

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF031919\
 Data File : VF062015.D
 Acq On : 19 Mar 2019 13:47
 Operator : VA/AP
 Sample : K1693-09
 Misc : 6.50g/5mL/MSVOA-F/SOIL
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 OR-03-031819-A

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\82F031619S.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.637	96	102	103	rBV3	4159	11157	1.22%	0.111%
2	2.160	151	155	158	rBV4	14432	36601	4.00%	0.366%
3	2.229	158	162	167	rVV3	21909	73092	7.98%	0.730%
4	2.367	170	176	182	rVB4	6739	26103	2.85%	0.261%
5	2.574	192	197	199	rBV3	4095	10183	1.11%	0.102%
6	3.994	336	341	345	rBV3	4002	14486	1.58%	0.145%
7	4.083	345	350	363	rVB	137615	444550	48.53%	4.443%
8	4.832	418	426	438	rBV2	247924	916039	100.00%	9.154%
9	5.434	484	487	493	rBV5	3101	10331	1.13%	0.103%
10	5.572	493	501	510	rVV	238214	643692	70.27%	6.433%
11	7.534	695	700	704	rVV	48355	123118	13.44%	1.230%
12	7.593	704	706	713	rVB	10886	24594	2.68%	0.246%
13	8.195	762	767	772	rBV3	4173	14643	1.60%	0.146%
14	9.753	919	925	931	rBV	131963	295743	32.28%	2.955%
15	9.871	934	937	942	rVB2	11242	25384	2.77%	0.254%
16	10.108	958	961	966	rVB4	8981	21451	2.34%	0.214%
17	10.197	966	970	974	rBV3	4034	11659	1.27%	0.117%
18	10.483	997	999	1005	rVB3	4237	9750	1.06%	0.097%
19	10.690	1012	1020	1024	rBV	33196	72001	7.86%	0.720%
20	10.828	1029	1034	1037	rVB5	5889	11618	1.27%	0.116%
21	11.035	1051	1055	1057	rBV3	4863	9894	1.08%	0.099%
22	11.104	1059	1062	1064	rVV3	6379	11077	1.21%	0.111%
23	11.252	1074	1077	1080	rVB3	8020	14408	1.57%	0.144%
24	11.390	1085	1091	1097	rBV	311993	550774	60.13%	5.504%
25	11.469	1097	1099	1103	rVB4	10925	20800	2.27%	0.208%
26	11.538	1103	1106	1108	rBV2	14347	25769	2.81%	0.258%
27	11.577	1108	1110	1114	rVB2	17323	32355	3.53%	0.323%
28	11.705	1117	1123	1126	rBV	49595	97226	10.61%	0.972%
29	11.804	1130	1133	1137	rVB	51833	76587	8.36%	0.765%
30	12.001	1147	1153	1157	rVB	64028	119706	13.07%	1.196%
31	12.139	1163	1167	1168	rBV3	6681	14238	1.55%	0.142%
32	12.188	1168	1172	1176	rVB2	37550	59326	6.48%	0.593%
33	12.287	1179	1182	1186	rVB2	11177	25014	2.73%	0.250%
34	12.415	1192	1195	1197	rVV	27642	40069	4.37%	0.400%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	12.445	1197	1198	1203	rVB4	21590	29041	3.17%	0.290%
36	12.534	1203	1207	1210	rBV	392787	581514	63.48%	5.811%
37	12.593	1210	1213	1217	rVB	129911	192290	20.99%	1.922%
38	12.701	1221	1224	1227	rBV2	91496	168030	18.34%	1.679%
39	12.760	1227	1230	1231	rVV	72397	114353	12.48%	1.143%
40	12.810	1231	1235	1241	rVB3	98994	243116	26.54%	2.430%
41	12.918	1242	1246	1248	rBV4	22609	40692	4.44%	0.407%
42	12.987	1250	1253	1257	rVB	64070	99016	10.81%	0.989%
43	13.076	1257	1262	1266	rBV	64699	159037	17.36%	1.589%
44	13.135	1266	1268	1270	rBV	104326	156401	17.07%	1.563%
45	13.214	1274	1276	1280	rVV2	83496	171467	18.72%	1.714%
46	13.273	1280	1282	1289	rVB5	20185	60423	6.60%	0.604%
47	13.362	1289	1291	1293	rVB2	20544	25788	2.82%	0.258%
48	13.401	1293	1295	1301	rBV	35003	70510	7.70%	0.705%
49	13.480	1301	1303	1305	rBV	60718	71883	7.85%	0.718%
50	13.520	1305	1307	1310	rBV	90042	155038	16.92%	1.549%
51	13.559	1310	1311	1314	rVV	40790	42473	4.64%	0.424%
52	13.599	1314	1315	1317	rVV2	18887	29133	3.18%	0.291%
53	13.687	1321	1324	1328	rBV2	175228	310685	33.92%	3.105%
54	13.746	1328	1330	1333	rVB3	37672	58996	6.44%	0.590%
55	13.796	1333	1335	1337	rBV	44440	61739	6.74%	0.617%
56	13.835	1337	1339	1343	rVV2	266292	437908	47.80%	4.376%
57	13.904	1343	1346	1348	rVB3	53436	91347	9.97%	0.913%
58	13.954	1348	1351	1357	rVB	204612	283263	30.92%	2.831%
59	14.032	1357	1359	1360	rBV	34613	49040	5.35%	0.490%
60	14.062	1360	1362	1365	rVB	76724	106269	11.60%	1.062%
61	14.141	1365	1370	1372	rBV4	95361	210735	23.01%	2.106%
62	14.200	1372	1376	1382	rVB2	95458	218628	23.87%	2.185%
63	14.279	1382	1384	1386	rBV4	8869	14043	1.53%	0.140%
64	14.348	1387	1391	1392	rBV4	12018	22975	2.51%	0.230%
65	14.397	1392	1396	1399	rVV3	78495	148511	16.21%	1.484%
66	14.447	1399	1401	1409	rVB3	95778	220994	24.12%	2.208%
67	14.545	1409	1411	1412	rBV	22707	25067	2.74%	0.251%
68	14.575	1412	1414	1417	rVB	74484	95062	10.38%	0.950%
69	14.723	1426	1429	1434	rVB	272916	367696	40.14%	3.674%
70	14.871	1436	1444	1448	rBV3	40912	152787	16.68%	1.527%
71	14.920	1448	1449	1452	rVB2	20076	21633	2.36%	0.216%

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72	14.979	1452	1455	1458	rBV	129029	182732	19.95%	1.826%
73	15.097	1464	1467	1469	rBV	103180	157772	17.22%	1.577%
74	15.226	1477	1480	1481	rBV3	8688	14110	1.54%	0.141%
75	15.265	1481	1484	1486	rVV	44417	72826	7.95%	0.728%
76	15.334	1486	1491	1494	rVV2	44687	111932	12.22%	1.119%
77	15.393	1494	1497	1502	rVV2	47115	83357	9.10%	0.833%
78	15.512	1504	1509	1511	rVB4	7632	22163	2.42%	0.221%
79	15.620	1517	1520	1525	rVB	45408	86009	9.39%	0.860%
80	15.699	1525	1528	1531	rBV3	8860	19252	2.10%	0.192%
81	15.807	1537	1539	1542	rVB4	7956	11157	1.22%	0.111%
82	15.985	1553	1557	1564	rVB6	13231	44372	4.84%	0.443%

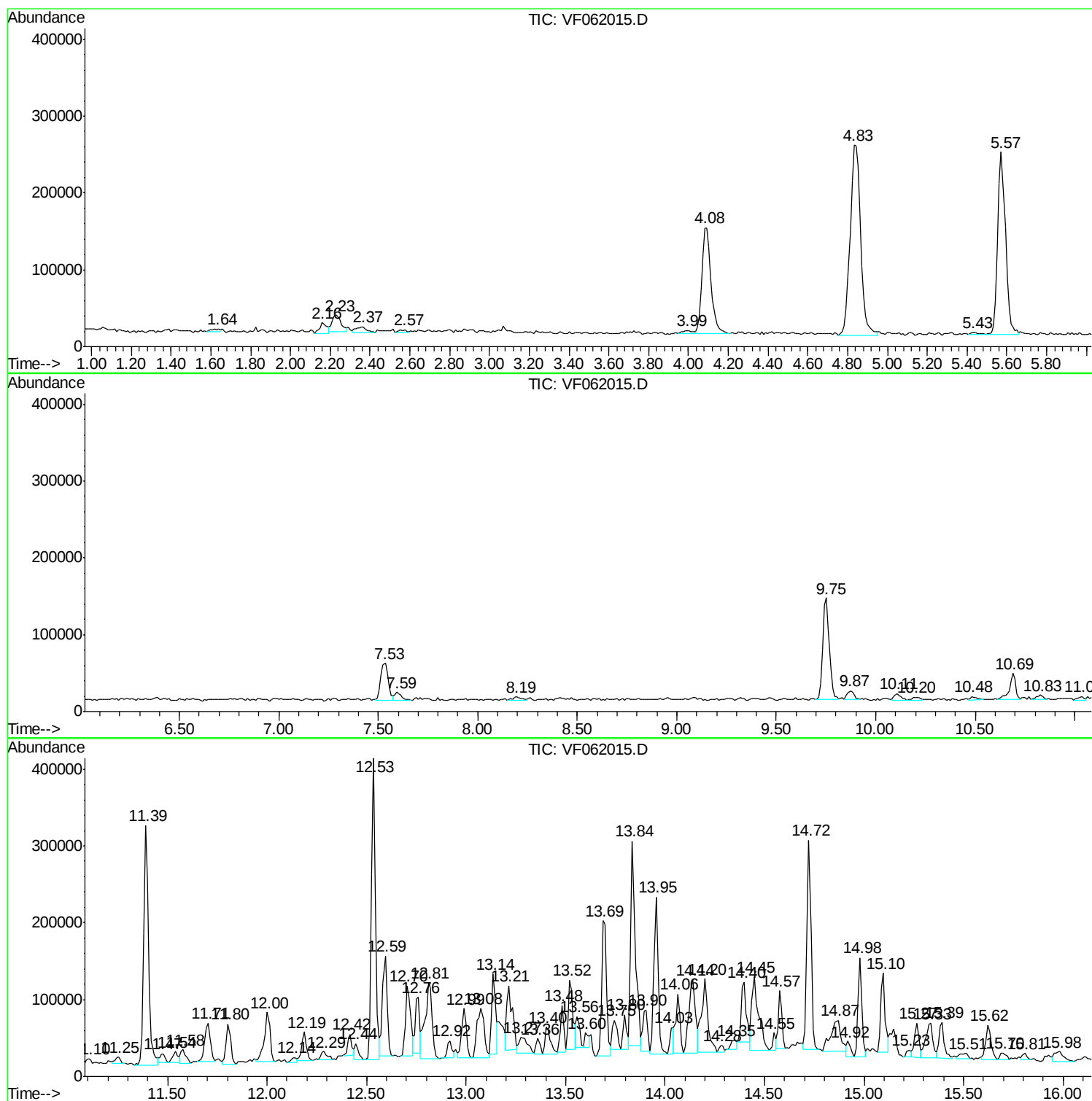
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F031619S.M
Quant Title : SW846 8260

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



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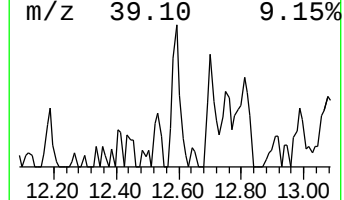
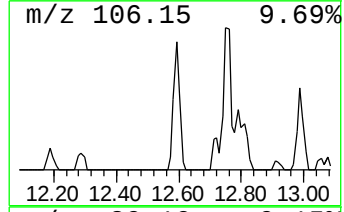
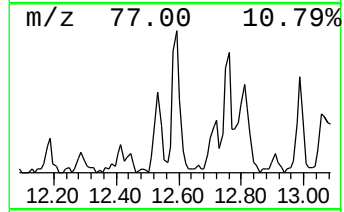
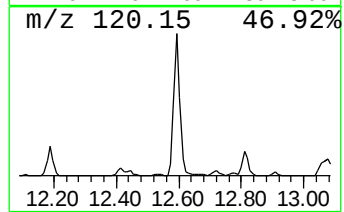
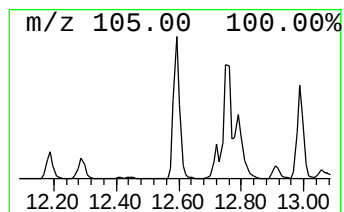
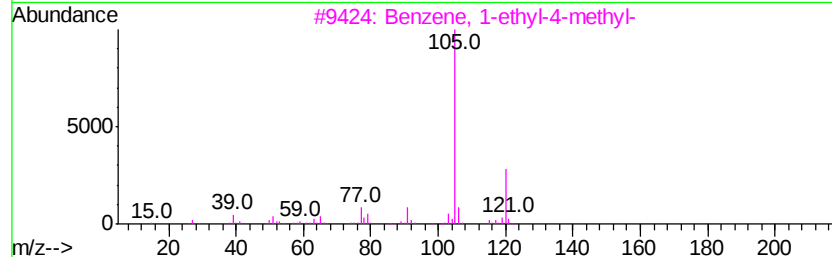
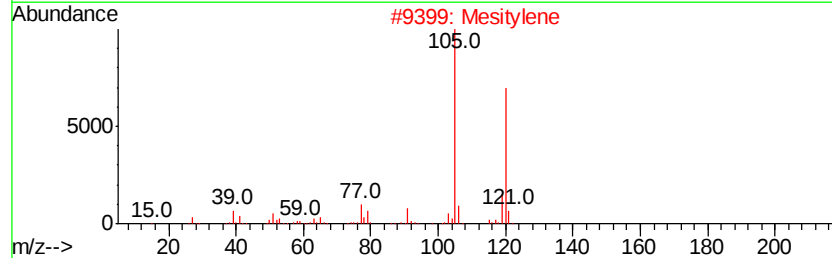
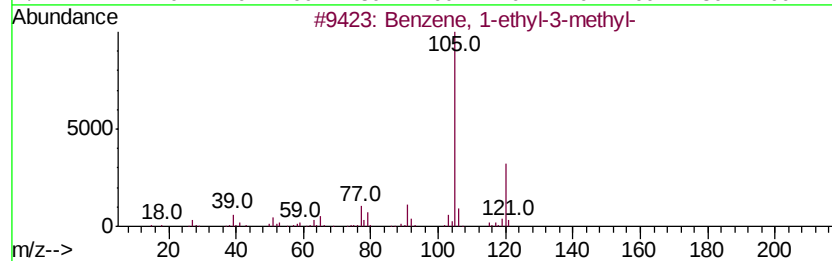
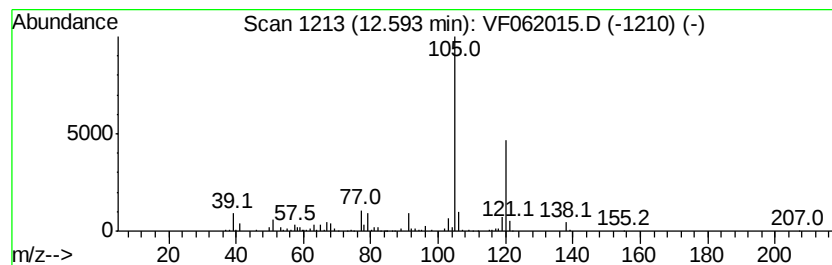
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F031619S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.59	16.53 ug/l	192290	1,4-Dichlorobenzene-d4	12.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
2		Mesitylene	120	C9H12	000108-67-8	91
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91



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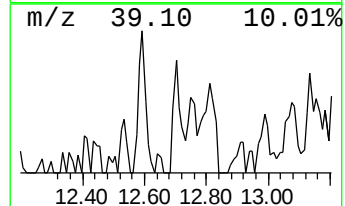
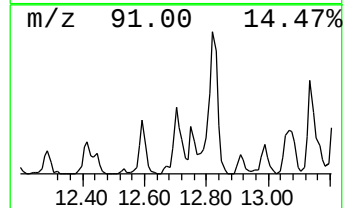
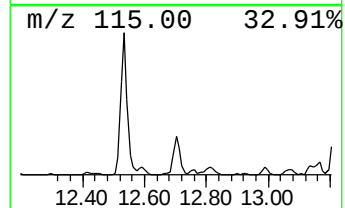
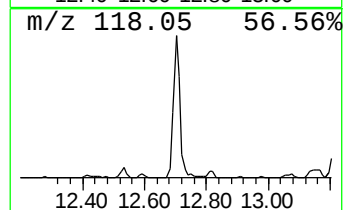
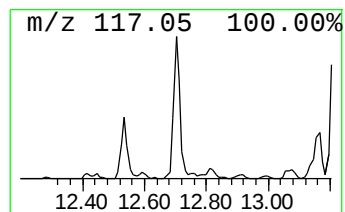
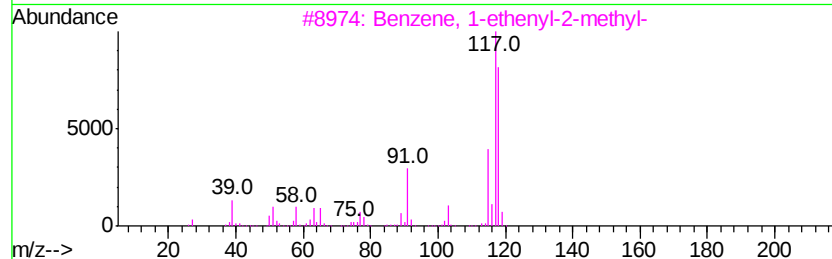
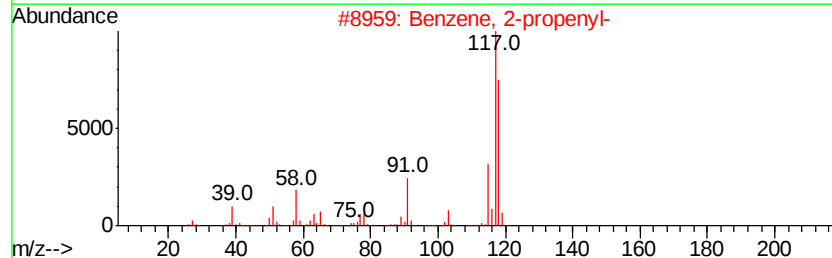
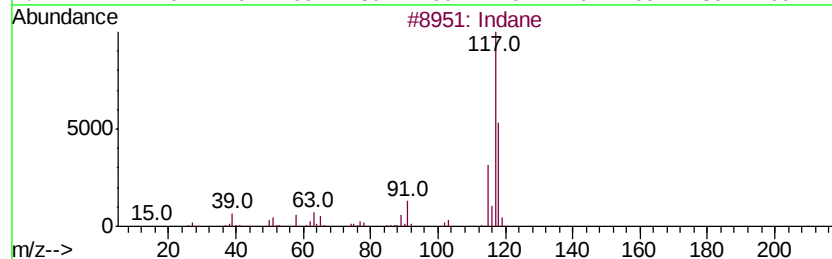
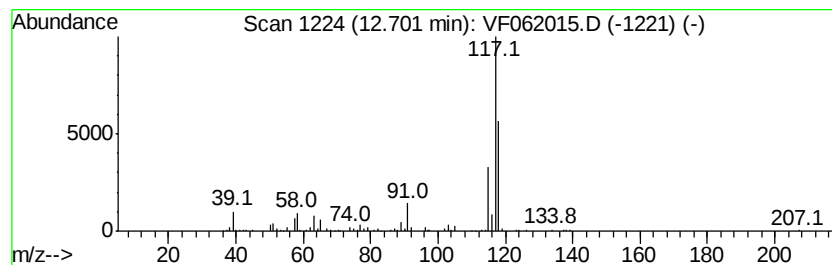
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Indane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	14.45 ug/l	168030	1,4-Dichlorobenzene-d4	12.53

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	91
2	Benzene, 2-propenyl-		118	C9H10	000300-57-2	83
3	Benzene, 1-ethenyl-2-methyl-		118	C9H10	000611-15-4	72
4	Benzene, cyclopropyl-		118	C9H10	000873-49-4	68
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...		118	C9H10	1000191-13-7	64



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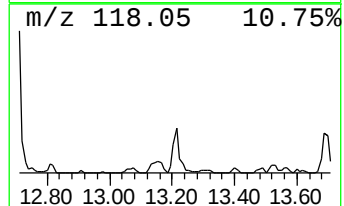
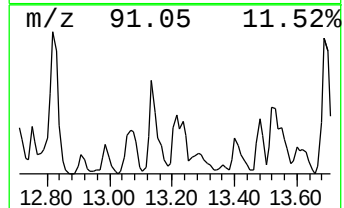
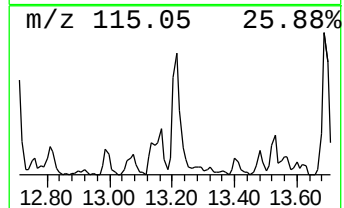
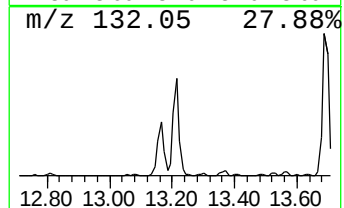
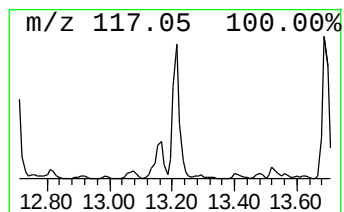
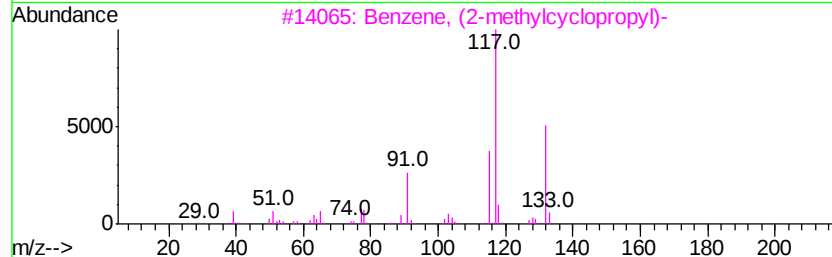
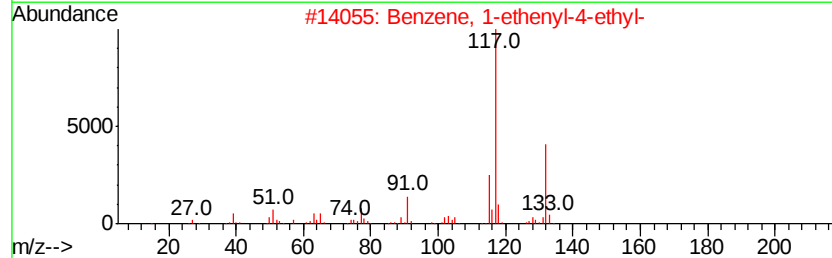
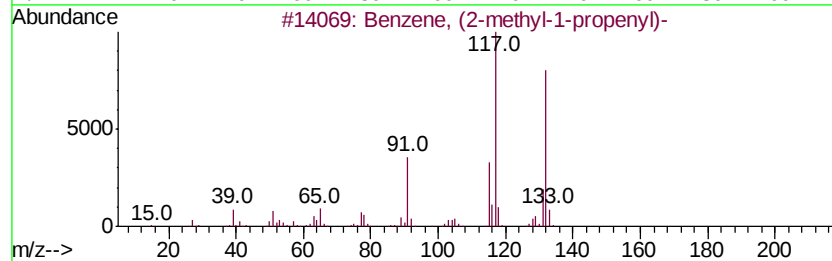
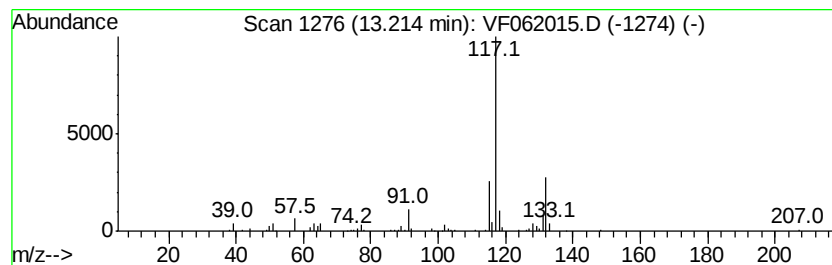
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, (2-methyl-1-propen... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.21	14.74 ug/l	171467	1,4-Dichlorobenzene-d4	12.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
3		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
4		Indan, 1-methyl-	132	C10H12	000767-58-8	87
5		Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	83



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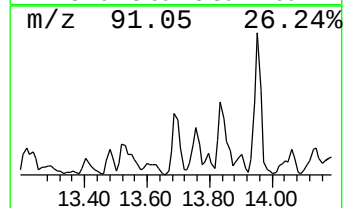
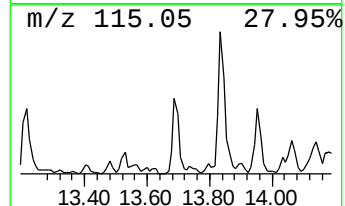
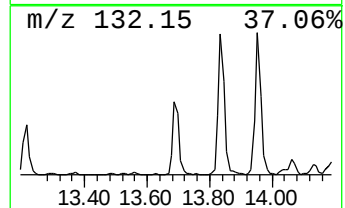
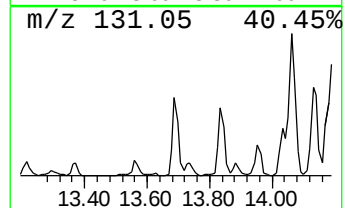
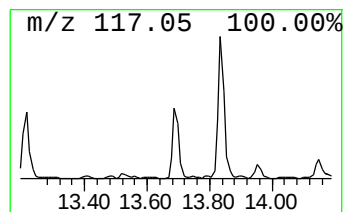
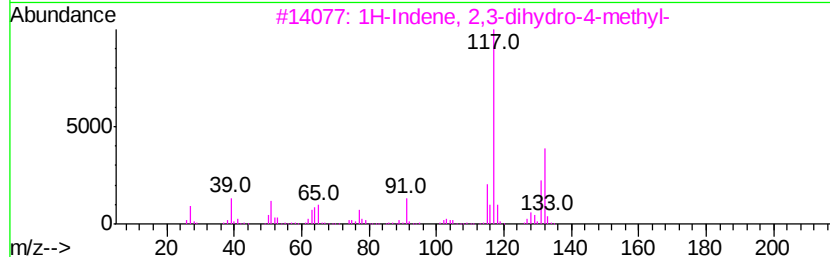
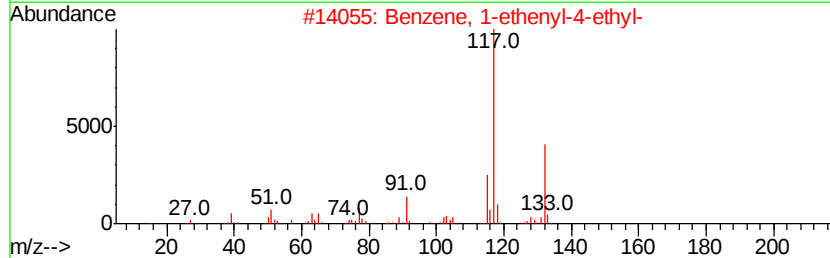
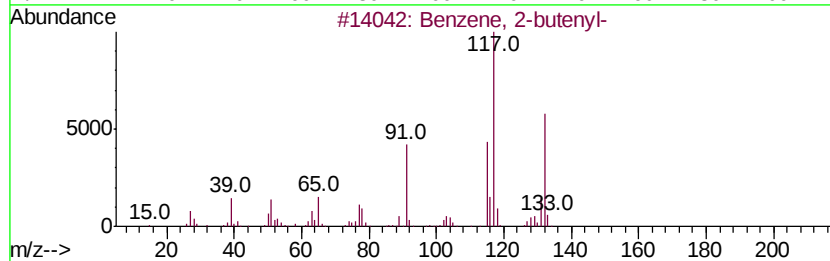
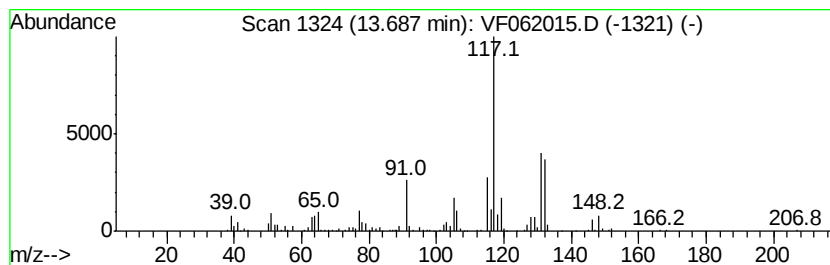
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 2-butenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.69	26.71 ug/l	310685	1,4-Dichlorobenzene-d4	12.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	83
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	58
4		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	58
5		Indan, 1-methyl-	132	C10H12	000767-58-8	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF031919\
Data File : VF062015.D
Acq On : 19 Mar 2019 13:47
Operator : VA/AP
Sample : K1693-09
Misc : 6.50µ/5mL/MSVOA-F/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
OR-03-031819-A

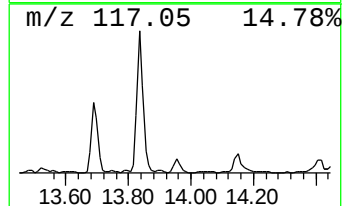
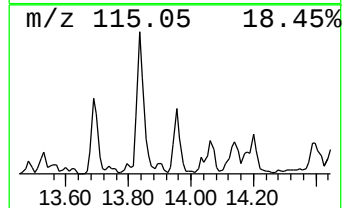
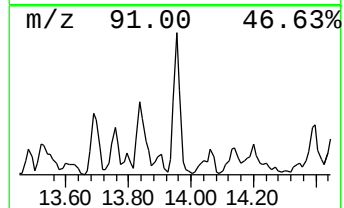
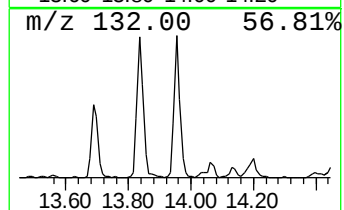
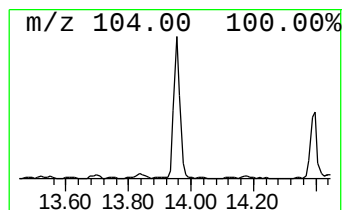
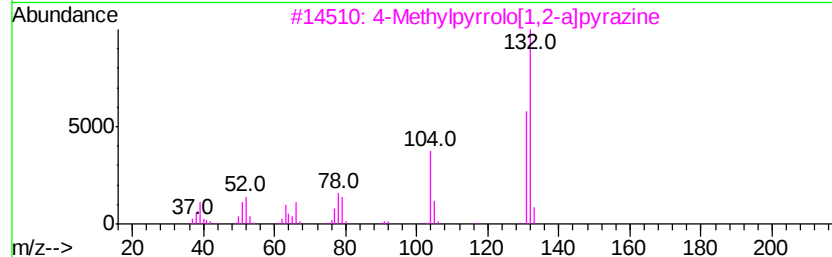
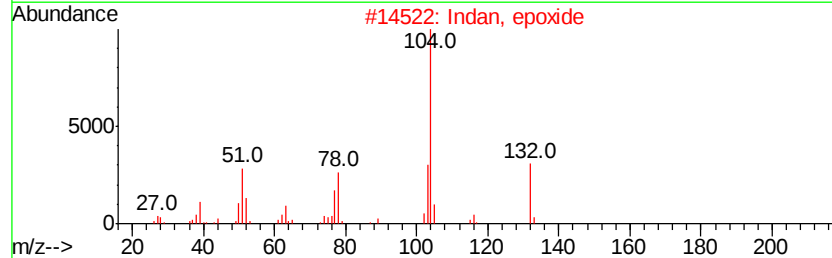
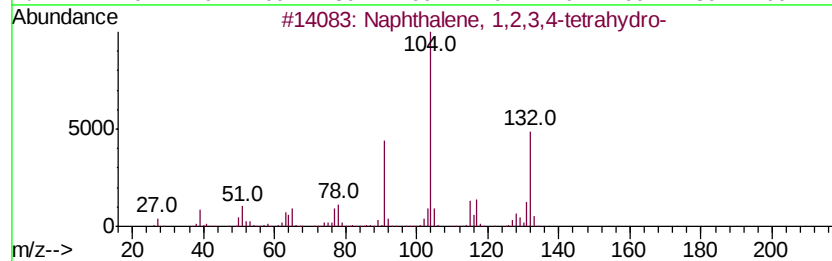
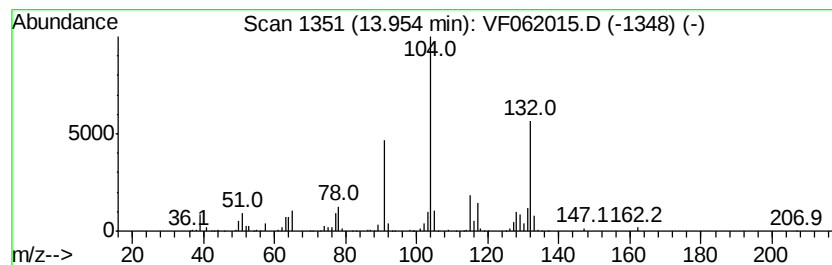
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Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	24.36 ug/l	283263	1,4-Dichlorobenzene-d4	12.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	97
2		Indan, epoxide	132	C9H8O	000768-22-9	58
3		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	53
4		Benzene, cyclobutyl-	132	C10H12	004392-30-7	50
5		Styrene	104	C8H8	000100-42-5	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF031919\
Data File : VF062015.D
Acq On : 19 Mar 2019 13:47
Operator : VA/AP
Sample : K1693-09
Misc : 6.50g/5mL/MSVOA-F/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
OR-03-031819-A

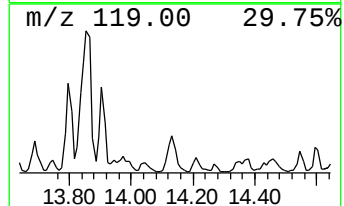
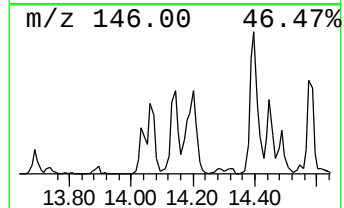
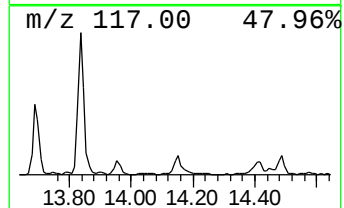
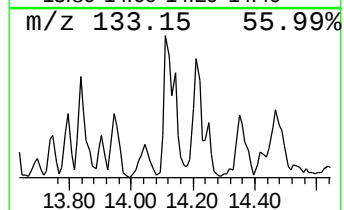
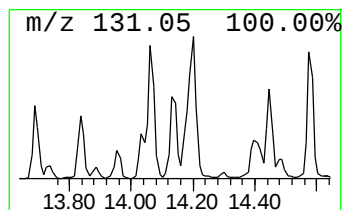
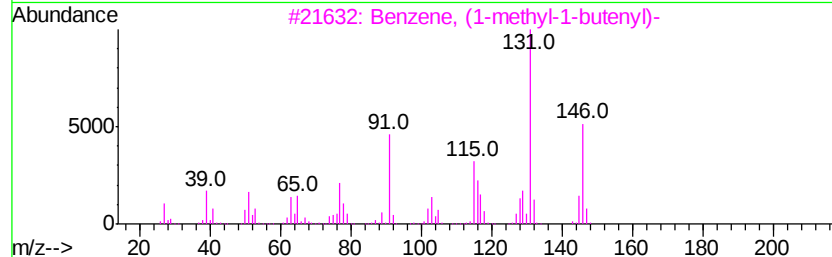
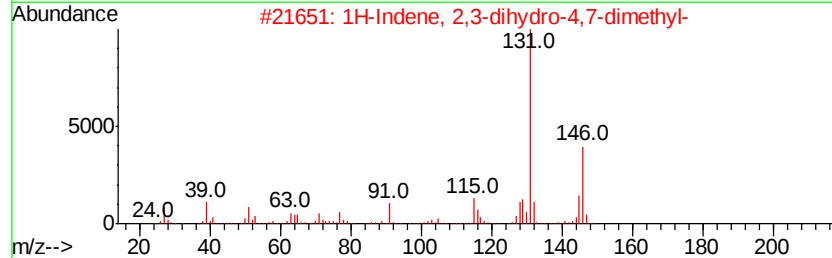
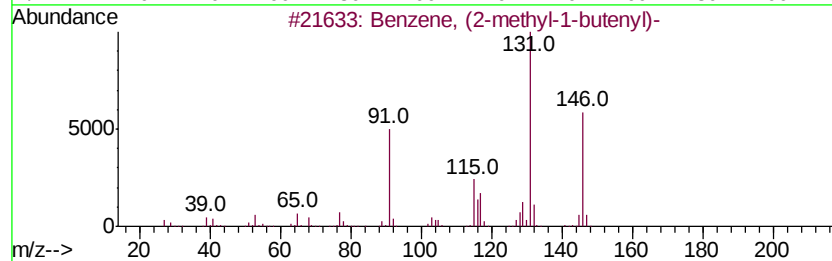
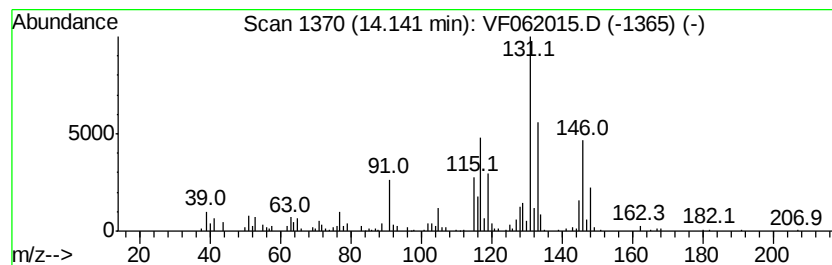
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, (2-methyl-1-butenyl)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.14	18.12 ug/l	210735	1,4-Dichlorobenzene-d4	12.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	90
2		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87
3		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	78
4		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	55
5		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	55



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF031919\
Data File : VF062015.D
Acq On : 19 Mar 2019 13:47
Operator : VA/AP
Sample : K1693-09
Misc : 6.50g/5mL/MSVOA-F/SOIL
ALS Vial : 7 Sample Multiplier: 1

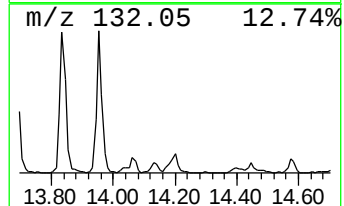
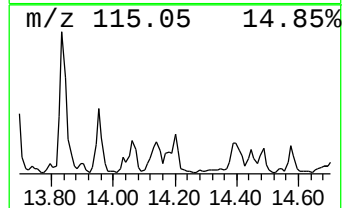
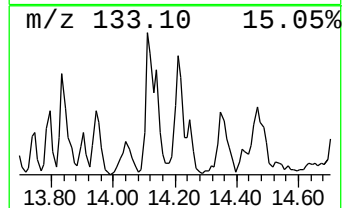
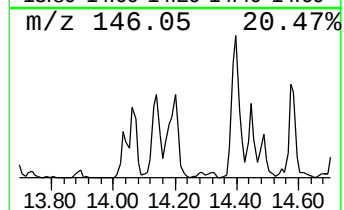
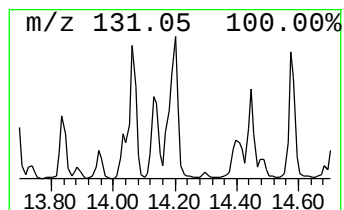
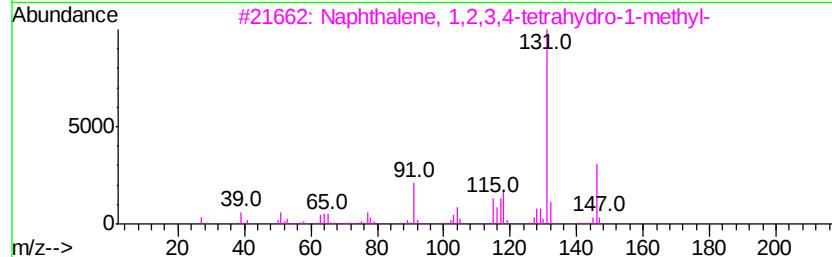
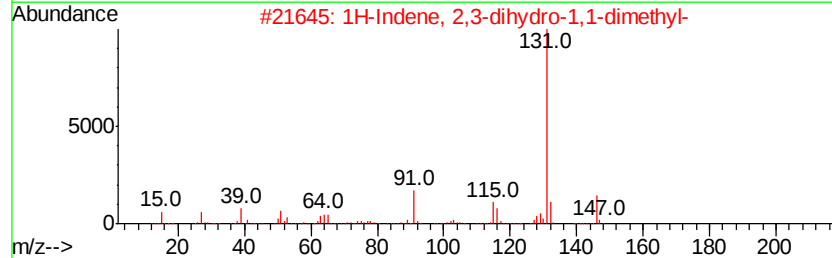
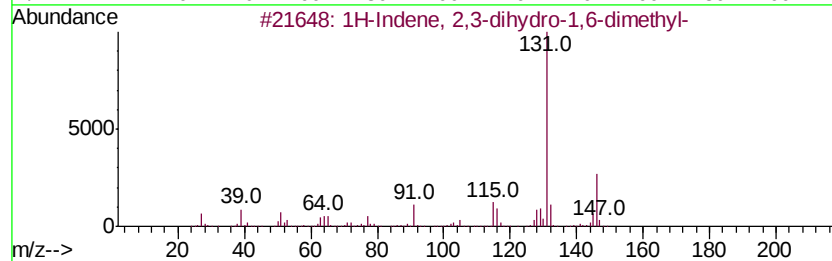
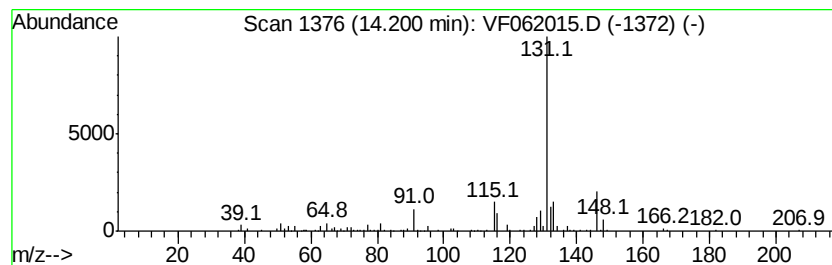
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ClientSampleId :
OR-03-031819-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F031619S.M
Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.20	18.80 ug/l	218628	1,4-Dichlorobenzene-d4	12.53	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	87
2	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	83
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	80
4	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	80
5	3-Phenyl-4,5-dimethyl-2,1-oxabor...	202	C13H19BO	1000062-26-1	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF031919\
Data File : VF062015.D
Acq On : 19 Mar 2019 13:47
Operator : VA/AP
Sample : K1693-09
Misc : 6.50g/5mL/MSVOA-F/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
OR-03-031819-A

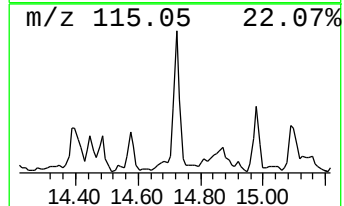
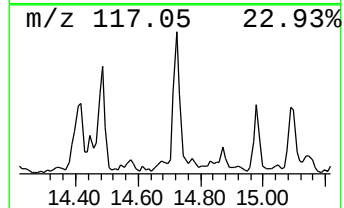
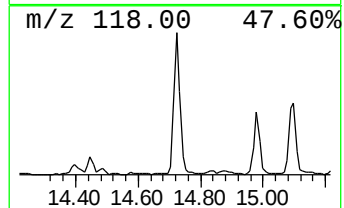
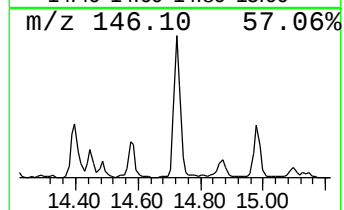
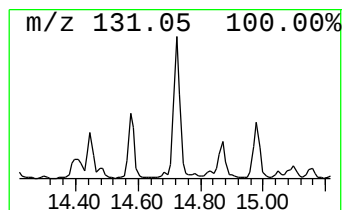
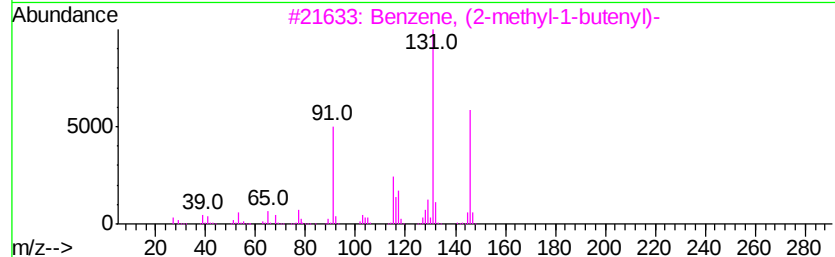
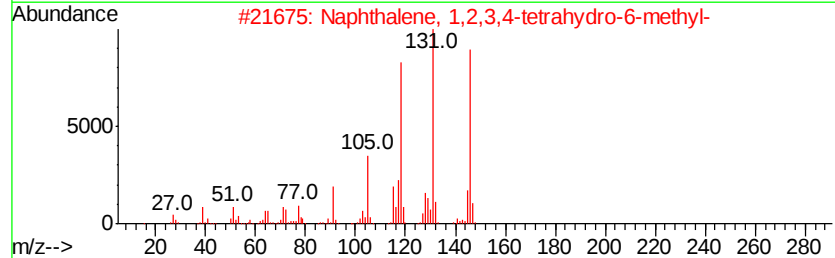
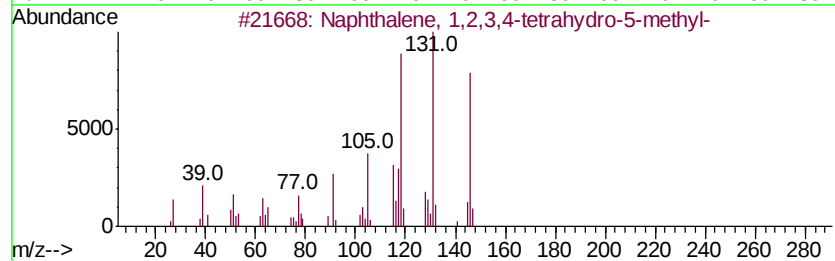
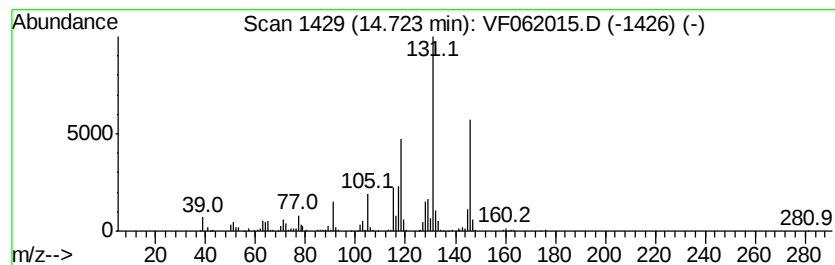
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F031619S.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.72	31.62 ug/l	367696	1,4-Dichlorobenzene-d4	12.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	96
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	93
3		Benzene, (2-methyl-1-butenvl)-	146	C11H14	056253-64-6	76
4		Benzene, 2-ethenvl-1,3,5-trimethyl-	146	C11H14	000769-25-5	70
5		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF031919\
Data File : VF062015.D
Acq On : 19 Mar 2019 13:47
Operator : VA/AP
Sample : K1693-09
Misc : 6.50g/5mL/MSVOA-F/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
OR-03-031819-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F031619S.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-...	12.59	16.5	ug/l	192290	4	12.53	581514	50.0
Indane	12.70	14.4	ug/l	168030	4	12.53	581514	50.0
Benzene, (2-methy...	13.21	14.7	ug/l	171467	4	12.53	581514	50.0
Benzene, 2-butenyl-	13.69	26.7	ug/l	310685	4	12.53	581514	50.0
Naphthalene, 1,2,...	13.95	24.4	ug/l	283263	4	12.53	581514	50.0
Benzene, (2-methy...	14.14	18.1	ug/l	210735	4	12.53	581514	50.0
1H-Indene, 2,3-di...	14.20	18.8	ug/l	218628	4	12.53	581514	50.0
Naphthalene, 1,2,...	14.72	31.6	ug/l	367696	4	12.53	581514	50.0