

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_F\DATA\VF121418\
 Data File : VF061023.D
 Acq On : 14 Dec 2018 16:41
 Operator : VA/AP
 Sample : J6403-21
 Misc : 5.31g/5mL/MSVOA-F/SOIL
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 TP-3-B

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\82F112618S.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	4.105	345	352	360	rVB3	8931	31965	1.44%	0.132%
2	4.854	422	428	437	rBV3	24076	82297	3.72%	0.339%
3	5.594	496	503	512	rVV	28374	77850	3.52%	0.321%
4	7.547	697	701	705	rBV2	31017	70143	3.17%	0.289%
5	7.616	705	708	714	rVB2	14600	31787	1.44%	0.131%
6	8.248	764	772	778	rBV3	18267	55096	2.49%	0.227%
7	8.613	803	809	815	rBV4	13395	45632	2.06%	0.188%
8	8.958	840	844	847	rBV3	10722	26423	1.19%	0.109%
9	9.283	873	877	881	rBV2	10033	24241	1.10%	0.100%
10	9.767	921	926	930	rBV3	31110	73907	3.34%	0.305%
11	9.895	934	939	944	rVB	86646	184382	8.33%	0.760%
12	10.122	958	962	966	rBV	31279	65051	2.94%	0.268%
13	10.220	967	972	981	rVB6	20857	67401	3.05%	0.278%
14	10.507	996	1001	1006	rVB5	22396	55555	2.51%	0.229%
15	10.664	1012	1017	1019	rBV3	56263	121781	5.50%	0.502%
16	10.704	1019	1021	1026	rVB	66704	114934	5.19%	0.474%
17	10.832	1030	1034	1041	rVB4	30891	89085	4.03%	0.367%
18	10.980	1046	1049	1051	rVV3	13741	26847	1.21%	0.111%
19	11.049	1051	1056	1059	rVV4	21440	74223	3.35%	0.306%
20	11.118	1059	1063	1066	rVV	124553	279351	12.63%	1.152%
21	11.207	1070	1072	1075	rVV3	51325	120950	5.47%	0.499%
22	11.266	1075	1078	1084	rVB2	75800	152835	6.91%	0.630%
23	11.424	1091	1094	1097	rBV4	13081	29947	1.35%	0.123%
24	11.493	1097	1101	1104	rVB	68002	109924	4.97%	0.453%
25	11.552	1104	1107	1108	rBV	43641	72356	3.27%	0.298%
26	11.611	1110	1113	1116	rVV3	56689	105920	4.79%	0.437%
27	11.651	1116	1117	1120	rVB2	36520	41030	1.85%	0.169%
28	11.720	1120	1124	1127	rVB	336583	546471	24.70%	2.253%
29	11.818	1131	1134	1138	rVB	254564	362968	16.40%	1.497%
30	11.947	1142	1147	1148	rBV5	24449	52588	2.38%	0.217%
31	12.016	1148	1154	1157	rBV	338507	652001	29.47%	2.688%
32	12.154	1165	1168	1169	rBV3	27099	47432	2.14%	0.196%
33	12.203	1169	1173	1176	rVB	626022	1059713	47.90%	4.369%
34	12.282	1178	1181	1184	rVB4	30885	53973	2.44%	0.223%

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35	12.331	1184	1186	1188	rBV2	23841	32989	1.49%	0.136%
36	12.381	1188	1191	1193	rBV2	35258	68388	3.09%	0.282%
37	12.420	1193	1195	1197	rBV	220843	342899	15.50%	1.414%
38	12.450	1197	1198	1204	rVB	211379	308459	13.94%	1.272%
39	12.529	1204	1206	1210	rVB3	34690	50057	2.26%	0.206%
40	12.608	1210	1214	1218	rBV	423012	692036	31.28%	2.853%
41	12.716	1221	1225	1229	rBV2	918356	1811178	81.86%	7.467%
42	12.795	1229	1233	1234	rVV3	174202	399286	18.05%	1.646%
43	12.825	1234	1236	1241	rVB	493486	722829	32.67%	2.980%
44	12.894	1241	1243	1244	rBV2	18411	23612	1.07%	0.097%
45	12.923	1244	1246	1251	rVV4	46824	93997	4.25%	0.388%
46	13.002	1251	1254	1258	rVB2	87704	142366	6.43%	0.587%
47	13.091	1258	1263	1265	rBV	304357	785619	35.51%	3.239%
48	13.150	1265	1269	1274	rVV	905283	2080284	94.02%	8.577%
49	13.219	1274	1276	1281	rVV	655770	1111582	50.24%	4.583%
50	13.278	1281	1282	1285	rVV	126130	228517	10.33%	0.942%
51	13.328	1285	1287	1290	rVV2	161485	325074	14.69%	1.340%
52	13.377	1290	1292	1294	rVV2	153601	269950	12.20%	1.113%
53	13.416	1294	1296	1301	rVV	200006	334153	15.10%	1.378%
54	13.495	1301	1304	1305	rVV2	138966	228149	10.31%	0.941%
55	13.535	1305	1308	1310	rVV	338581	510659	23.08%	2.105%
56	13.574	1310	1312	1315	rVV2	155720	279270	12.62%	1.151%
57	13.634	1316	1318	1322	rVV	145445	281798	12.74%	1.162%
58	13.703	1322	1325	1333	rVV	707422	1174409	53.08%	4.842%
59	13.801	1333	1335	1337	rVV	67021	119216	5.39%	0.492%
60	13.851	1337	1340	1345	rVV3	593362	1413673	63.89%	5.829%
61	13.910	1345	1346	1348	rVV	53308	72385	3.27%	0.298%
62	13.969	1348	1352	1356	rVV	291067	526617	23.80%	2.171%
63	14.038	1356	1359	1361	rVV	164028	318432	14.39%	1.313%
64	14.077	1361	1363	1366	rVV	536085	856780	38.72%	3.532%
65	14.146	1366	1370	1374	rVV3	116341	260525	11.77%	1.074%
66	14.215	1374	1377	1382	rVB	122750	197602	8.93%	0.815%
67	14.304	1383	1386	1389	rBV4	26417	56777	2.57%	0.234%
68	14.354	1389	1391	1394	rVB3	32025	49166	2.22%	0.203%
69	14.452	1394	1401	1407	rBV	1307862	2212549	100.00%	9.122%
70	14.580	1412	1414	1417	rVB3	22182	38404	1.74%	0.158%
71	14.679	1422	1424	1427	rBV2	44862	65152	2.94%	0.269%

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72	14.728	1427	1429	1434	rVB	71429	123399	5.58%	0.509%
73	14.906	1445	1447	1454	rVB3	49367	90750	4.10%	0.374%
74	15.113	1464	1468	1471	rVV5	16332	52261	2.36%	0.215%
75	15.162	1471	1473	1477	rVV4	24313	59008	2.67%	0.243%
76	15.281	1480	1485	1492	rVV	112462	236978	10.71%	0.977%
77	15.399	1494	1497	1506	rVB	85478	169676	7.67%	0.700%
78	15.991	1553	1557	1561	rVB2	14273	31440	1.42%	0.130%

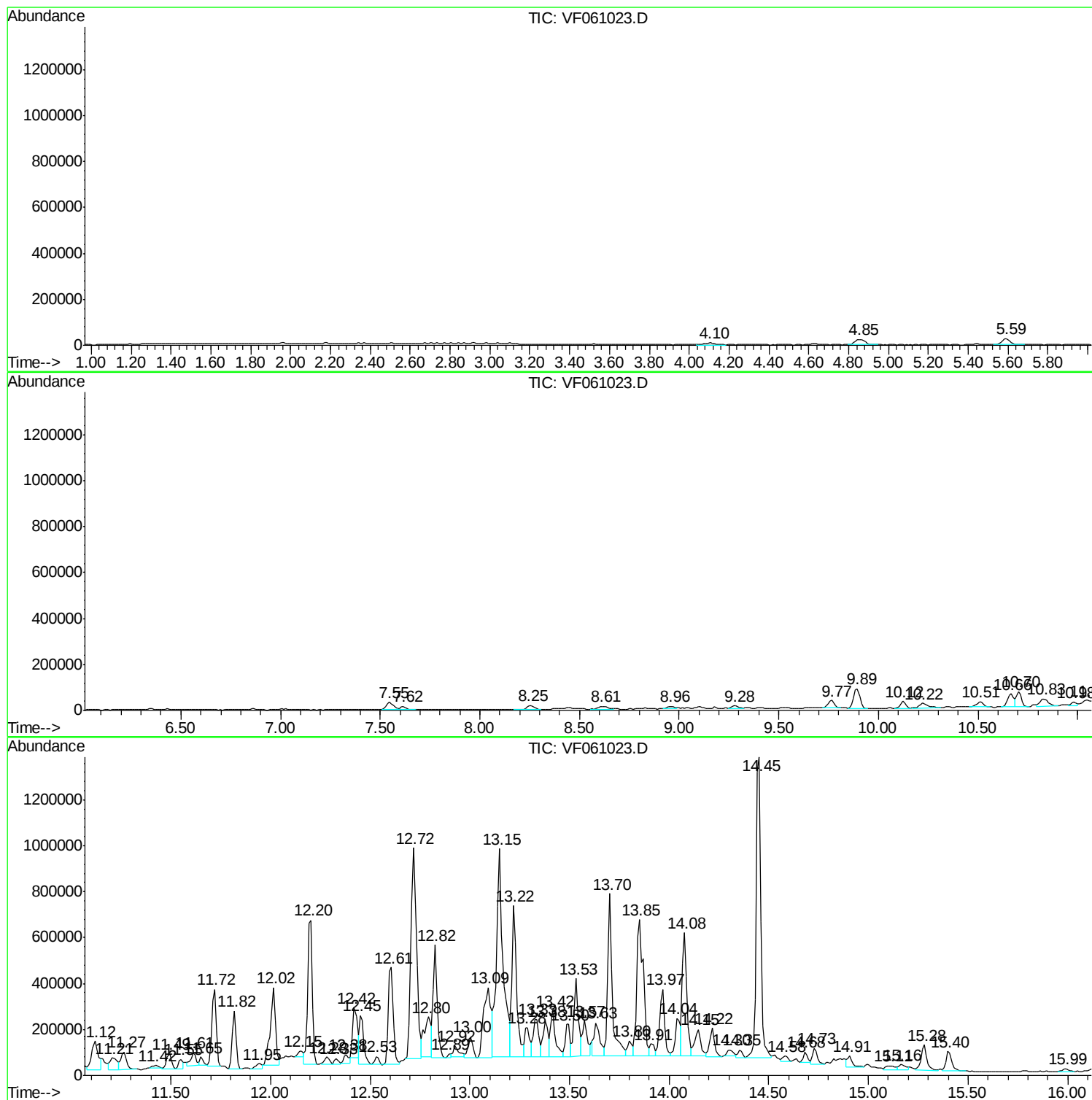
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Data Path : Z:\V0ASRV\HPCHEM1\MSV0A_F\DATA\VF121418\
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Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_F\METHODS\82F112618S.M
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TIC Library : C:\DATABASE\NIST11.L
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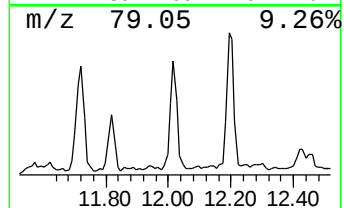
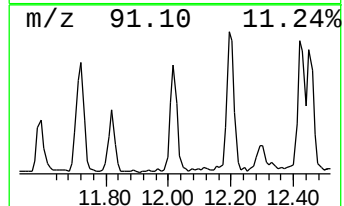
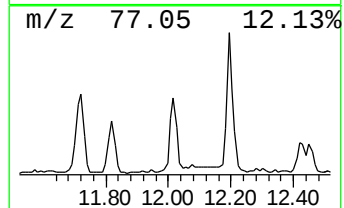
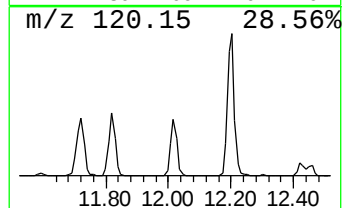
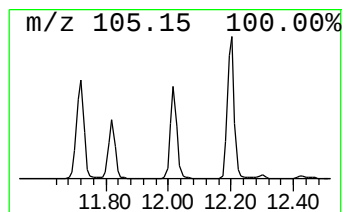
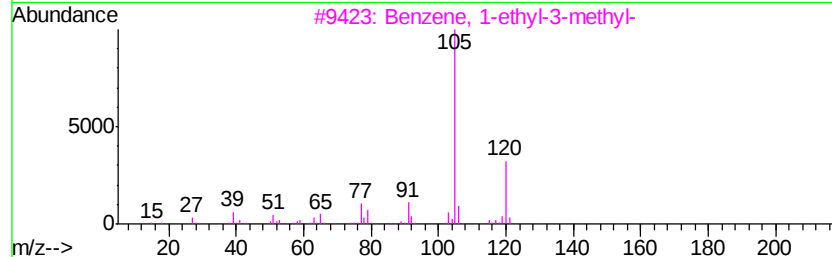
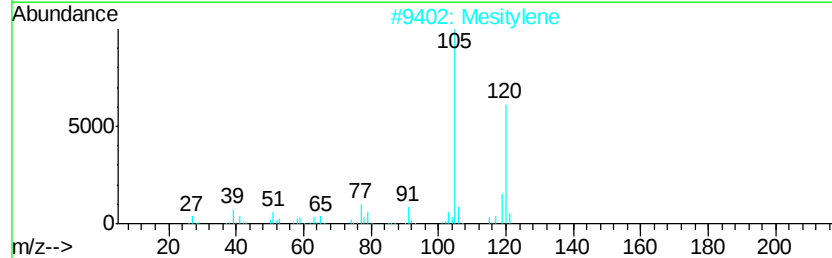
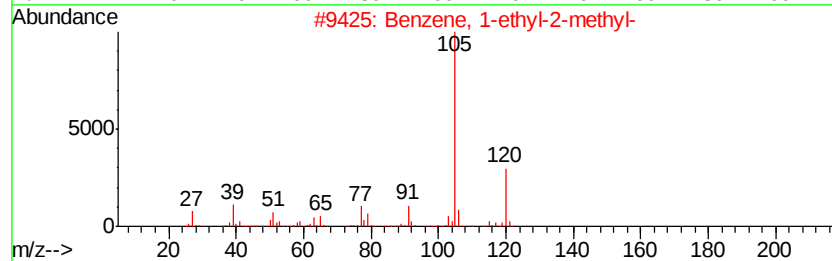
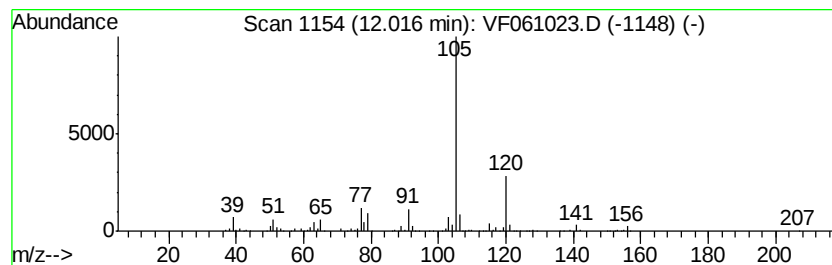
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F112618S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.02	651.26 ug/l	652001	1,4-Dichlorobenzene-d4	12.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Mesitylene	120	C9H12	000108-67-8	91
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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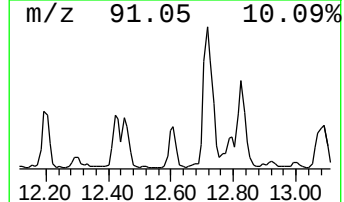
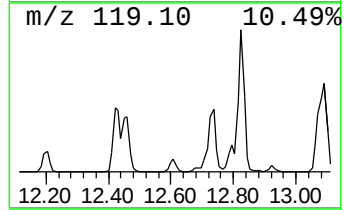
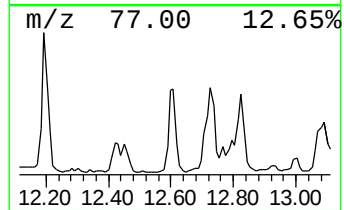
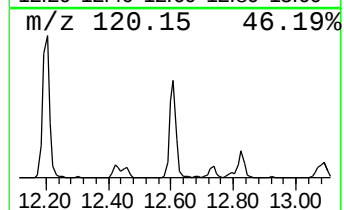
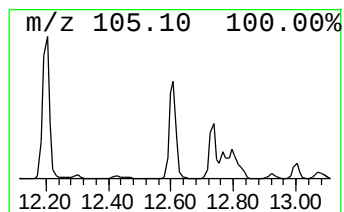
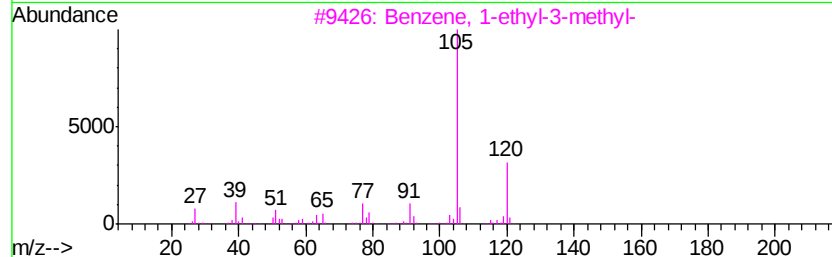
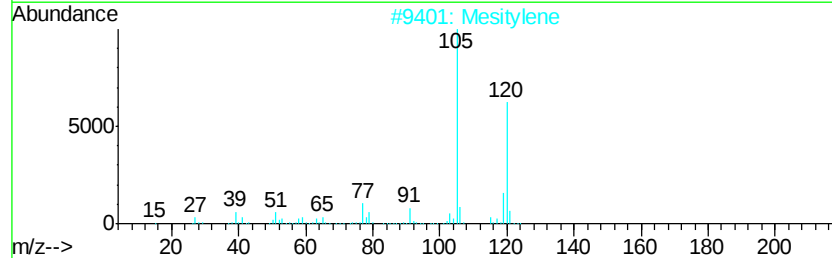
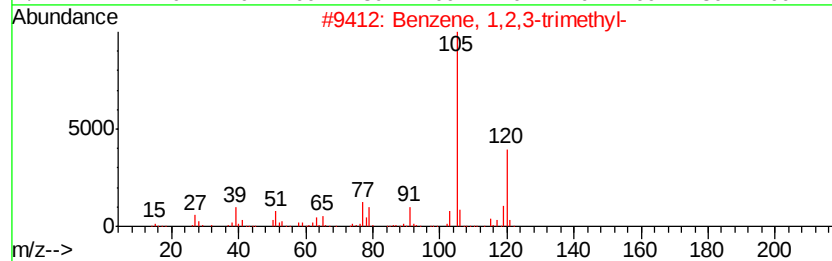
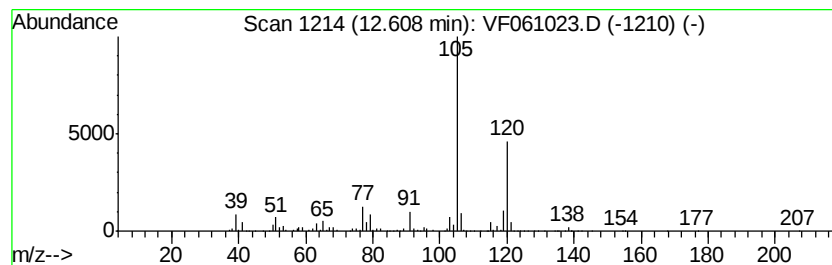
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Benzene, 1,2,3-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.61	691.25 ug/l	692036	1,4-Dichlorobenzene-d4	12.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Mesitylene	120	C9H12	000108-67-8	94
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90



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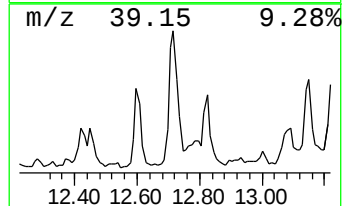
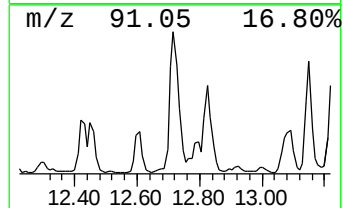
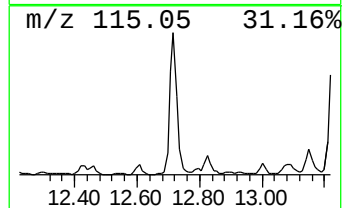
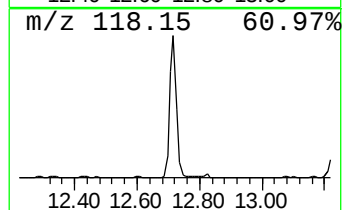
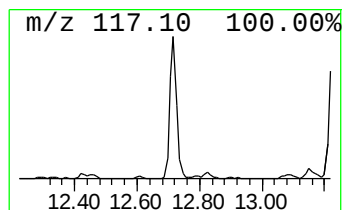
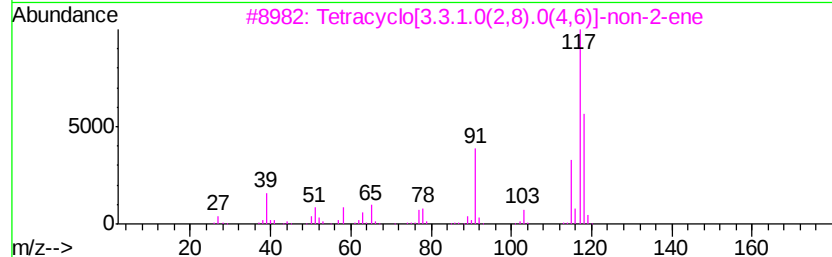
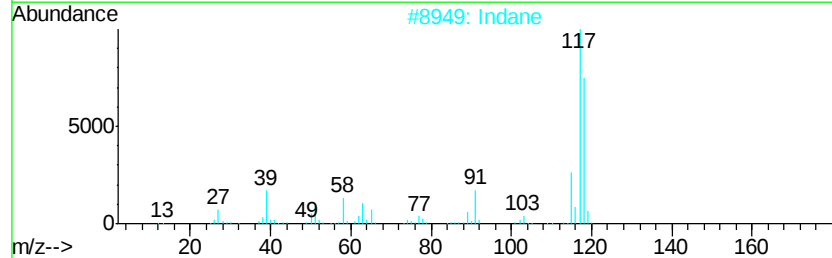
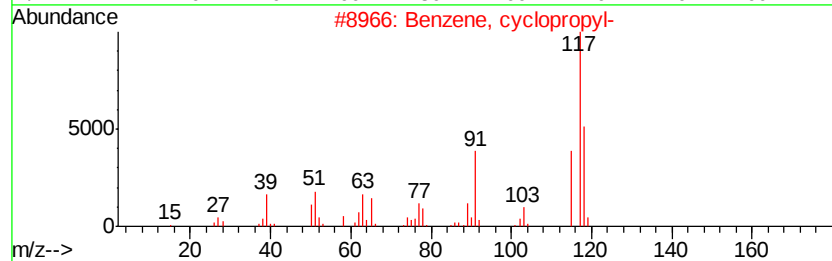
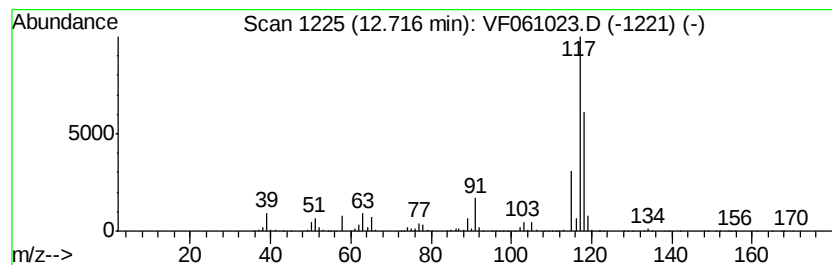
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, cyclopropyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.72	1809.12 ug/l	1811180	1,4-Dichlorobenzene-d4	12.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	94
2		Indane	118	C9H10	000496-11-7	93
3		Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	83
4		Benzene, 2-propenyl-	118	C9H10	000300-57-2	81
5		Benzene, 1-propenyl-	118	C9H10	000637-50-3	81



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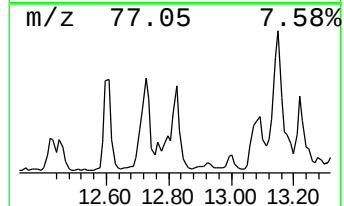
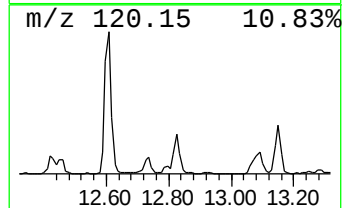
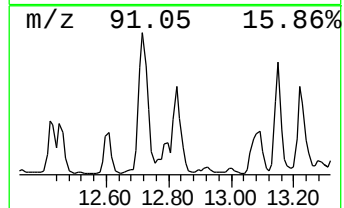
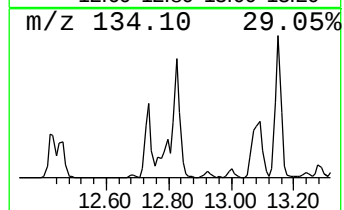
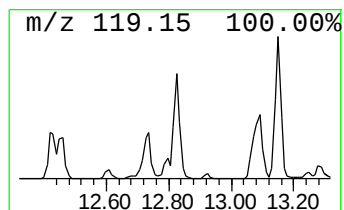
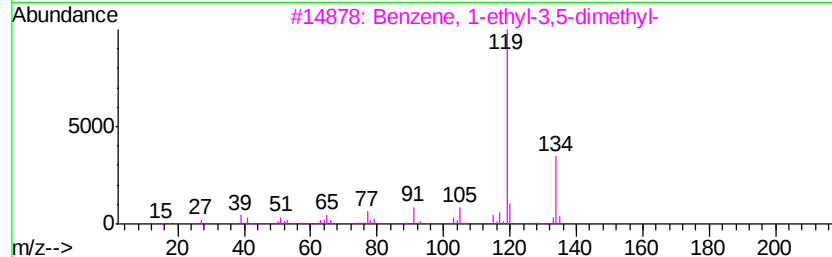
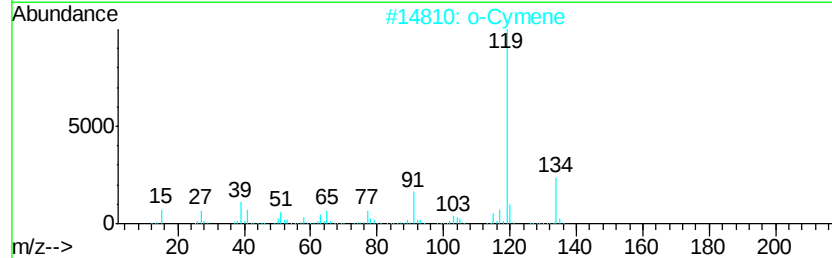
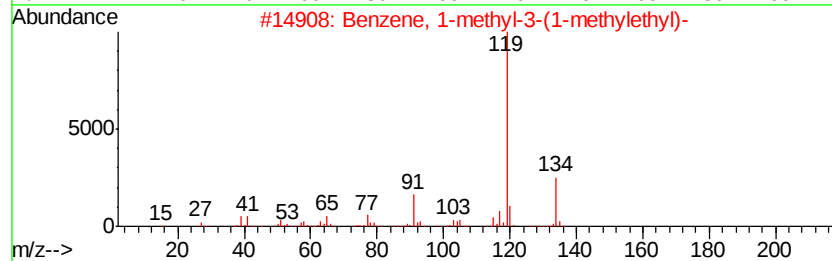
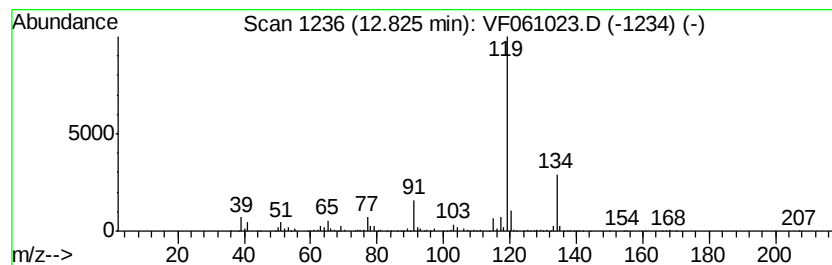
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-methyl-3-(1-meth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.82	722.01 ug/l	722829	1,4-Dichlorobenzene-d4	12.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
2		o-Cymene	134	C10H14	000527-84-4	95
3		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91
5		p-Cymene	134	C10H14	000099-87-6	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF121418\
Data File : VF061023.D
Acq On : 14 Dec 2018 16:41
Operator : VA/AP
Sample : J6403-21
Misc : 5.31g/5mL/MSVOA-F/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
TP-3-B

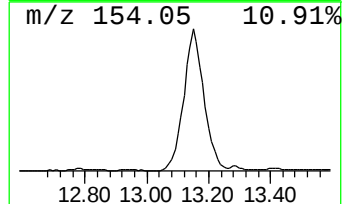
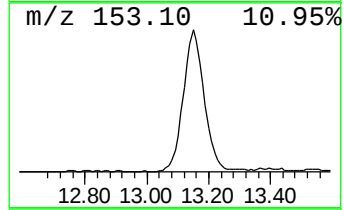
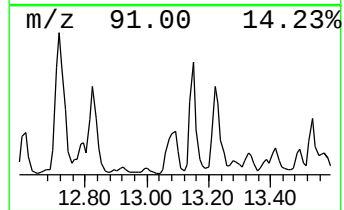
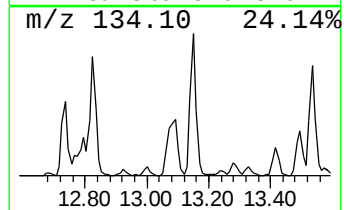
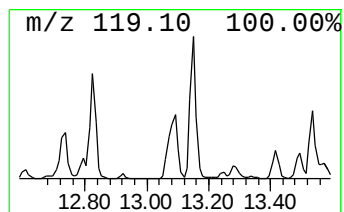
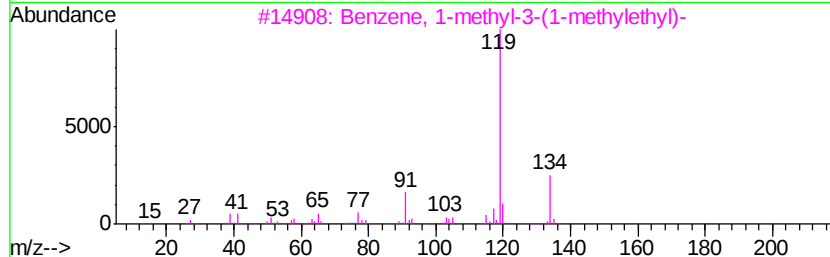
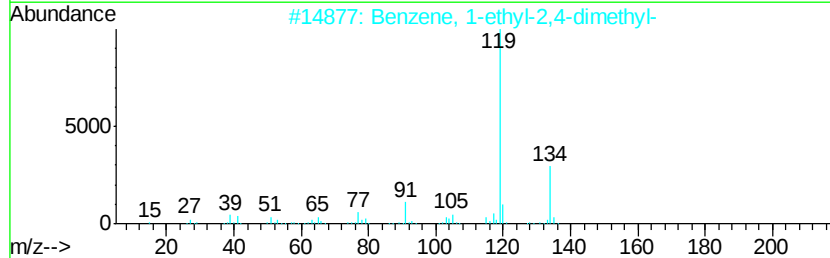
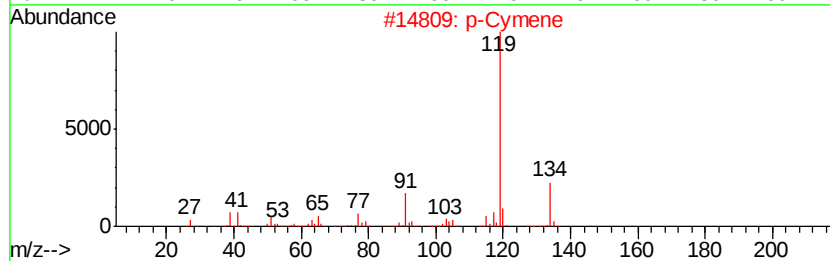
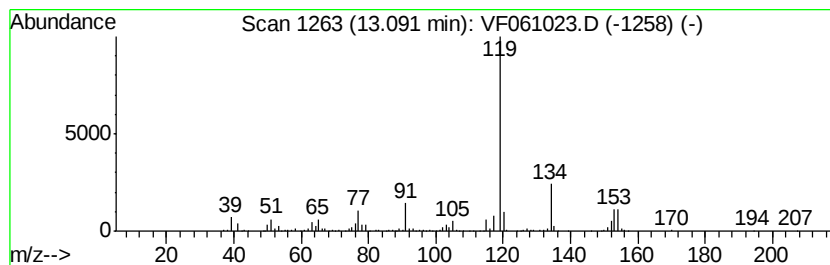
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Quant Title : SW846 8260

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

Peak Number 5 p-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.09	784.72 ug/l	785619	1,4-Dichlorobenzene-d4	12.54

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene		134	C10H14	000099-87-6	93
2	Benzene, 1-ethyl-2,4-dimethyl-		134	C10H14	000874-41-9	93
3	Benzene, 1-methyl-3-(1-methyleth...		134	C10H14	000535-77-3	93
4	o-Cymene		134	C10H14	000527-84-4	87
5	Benzene, 1,2,3,4-tetramethyl-		134	C10H14	000488-23-3	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF121418\
Data File : VF061023.D
Acq On : 14 Dec 2018 16:41
Operator : VA/AP
Sample : J6403-21
Misc : 5.31g/5mL/MSVOA-F/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
TP-3-B

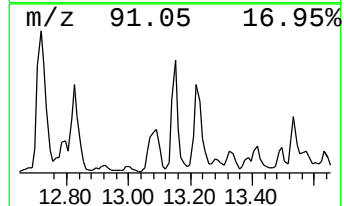
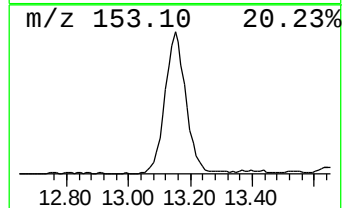
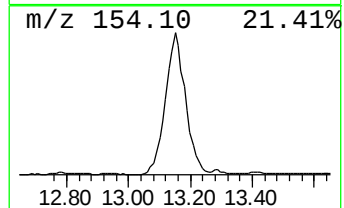
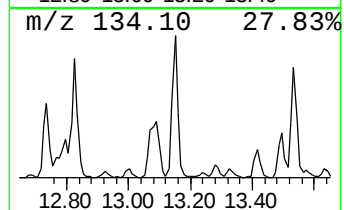
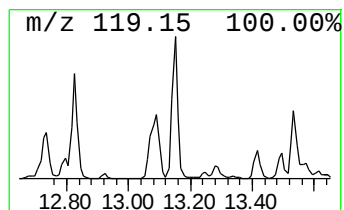
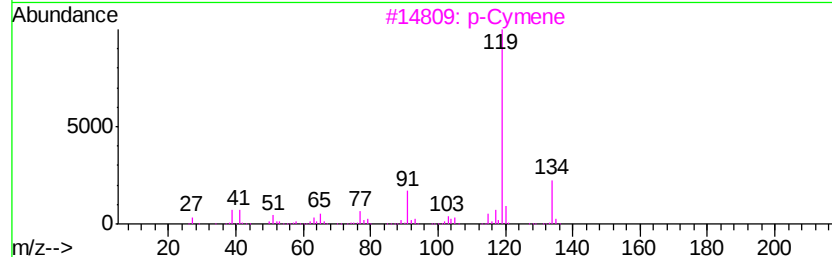
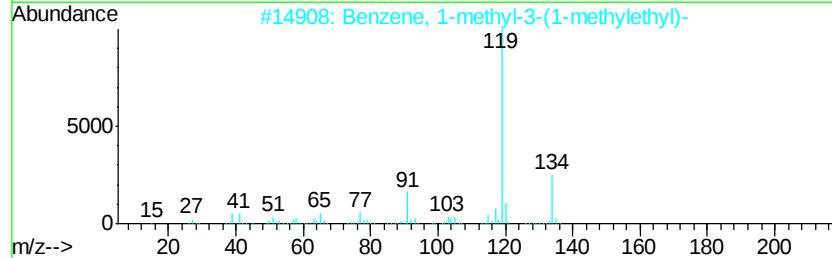
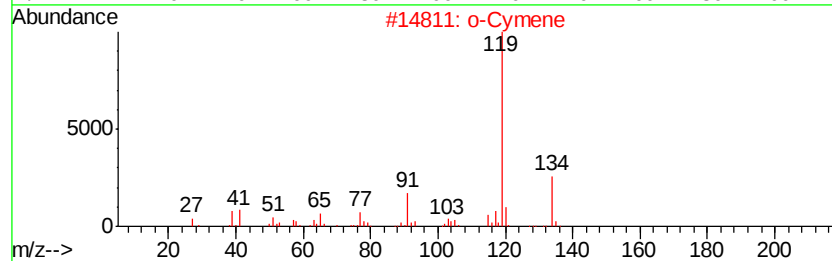
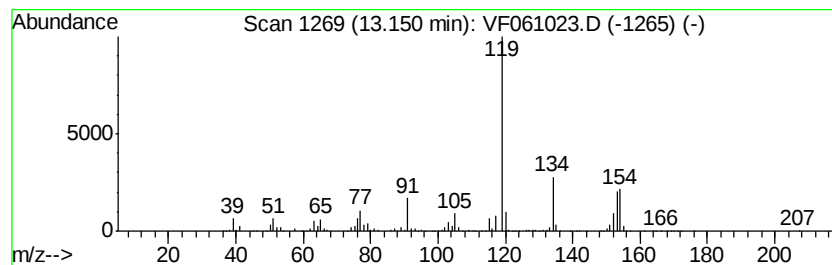
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Quant Title : SW846 8260

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

Peak Number 6 o-Cymene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.15	2077.92 ug/l	2080280	1,4-Dichlorobenzene-d4	12.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	92
2		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91
3		p-Cymene	134	C10H14	000099-87-6	91
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	86
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF121418\
Data File : VF061023.D
Acq On : 14 Dec 2018 16:41
Operator : VA/AP
Sample : J6403-21
Misc : 5.31g/5mL/MSVOA-F/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
TP-3-B

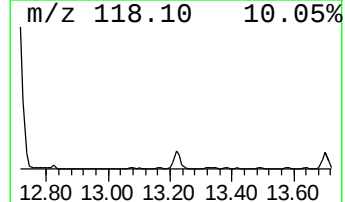
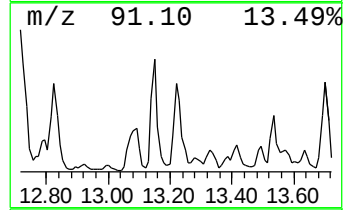
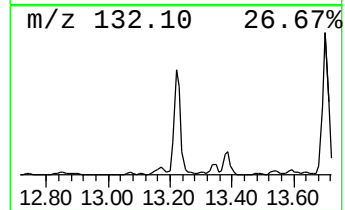
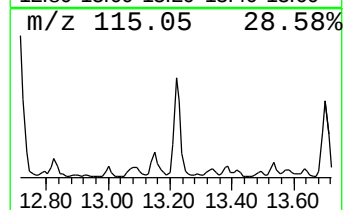
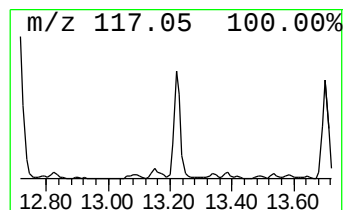
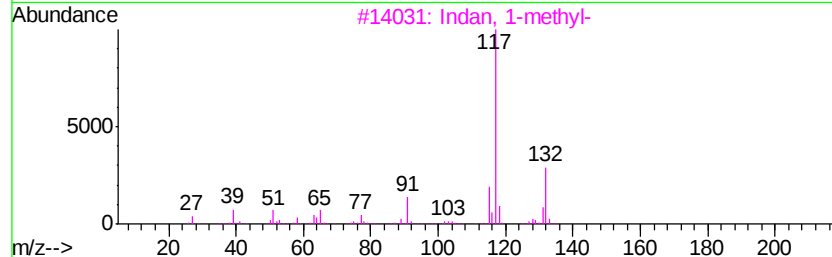
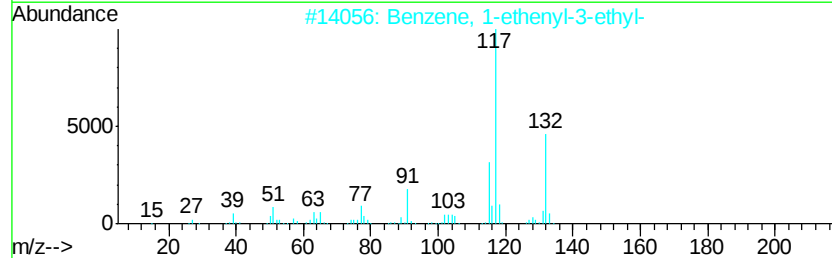
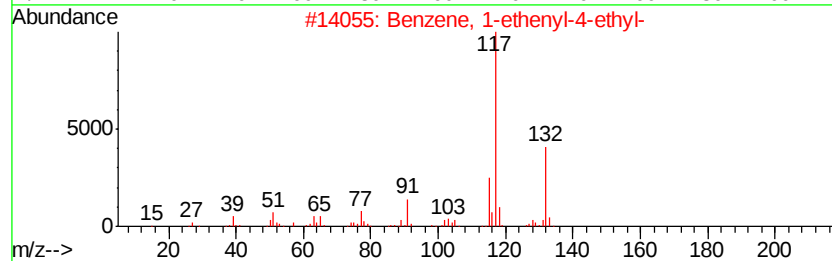
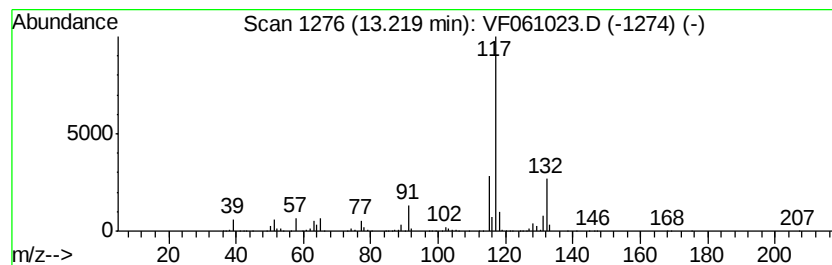
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	1110.32 ug/l	1111580	1,4-Dichlorobenzene-d4	12.54

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	91
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
3	Indan, 1-methyl-	132	C10H12	000767-58-8	90
4	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF121418\
Data File : VF061023.D
Acq On : 14 Dec 2018 16:41
Operator : VA/AP
Sample : J6403-21
Misc : 5.31g/5mL/MSVOA-F/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TP-3-B

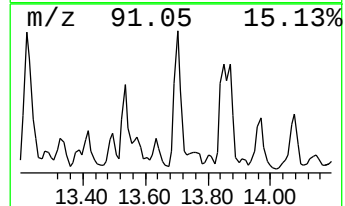
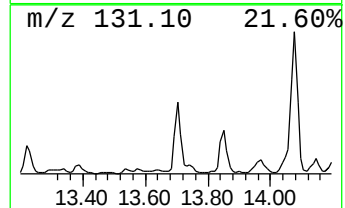
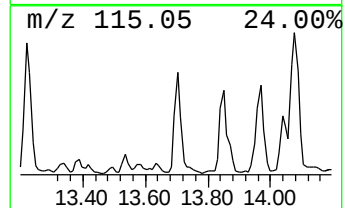
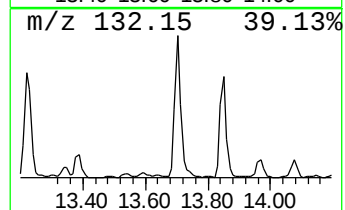
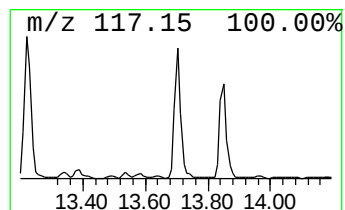
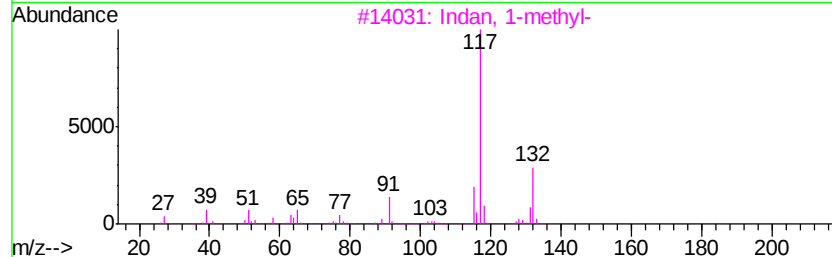
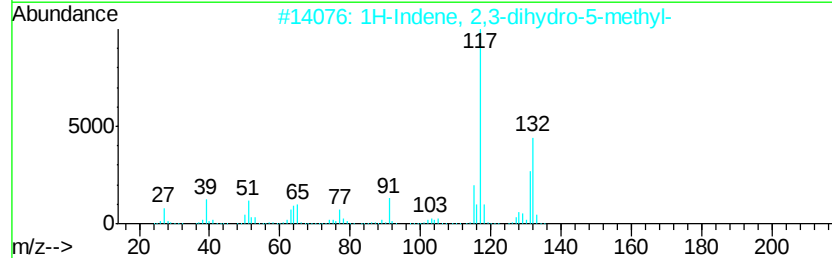
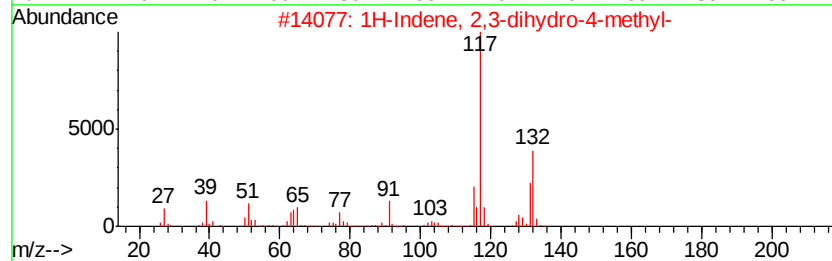
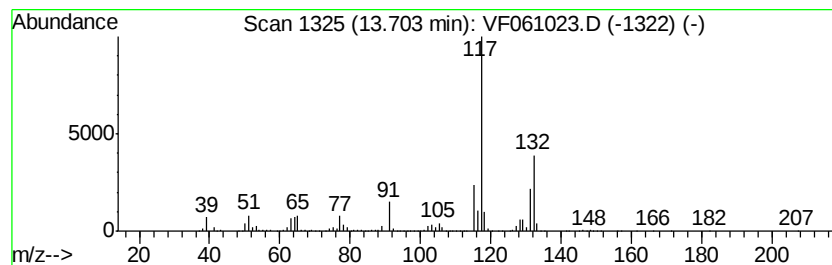
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Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.70	1173.07 ug/l	1174410	1,4-Dichlorobenzene-d4	12.54

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	95
3		Indan, 1-methyl-	132	C10H12	000767-58-8	90
4		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF121418\
Data File : VF061023.D
Acq On : 14 Dec 2018 16:41
Operator : VA/AP
Sample : J6403-21
Misc : 5.31g/5mL/MSVOA-F/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
TP-3-B

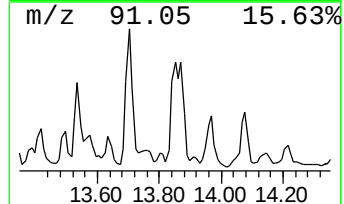
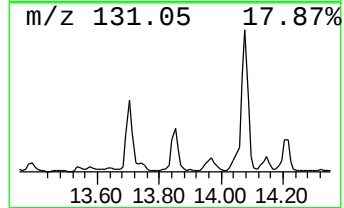
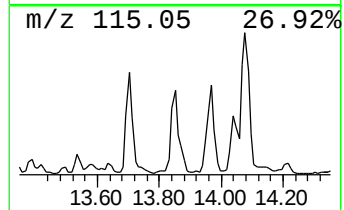
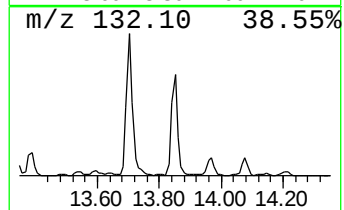
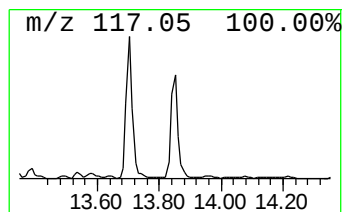
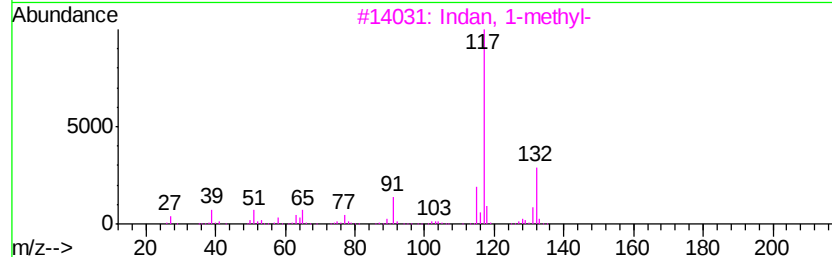
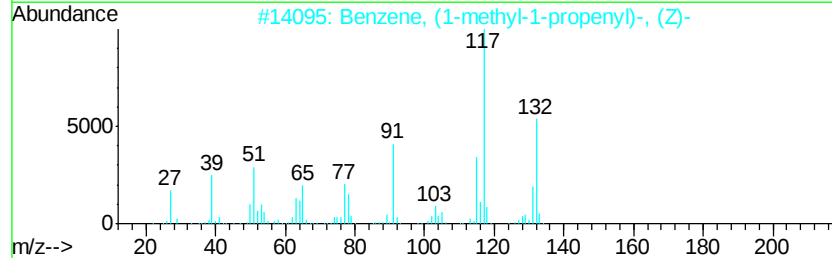
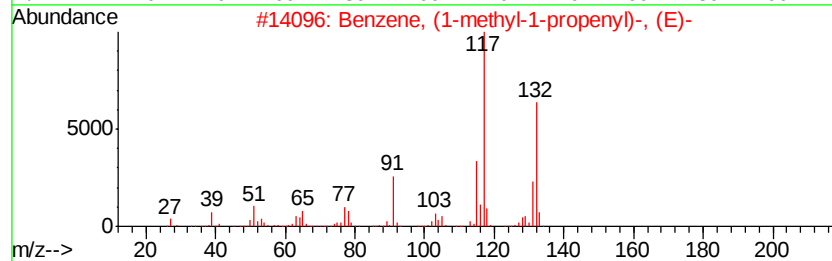
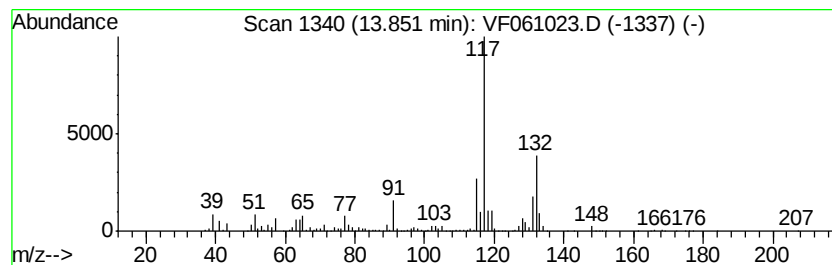
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, (1-methyl-1-propen... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.85	1412.06 ug/l	1413670	1,4-Dichlorobenzene-d4	12.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
2		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	87
3		Indan, 1-methyl-	132	C10H12	000767-58-8	87
4		1-Phenyl-1-butene	132	C10H12	000824-90-8	87
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF121418\
Data File : VF061023.D
Acq On : 14 Dec 2018 16:41
Operator : VA/AP
Sample : J6403-21
Misc : 5.31g/5mL/MSVOA-F/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TP-3-B

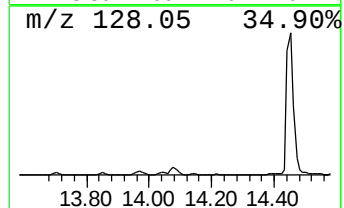
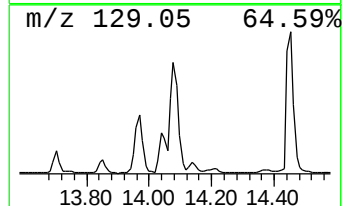
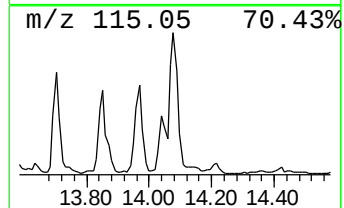
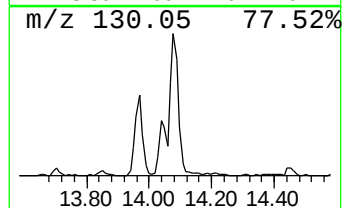
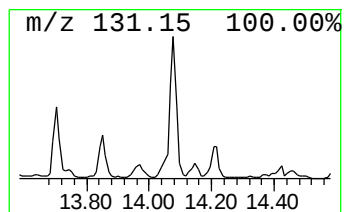
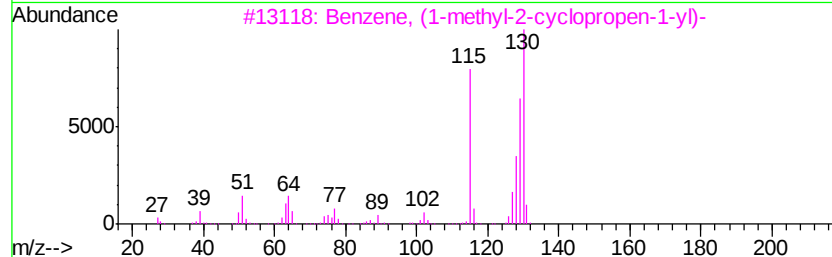
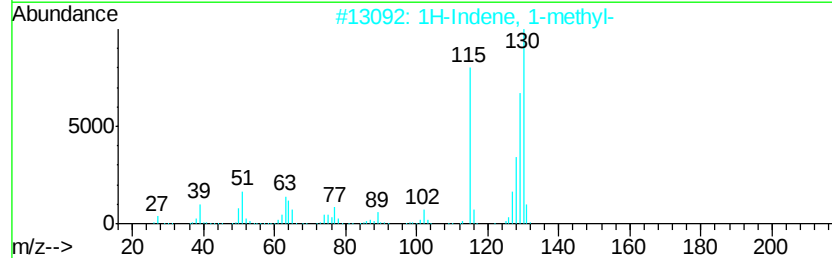
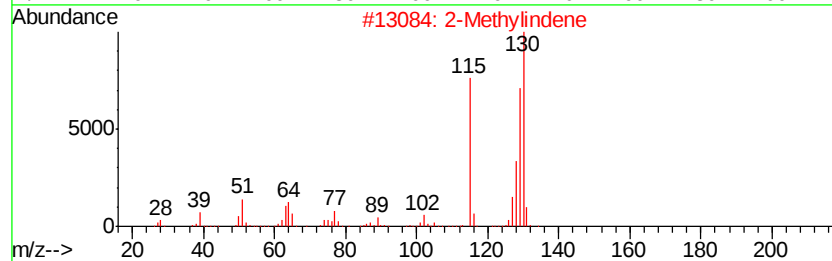
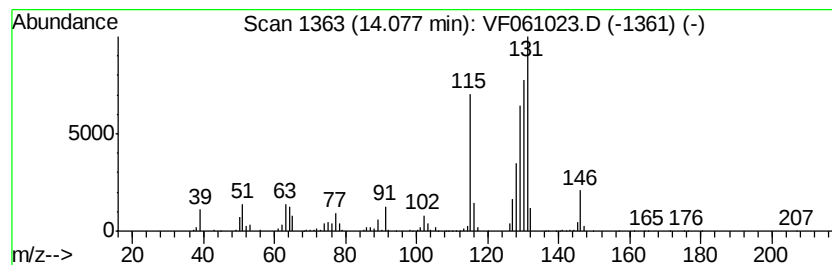
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Quant Title : SW846 8260

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

Peak Number 10 2-Methylindene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.08	855.80 ug/l	856780	1,4-Dichlorobenzene-d4	12.54

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methylindene		130	C10H10	002177-47-1	91
2	1H-Indene, 1-methyl-		130	C10H10	000767-59-9	91
3	Benzene, (1-methyl-2-cyclopropen...		130	C10H10	065051-83-4	90
4	Cycloprop[a]indene, 1,1a,6,6a-te...		130	C10H10	015677-15-3	90
5	Naphthalene, 1,2-dihydro-		130	C10H10	000447-53-0	78



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_F\DATA\VF121418\
 Data File : VF061023.D
 Acq On : 14 Dec 2018 16:41
 Operator : VA/AP
 Sample : J6403-21
 Misc : 5.31g/5mL/MSVOA-F/SOIL
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 TP-3-B

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F112618S.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
Benzene, 1-ethyl-...	12.02	651.3	ug/l	652001	4	12.54	50057	50.0
Benzene, 1,2,3-tr...	12.61	691.3	ug/l	692036	4	12.54	50057	50.0
Benzene, cyclopro...	12.72	1809.1	ug/l	1811180	4	12.54	50057	50.0
Benzene, 1-methyl...	12.82	722.0	ug/l	722829	4	12.54	50057	50.0
p-Cymene	13.09	784.7	ug/l	785619	4	12.54	50057	50.0
o-Cymene	13.15	2077.9	ug/l	2080280	4	12.54	50057	50.0
Benzene, 1-etheny...	13.22	1110.3	ug/l	1111580	4	12.54	50057	50.0
1H-Indene, 2,3-di...	13.70	1173.1	ug/l	1174410	4	12.54	50057	50.0
Benzene, (1-methy...	13.85	1412.1	ug/l	1413670	4	12.54	50057	50.0
2-Methylindene	14.08	855.8	ug/l	856780	4	12.54	50057	50.0