157%

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Integrator: RTE
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35	11.940	1154	1157	1160	rVB	5389545	9140884	13.72%	1.340%
36	11.999	1160	1163	1166	rBV	1609676	3212573	4.82%	0.471%
37	12.068	1166	1170	1171	rBV2	2190660	4956823	7.44%	0.726%
38	12.118	1171	1175	1179	rVB2	11247872	25056226	37.61%	3.672%
39	12.187	1179	1182	1186	rVB2	2384096	4912961	7.37%	0.720%
40	12.265	1186	1190	1191	rBV2	3632626	6919013	10.39%	1.014%
41	12.315	1191	1195	1199	rVB	13425167	26023171	39.06%	3.814%
42	12.403	1200	1204	1206	rBV3	3358656	6351735	9.53%	0.931%
43	12.492	1209	1213	1215	rBV	3947895	8883393	13.33%	1.302%
44	12.532	1215	1217	1218	rVV	5400232	8465781	12.71%	1.241%
45	12.561	1218	1220	1226	rVV3	6095312	16079266	24.13%	2.357%
46	12.650	1226	1229	1232	rVB	5905337	10573391	15.87%	1.550%
47	12.709	1232	1235	1239	rBV2	9360928	17302307	25.97%	2.536%
48	12.827	1244	1247	1249	rBV3	2554396	5628699	8.45%	0.825%
49	12.877	1249	1252	1253	rBV	5067126	9483847	14.24%	1.390%
50	12.946	1253	1259	1264	rVB5	21528545	66622338	100.00%	9.764%
51	13.064	1264	1271	1273	rBV6	3545738	11597825	17.41%	1.700%
52	13.113	1273	1276	1280	rVB2	6478381	13731182	20.61%	2.012%
53	13.202	1280	1285	1286	rBV	5790452	17447412	26.19%	2.557%
54	13.251	1289	1290	1294	rVB2	4361728	5978280	8.97%	0.876%
55	13.320	1294	1297	1302	rVB4	5240707	14021165	21.05%	2.055%
56	13.389	1302	1304	1307	rBV3	5766409	11705076	17.57%	1.716%
57	13.438	1307	1309	1311	rVB2	3547068	5006774	7.52%	0.734%
58	13.488	1311	1314	1315	rBV2	5187142	7141506	10.72%	1.047%
59	13.517	1315	1317	1319	rVB	6598940	7962671	11.95%	1.167%
60	13.586	1321	1324	1326	rVB2	7335625	9778248	14.68%	1.433%
61	13.636	1326	1329	1331	rBV	6374039	10977915	16.48%	1.609%
62	13.803	1342	1346	1349	rBV2	16843492	30847714	46.30%	4.521%
63	13.902	1354	1356	1358	rBV	3431116	4873982	7.32%	0.714%
64	13.941	1358	1360	1364	rVV4	10923787	24653834	37.01%	3.613%
65	14.000	1364	1366	1369	rVB3	5201895	8493995	12.75%	1.245%
66	14.060	1369	1372	1374	rVB2	2708195	5118724	7.68%	0.750%
67	14.168	1377	1383	1385	rBV2	3713106	9712433	14.58%	1.423%
68	14.237	1385	1390	1393	rBV4	7026022	16854679	25.30%	2.470%
69	14.306	1393	1397	1402	rVB4	4922710	11917161	17.89%	1.747%
70	14.375	1402	1404	1407	rBV4	1387538	1794473	2.69%	0.263%
71	14.434	1408	1410	1415	rVB4	2091671	4158473	6.24%	0.609%

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72	14.533	1415	1420	1423	rBV5	3840848	9618860	14.44%	1.410%
73	14.631	1428	1430	1432	rBV2	1200269	1604470	2.41%	0.235%
74	14.671	1432	1434	1437	rVB2	2351523	3358484	5.04%	0.492%
75	14.720	1437	1439	1441	rVB3	511105	742205	1.11%	0.109%
76	14.819	1446	1449	1451	rVB3	2457786	3811637	5.72%	0.559%
77	14.947	1459	1462	1467	rVB3	960293	2853815	4.28%	0.418%
78	15.016	1467	1469	1472	rVB3	498195	764017	1.15%	0.112%
79	15.075	1472	1475	1480	rVB2	1047070	1738539	2.61%	0.255%
80	15.183	1483	1486	1489	rVB2	684005	1015794	1.52%	0.149%
81	15.242	1489	1492	1494	rBV2	457414	698636	1.05%	0.102%
82	15.361	1499	1504	1508	rBV	3787638	6055153	9.09%	0.887%
83	15.489	1514	1517	1524	rVB2	2563232	4368402	6.56%	0.640%

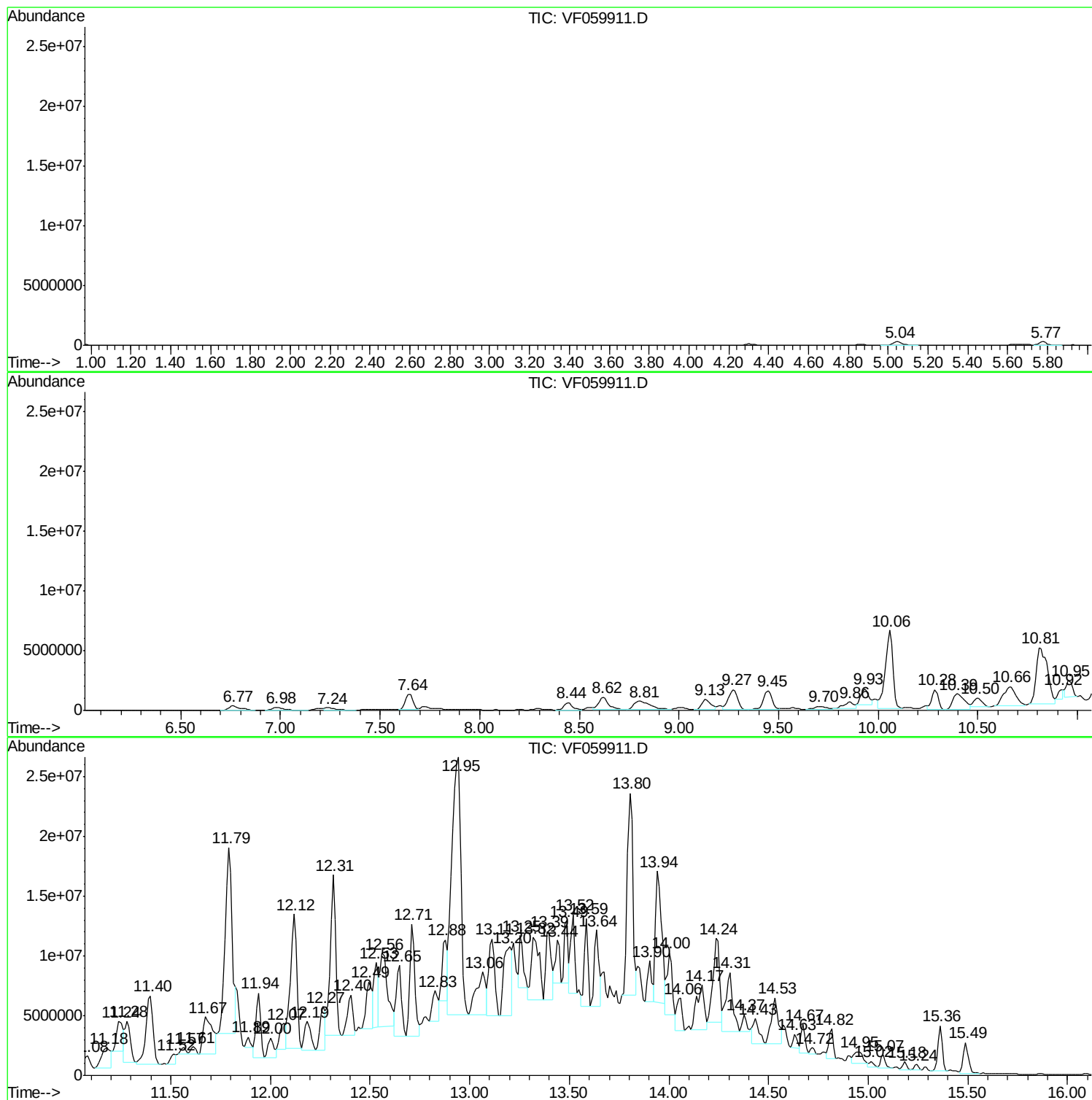
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Quant Title : SW846 8260

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



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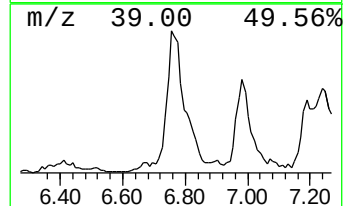
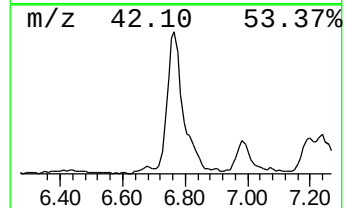
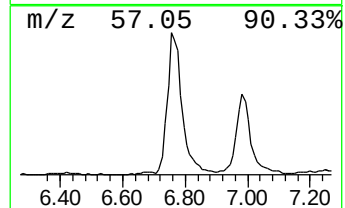
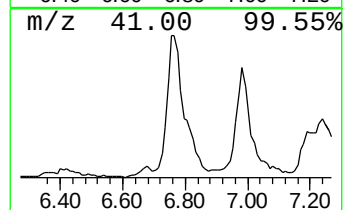
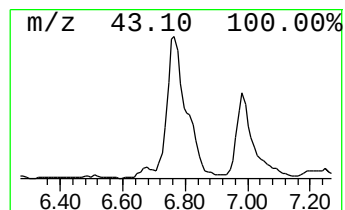
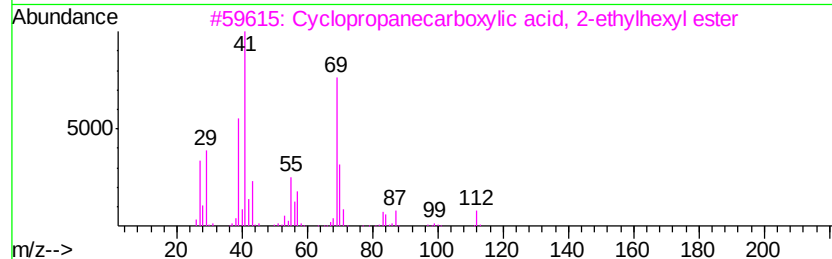
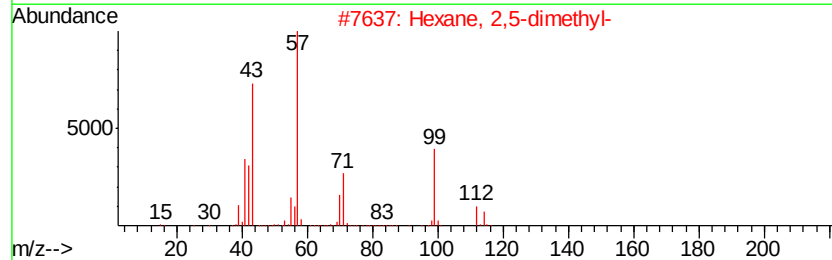
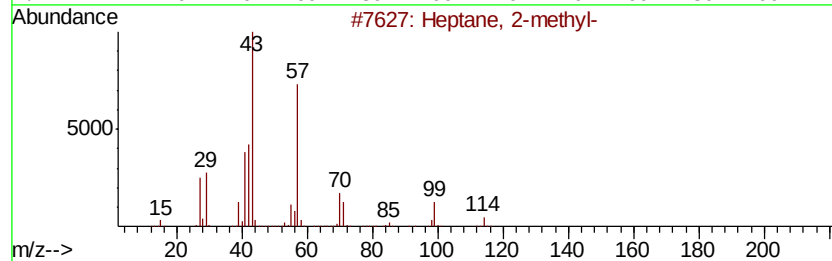
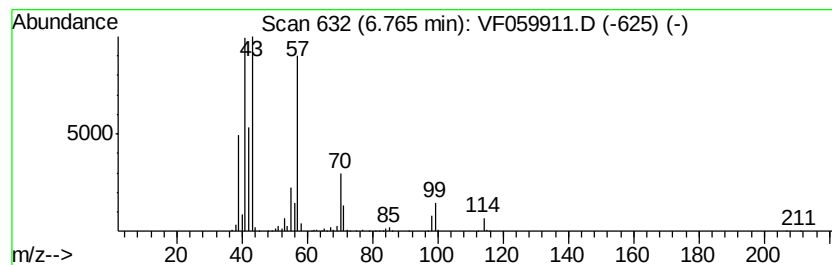
Instrument :
MSVOA_F
ClientSampleId :
TP-2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F080418S.M
Quant Title : SW846 8260

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

Peak Number 1 Heptane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
6.77	84.50 ug/l	1335100	1,4-Difluorobenzene	5.78	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2-methyl-	114	C8H18	000592-27-8	58
2	Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	49
3	Cyclopropanecarboxylic acid, 2-ethyl-	198	C12H22O2	103604-31-5	38
4	2-Pentenoic acid, 2-methyl-	114	C6H10O2	003142-72-1	38
5	Heptane, 2,3,4-trimethyl-	142	C10H22	052896-95-4	27



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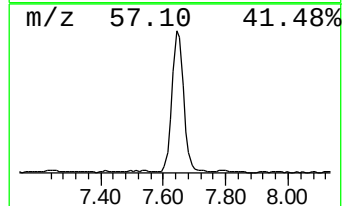
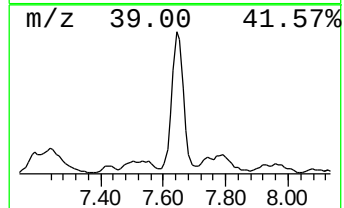
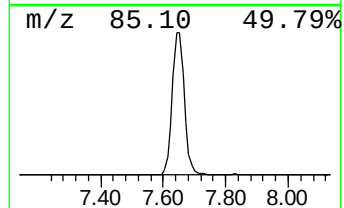
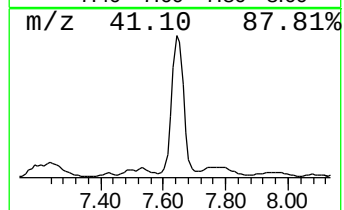
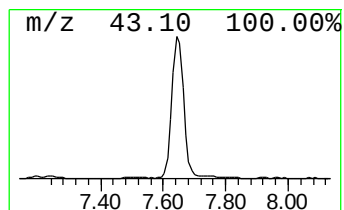
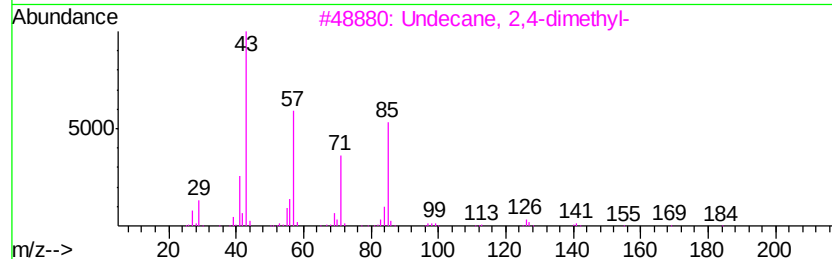
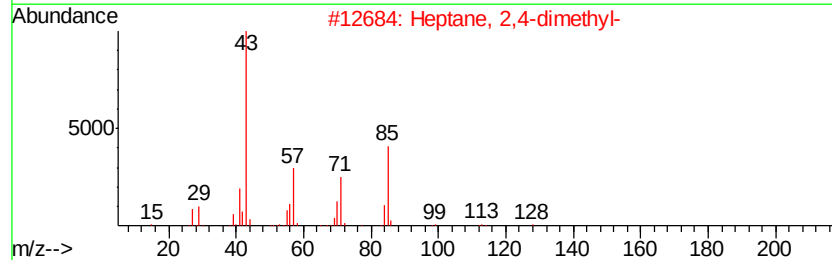
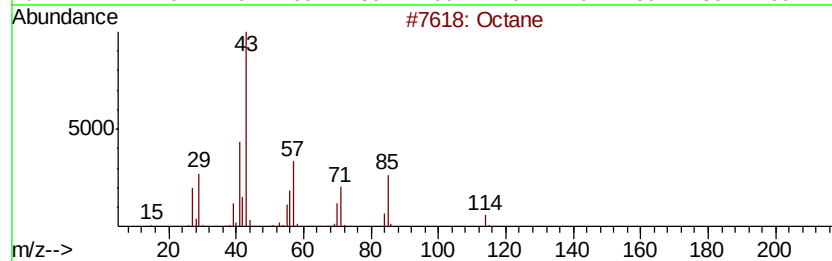
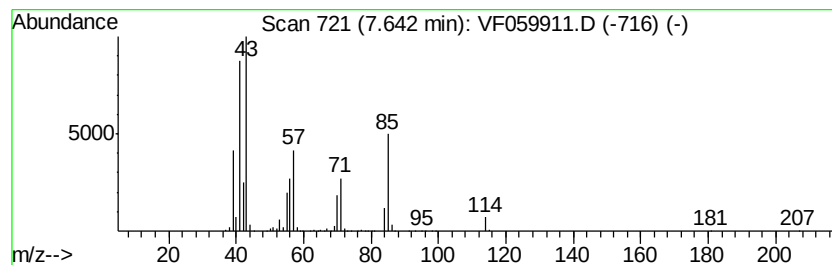
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Octane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	212.27 ug/l	3353980	1,4-Difluorobenzene	5.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane	114	C8H18	000111-65-9	64
2		Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	47
3		Undecane, 2,4-dimethyl-	184	C13H28	017312-80-0	43
4		Hexane, 2-methyl-	100	C7H16	000591-76-4	43
5		Sulfurous acid, hexyl pentadecyl...	376	C21H44O3S	1000309-13-7	43



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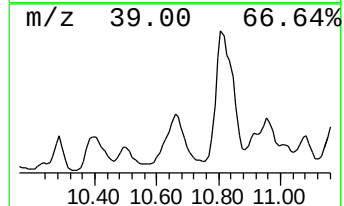
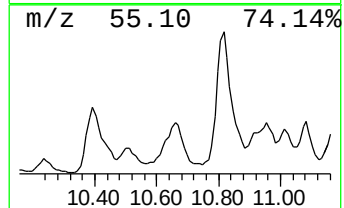
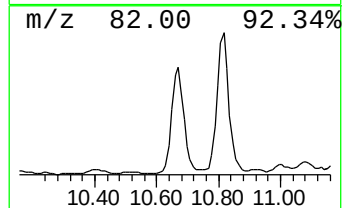
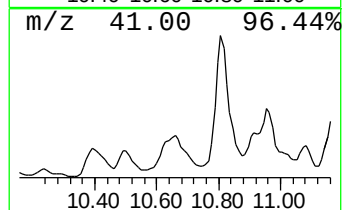
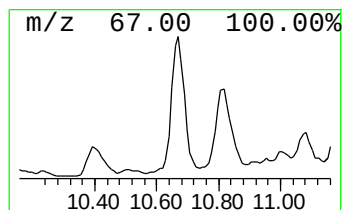
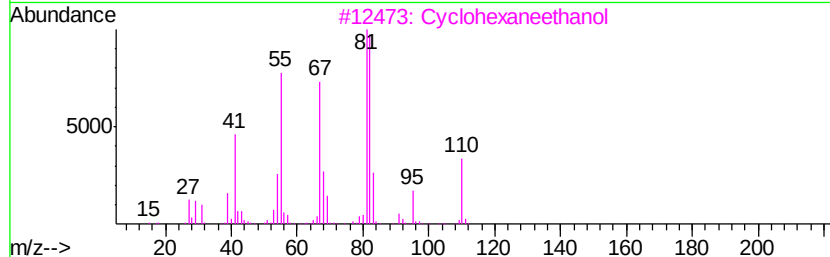
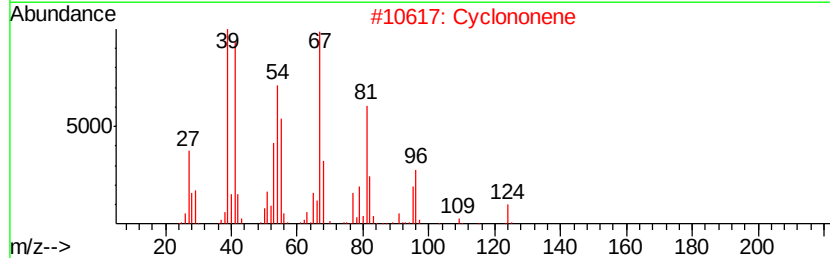
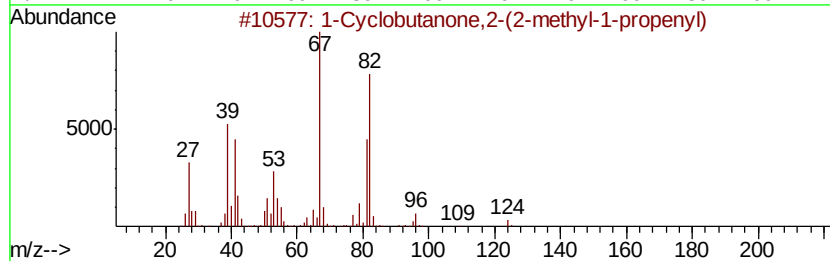
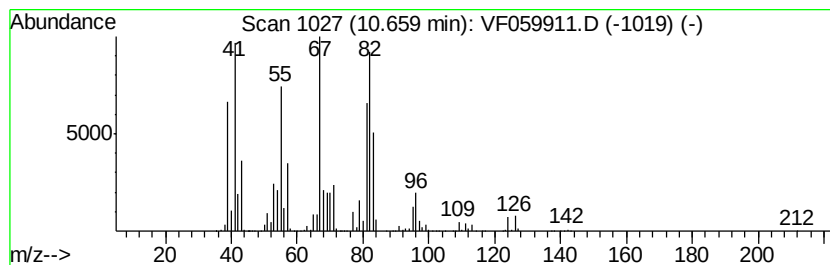
Instrument :
MSVOA_F
ClientSampled :
TP-2

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Quant Title : SW846 8260

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

Peak Number 3 1-Cyclobutanone,2-(2-methyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.66	92.83 ug/l	6922450	Chlorobenzene-d5	9.93		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Cyclobutanone,2-(2-methyl-1-pr...	124	C8H12O	091531-45-2	53
2		Cyclononene	124	C9H16	003618-11-9	47
3		Cyclohexaneethanol	128	C8H16O	004442-79-9	47
4		3-Hexyne	82	C6H10	000928-49-4	46
5		2,4-Hexadiene, (E,Z)-	82	C6H10	005194-50-3	43



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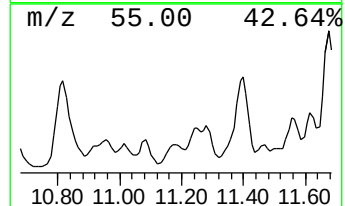
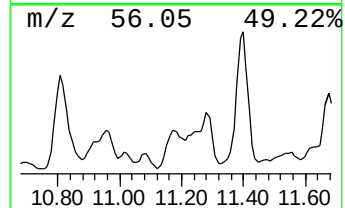
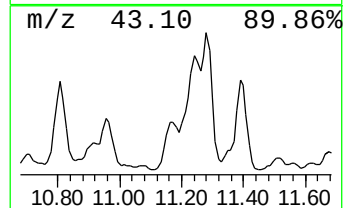
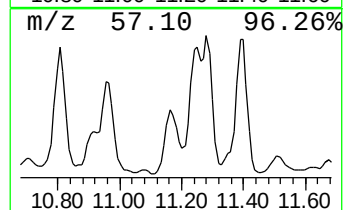
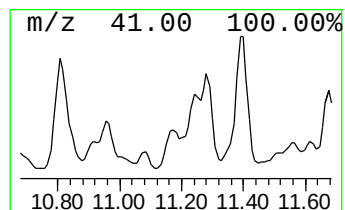
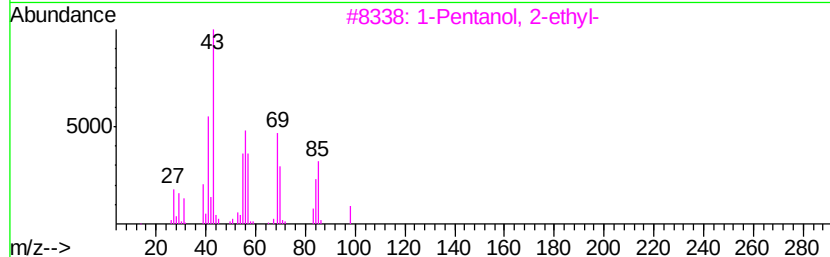
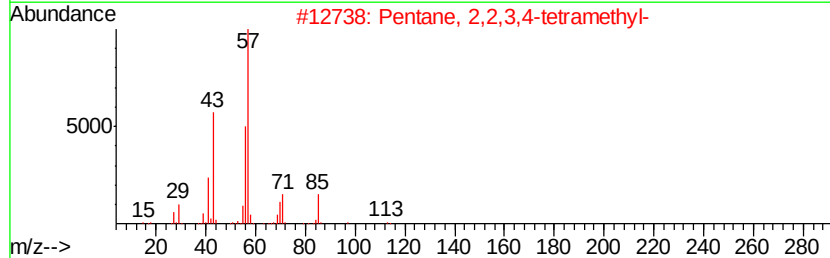
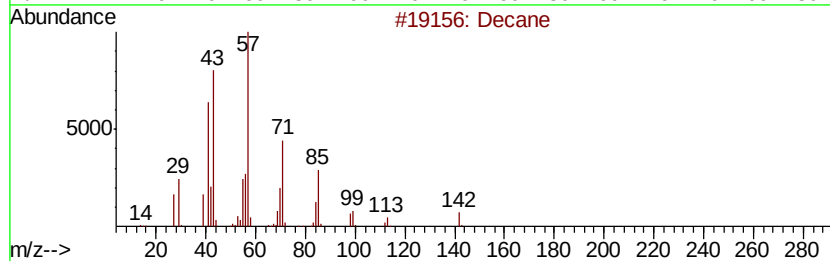
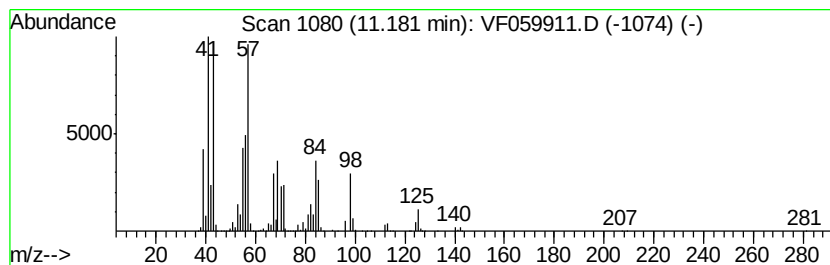
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Quant Title : SW846 8260

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

Peak Number 4 unknown11.18 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.18	74.09 ug/l	5524550	Chlorobenzene-d5	9.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decane	142	C10H22	000124-18-5	43
2		Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	38
3		1-Pentanol, 2-ethyl-	116	C7H16O	027522-11-8	38
4		2,2,6,6-Tetramethylheptane	156	C11H24	040117-45-1	35
5		1-Octene, 3,4-dimethyl-	140	C10H20	056728-11-1	35



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF080718\
Data File : VF059911.D
Acq On : 7 Aug 2018 9:48
Operator : VA/AP
Sample : J4297-02
Misc : 6.00µ/5mL/MSVOA F/SOIL
ALS Vial : 37 Sample Multiplier: 1

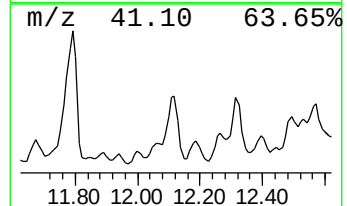
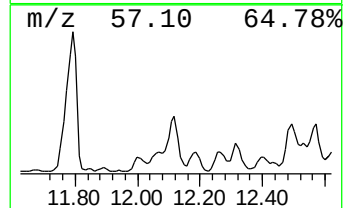
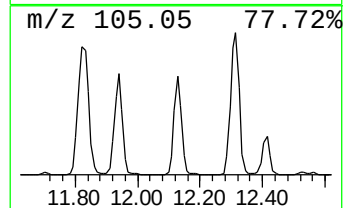
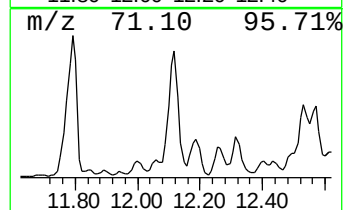
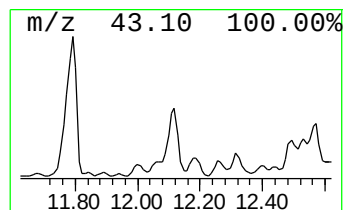
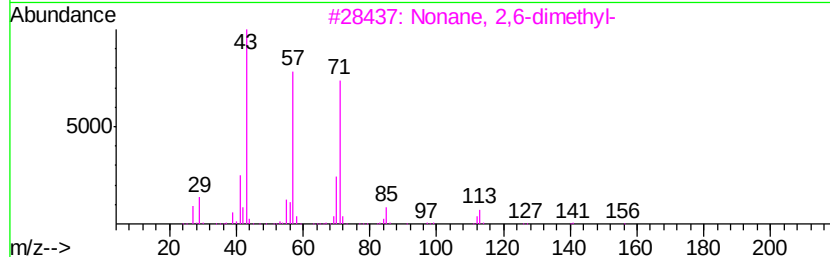
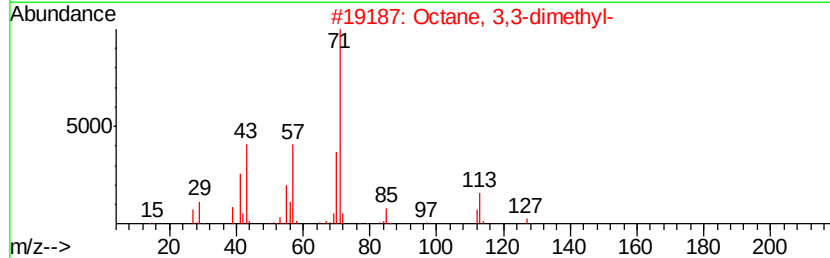
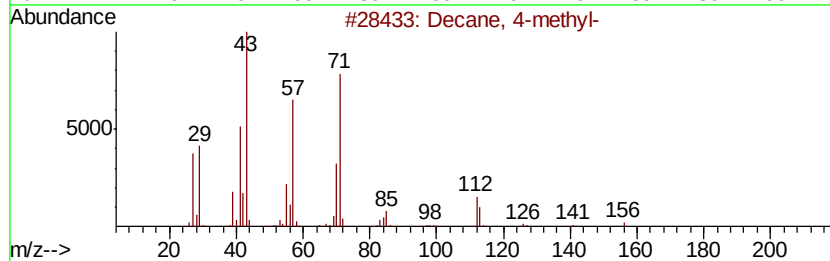
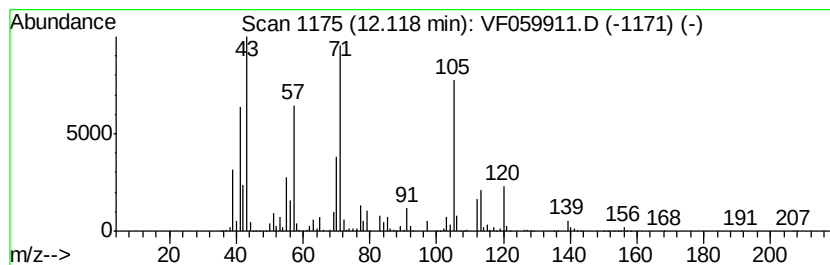
Instrument :
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ClientSampleId :
TP-2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F080418S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.12	118.49 ug/l	25056200	1,4-Dichlorobenzene-d4	12.65
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Decane, 4-methyl-	156 C11H24	002847-72-5	70
2	Octane, 3,3-dimethyl-	142 C10H22	004110-44-5	58
3	Nonane, 2,6-dimethyl-	156 C11H24	017302-28-2	53
4	Heptane, 3,3,5-trimethyl-	142 C10H22	007154-80-5	50
5	Sulfurous acid, octyl 2-pentyl e...	264 C13H28O3S	1000309-15-7	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF080718\
Data File : VF059911.D
Acq On : 7 Aug 2018 9:48
Operator : VA/AP
Sample : J4297-02
Misc : 6.00a/5mL/MSVOA F/SOIL
ALS Vial : 37 Sample Multiplier: 1

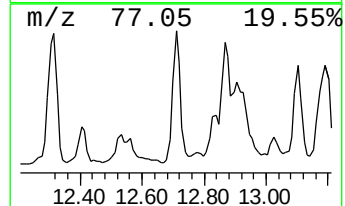
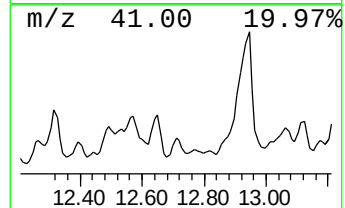
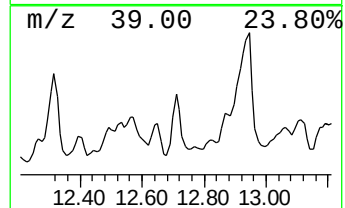
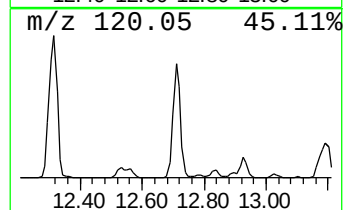
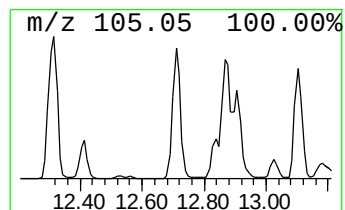
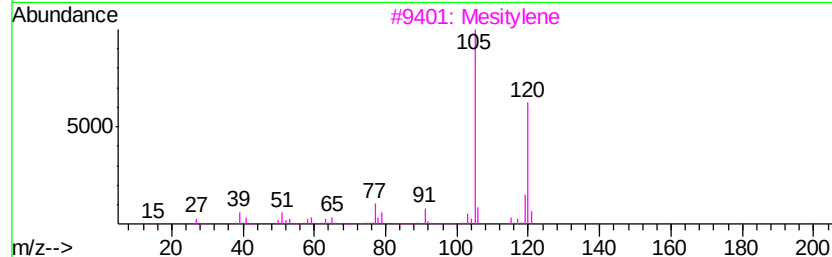
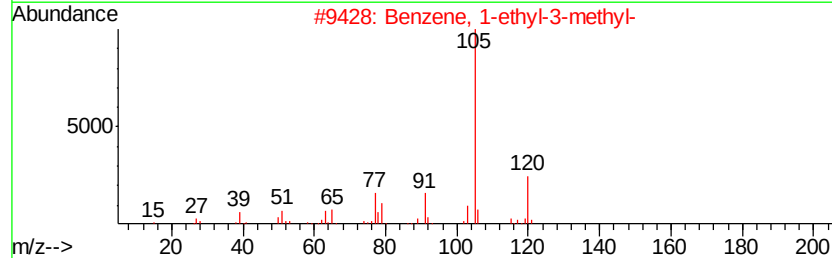
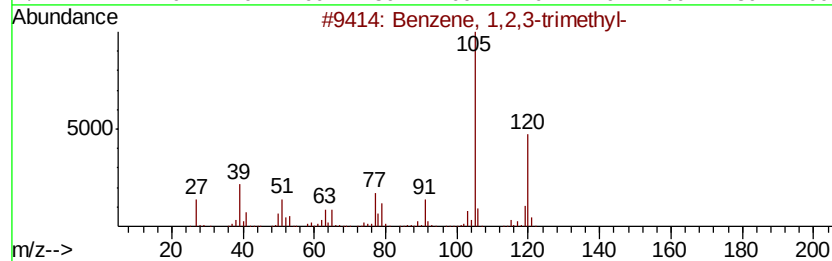
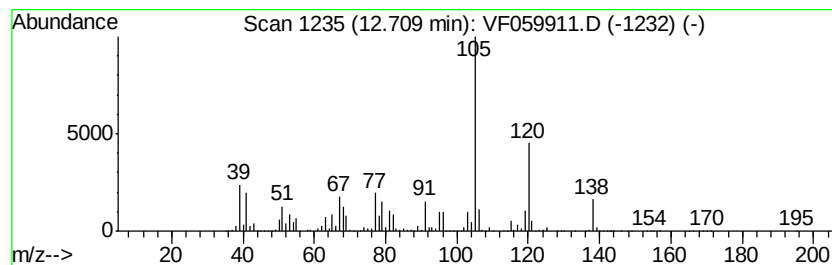
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F080418S.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.71	81.82 ug/l	17302300	1,4-Dichlorobenzene-d4	12.65
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2,3-trimethyl-	120 C9H12	000526-73-8 94
2		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 90
3		Mesitylene	120 C9H12	000108-67-8 90
4		Benzene, 1,2,4-trimethyl-	120 C9H12	000095-63-6 86
5		Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8 76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF080718\
Data File : VF059911.D
Acq On : 7 Aug 2018 9:48
Operator : VA/AP
Sample : J4297-02
Misc : 6.00µ/5mL/MSVOA F/SOIL
ALS Vial : 37 Sample Multiplier: 1

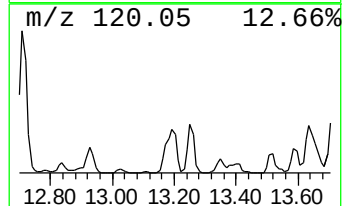
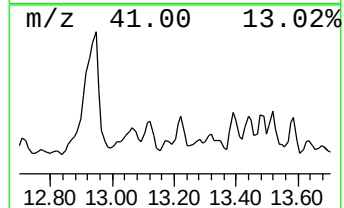
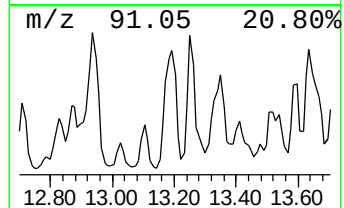
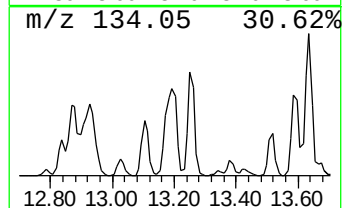
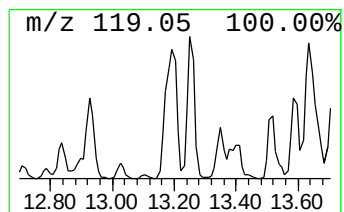
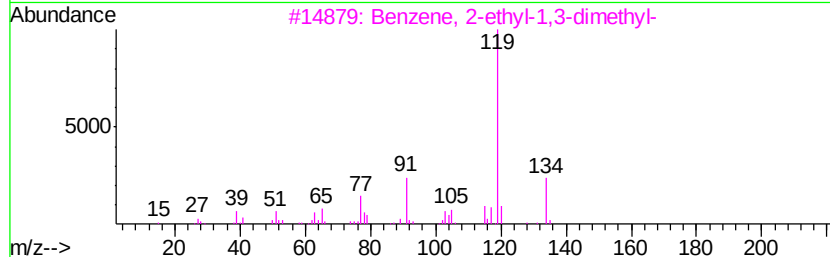
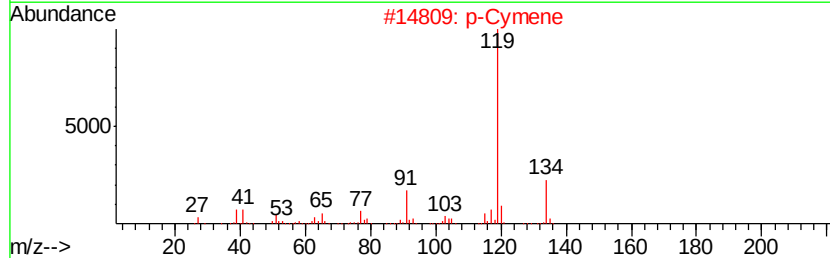
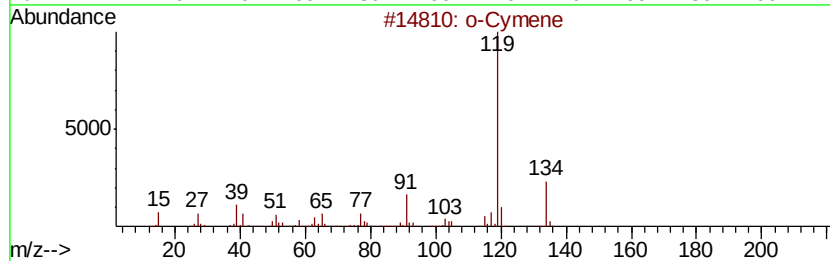
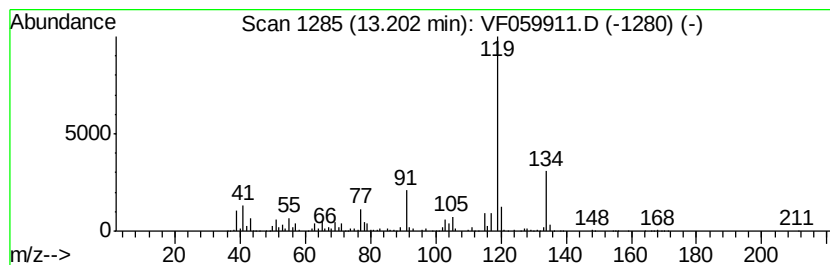
Instrument :
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ClientSampleId :
TP-2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F080418S.M
Quant Title : SW846 8260

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

Peak Number 7 o-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.20	82.51 ug/l	17447400	1,4-Dichlorobenzene-d4	12.65
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		o-Cymene	134 C10H14	000527-84-4 95
2		p-Cymene	134 C10H14	000099-87-6 95
3		Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4 94
4		Benzene, 2-ethyl-1,4-dimethyl-	134 C10H14	001758-88-9 94
5		Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5 91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF080718\
Data File : VF059911.D
Acq On : 7 Aug 2018 9:48
Operator : VA/AP
Sample : J4297-02
Misc : 6.00µ/5mL/MSVOA F/SOIL
ALS Vial : 37 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
TP-2

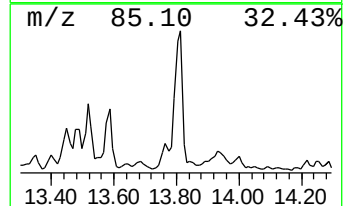
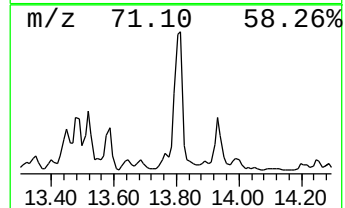
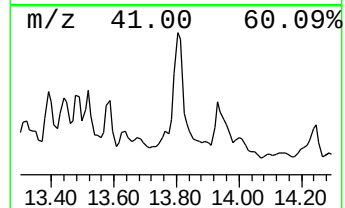
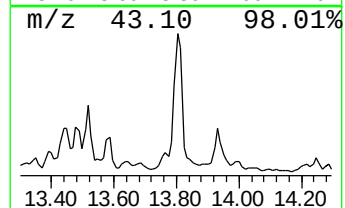
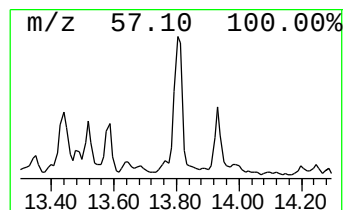
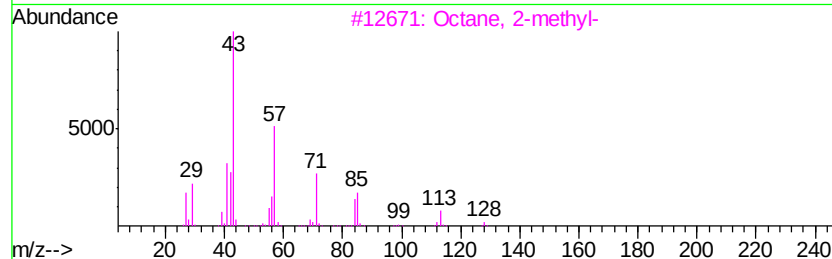
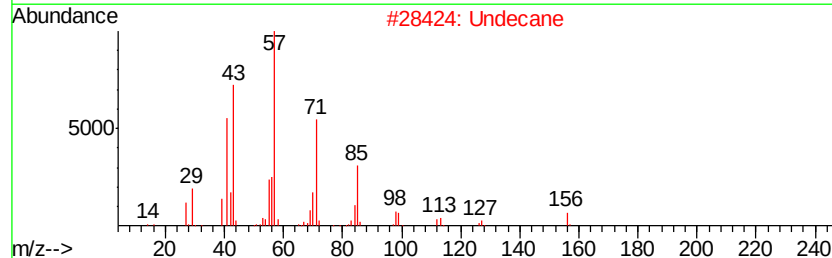
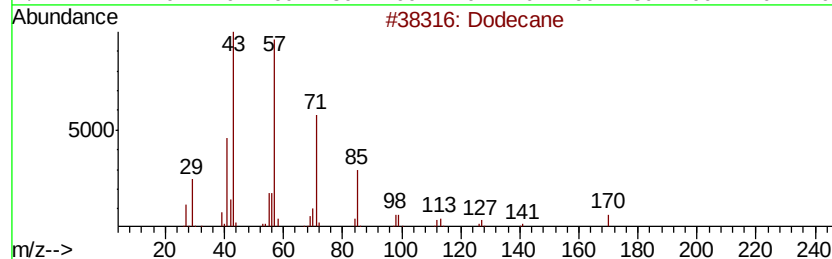
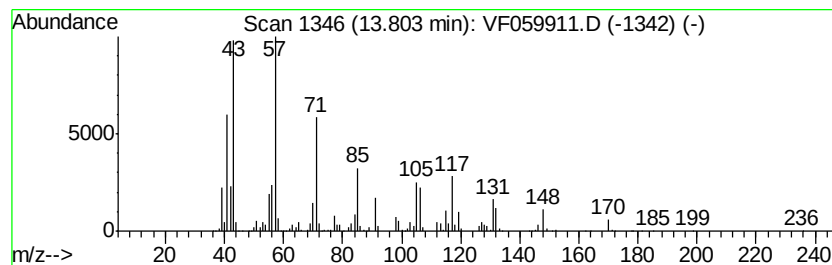
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Quant Title : SW846 8260

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

Peak Number 8 Dodecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.80	145.87 ug/l	30847700	1,4-Dichlorobenzene-d4	12.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane	170	C12H26	000112-40-3	83
2		Undecane	156	C11H24	001120-21-4	25
3		Octane, 2-methyl-	128	C9H20	003221-61-2	25
4		Tridecane	184	C13H28	000629-50-5	25
5		Undecane, 2-methyl-	170	C12H26	007045-71-8	22



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF080718\
Data File : VF059911.D
Acq On : 7 Aug 2018 9:48
Operator : VA/AP
Sample : J4297-02
Misc : 6.00µ/5mL/MSVOA F/SOIL
ALS Vial : 37 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampled :
TP-2

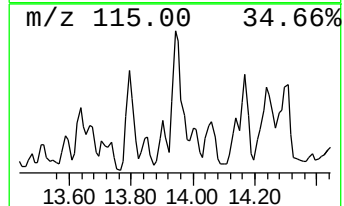
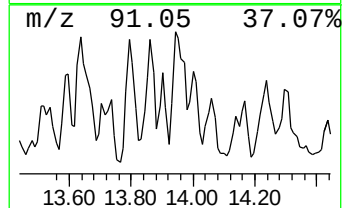
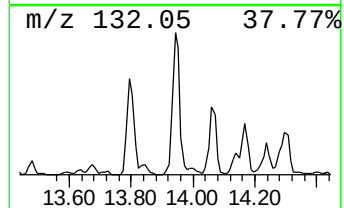
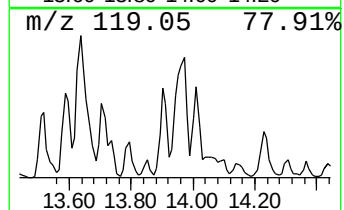
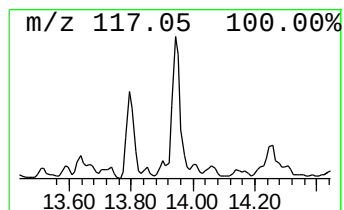
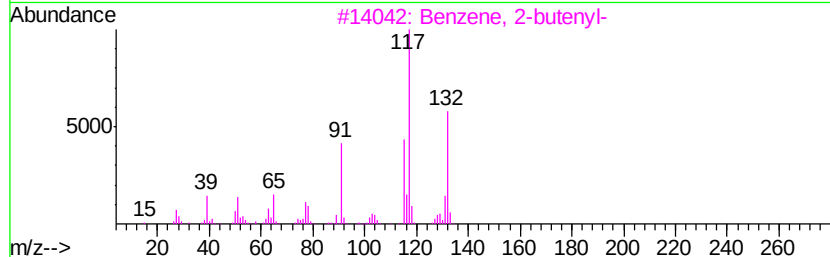
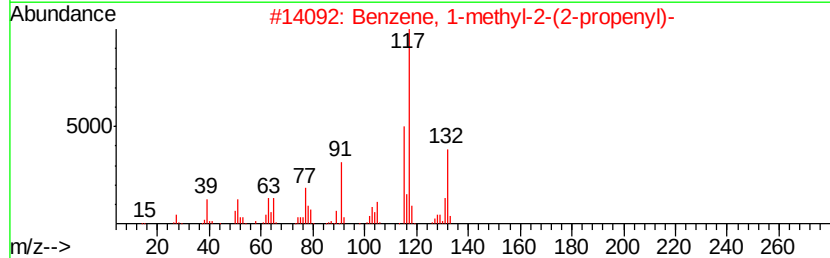
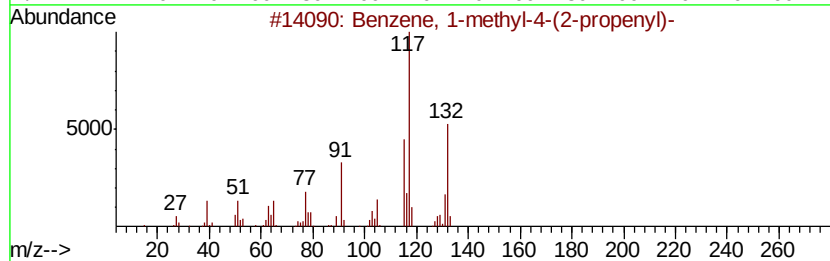
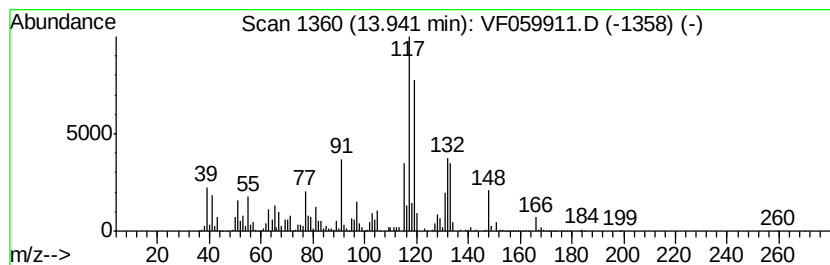
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-methyl-4-(2-prop... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.94	116.58 ug/l	24653800	1,4-Dichlorobenzene-d4	12.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	89
2			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	89
3			Benzene, 2-butenyl-	132	C10H12	001560-06-1	86
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	55
5			Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	55



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF080718\
Data File : VF059911.D
Acq On : 7 Aug 2018 9:48
Operator : VA/AP
Sample : J4297-02
Misc : 6.00µ/5mL/MSVOA F/SOIL
ALS Vial : 37 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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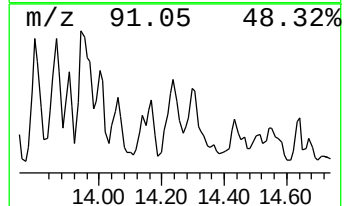
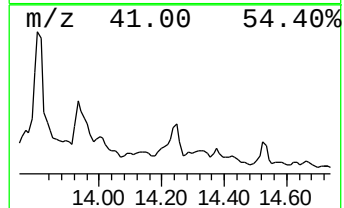
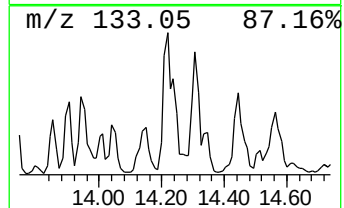
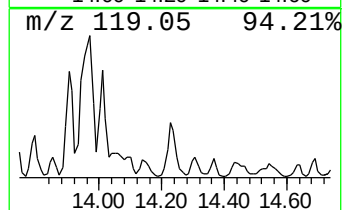
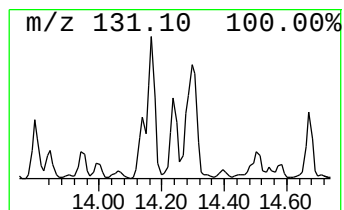
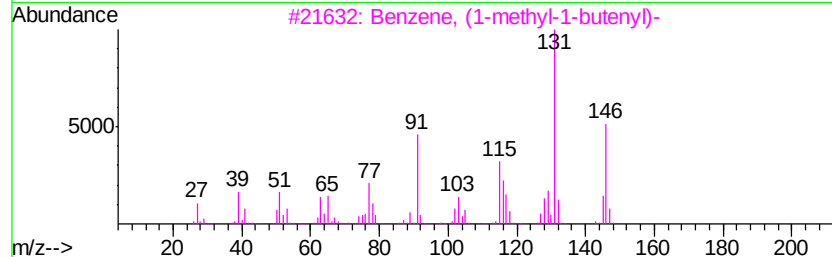
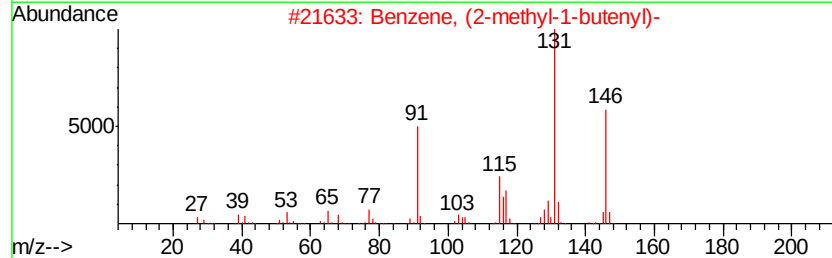
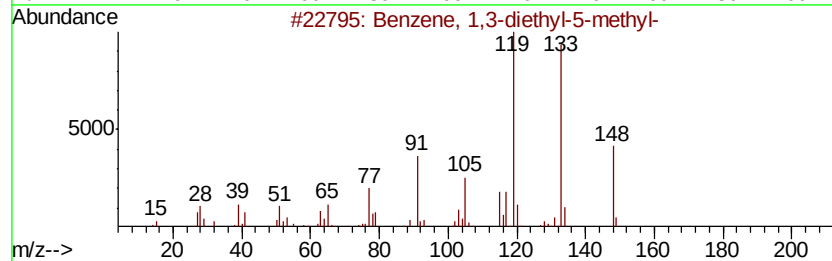
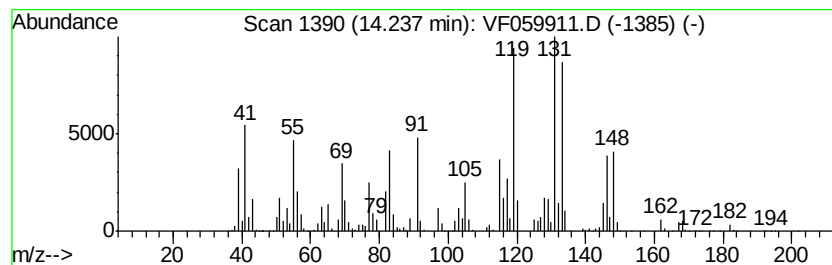
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1,3-diethyl-5-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.24	79.70 ug/l	16854700	1,4-Dichlorobenzene-d4	12.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	90
2		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	74
3		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	47
4		Benzene, 2,4-diethyl-1-methyl-	148	C11H16	001758-85-6	41
5		Benzene, (2-methyl-1-methylenepr...	146	C11H14	017498-71-4	38



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_F\DATA\VF080718\
 Data File : VF059911.D
 Acq On : 7 Aug 2018 9:48
 Operator : VA/AP
 Sample : J4297-02
 Misc : 6.00g/5mL/MSVOA_F/SOIL
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 TP-2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F080418S.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-	6.77	84.5	ug/l	1335100	2	5.78	790012	50.0
Octane	7.64	212.3	ug/l	3353980	2	5.78	790012	50.0
1-Cyclobutanone, 2...	10.66	92.8	ug/l	6922450	3	9.93	3728400	50.0
unknown11.18	11.18	74.1	ug/l	5524550	3	9.93	3728400	50.0
Decane, 4-methyl-	12.12	118.5	ug/l	25056200	4	12.65	10573400	50.0
Benzene, 1,2,3-tr...	12.71	81.8	ug/l	17302300	4	12.65	10573400	50.0
o-Cymene	13.20	82.5	ug/l	17447400	4	12.65	10573400	50.0
Dodecane	13.80	145.9	ug/l	30847700	4	12.65	10573400	50.0
Benzene, 1-methyl...	13.94	116.6	ug/l	24653800	4	12.65	10573400	50.0
Benzene, 1,3-diet...	14.24	79.7	ug/l	16854700	4	12.65	10573400	50.0