

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD110118\
 Data File : VD060270.D
 Acq On : 1 Nov 2018 18:57
 Operator : VA/AP
 Sample : J5200-19
 Misc : 6.83g/5ml/MSVOA D/SOIL
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 OR-03-103118-B

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

Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D100118S.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	2.822	128	132	138	rBV4	6375	18254	1.16%	0.104%
2	3.147	161	165	172	rBV	30372	81250	5.16%	0.464%
3	3.708	218	222	234	rVB2	21065	63379	4.03%	0.362%
4	5.617	411	416	426	rVB3	12934	46162	2.93%	0.264%
5	6.316	482	487	492	rBV	279173	728723	46.30%	4.160%
6	6.404	492	496	513	rVB	376136	950234	60.37%	5.424%
7	6.788	529	535	549	rVB	279466	716771	45.54%	4.092%
8	7.438	596	601	613	rBV	568091	1314820	83.54%	7.505%
9	8.038	657	662	668	rVB3	7373	18074	1.15%	0.103%
10	9.455	801	806	812	rBV	644592	1357809	86.27%	7.751%
11	9.554	812	816	824	rVB	117955	258597	16.43%	1.476%
12	10.193	876	881	889	rBV	39573	99378	6.31%	0.567%
13	11.217	982	985	988	rBV	24450	21929	1.39%	0.125%
14	11.286	988	992	1000	rVV	1020452	1573919	100.00%	8.984%
15	11.433	1004	1007	1013	rVV	22922	41874	2.66%	0.239%
16	11.561	1013	1020	1025	rVV	95385	188913	12.00%	1.078%
17	11.630	1025	1027	1033	rVB2	21660	39607	2.52%	0.226%
18	11.896	1050	1054	1056	rBV2	10032	19335	1.23%	0.110%
19	11.935	1056	1058	1065	rVB	59906	90951	5.78%	0.519%
20	12.092	1068	1074	1078	rBV4	10505	27641	1.76%	0.158%
21	12.191	1078	1084	1088	rVB3	18491	38329	2.44%	0.219%
22	12.279	1090	1093	1096	rVB2	12626	17010	1.08%	0.097%
23	12.427	1105	1108	1113	rBV	744392	1010343	64.19%	5.767%
24	12.585	1120	1124	1127	rVB2	13874	32225	2.05%	0.184%
25	12.644	1127	1130	1132	rVB	25980	35066	2.23%	0.200%
26	12.712	1132	1137	1139	rBV	89304	131013	8.32%	0.748%
27	12.742	1139	1140	1143	rVV	37442	41921	2.66%	0.239%
28	12.791	1143	1145	1147	rVV2	40929	72554	4.61%	0.414%
29	12.821	1147	1148	1152	rVB	71906	77322	4.91%	0.441%
30	12.949	1156	1161	1163	rBV	62803	89887	5.71%	0.513%
31	13.057	1169	1172	1173	rBV	14313	18853	1.20%	0.108%
32	13.096	1173	1176	1182	rVB	200189	248147	15.77%	1.416%
33	13.224	1185	1189	1193	rVB2	26076	40262	2.56%	0.230%
34	13.293	1193	1196	1200	rBV	50733	66189	4.21%	0.378%

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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

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35	13.372	1201	1204	1207	rBV	1104067	1419945	90.22%	8.105%
36	13.411	1207	1208	1212	rVB	105620	124314	7.90%	0.710%
37	13.500	1212	1217	1220	rVB2	19746	38769	2.46%	0.221%
38	13.559	1220	1223	1225	rBV	102295	167099	10.62%	0.954%
39	13.588	1225	1226	1228	rVB	81639	59223	3.76%	0.338%
40	13.628	1228	1230	1236	rBV3	110958	220849	14.03%	1.261%
41	13.736	1238	1241	1242	rBV	65366	80519	5.12%	0.460%
42	13.765	1242	1244	1248	rVB	56594	68630	4.36%	0.392%
43	13.824	1248	1250	1252	rBV	51202	73912	4.70%	0.422%
44	13.854	1252	1253	1256	rVB	54020	46254	2.94%	0.264%
45	13.903	1256	1258	1261	rVB	85987	91198	5.79%	0.521%
46	13.992	1264	1267	1272	rVB3	79172	154618	9.82%	0.883%
47	14.070	1272	1275	1278	rVB3	31413	51448	3.27%	0.294%
48	14.130	1278	1281	1283	rBV2	46620	92521	5.88%	0.528%
49	14.169	1283	1285	1286	rBV	15813	18974	1.21%	0.108%
50	14.189	1286	1287	1289	rVB	31106	34068	2.16%	0.194%
51	14.228	1289	1291	1294	rVB	82065	91703	5.83%	0.523%
52	14.277	1294	1296	1298	rBV2	66440	83284	5.29%	0.475%
53	14.317	1298	1300	1303	rVB3	26689	43458	2.76%	0.248%
54	14.415	1303	1310	1312	rBV3	98615	213807	13.58%	1.220%
55	14.454	1312	1314	1315	rVV	45500	53188	3.38%	0.304%
56	14.484	1315	1317	1319	rVV2	61277	121981	7.75%	0.696%
57	14.523	1319	1321	1325	rVV2	258816	363689	23.11%	2.076%
58	14.582	1325	1327	1330	rVB2	53330	67976	4.32%	0.388%
59	14.641	1330	1333	1337	rVB	206449	281623	17.89%	1.608%
60	14.749	1337	1344	1346	rBV2	68248	207285	13.17%	1.183%
61	14.779	1346	1347	1349	rVV	69055	54796	3.48%	0.313%
62	14.838	1349	1353	1357	rVB2	67580	139306	8.85%	0.795%
63	14.966	1363	1366	1370	rBV	319982	517227	32.86%	2.952%
64	15.025	1370	1372	1374	rVV	112285	150541	9.56%	0.859%
65	15.074	1374	1377	1379	rVV2	147515	235139	14.94%	1.342%
66	15.153	1382	1385	1388	rVV4	89897	153390	9.75%	0.876%
67	15.202	1388	1390	1394	rVV2	108882	162804	10.34%	0.929%
68	15.271	1394	1397	1399	rVV	22090	41863	2.66%	0.239%
69	15.350	1403	1405	1410	rVV	264283	408080	25.93%	2.329%
70	15.429	1410	1413	1415	rVV3	49519	89912	5.71%	0.513%
71	15.478	1415	1418	1420	rVV	87418	147353	9.36%	0.841%

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72	15.517	1420	1422	1424	rVV	32885	43516	2.76%	0.248%
73	15.586	1425	1429	1431	rVV	174733	254652	16.18%	1.454%
74	15.645	1431	1435	1438	rVV5	23108	67091	4.26%	0.383%
75	15.724	1438	1443	1445	rVV3	147973	276351	17.56%	1.577%
76	15.773	1445	1448	1453	rVV	188332	348282	22.13%	1.988%
77	15.842	1453	1455	1457	rVV2	41510	53499	3.40%	0.305%
78	15.911	1459	1462	1465	rVV	171646	274174	17.42%	1.565%
79	15.960	1465	1467	1469	rVV2	37543	53224	3.38%	0.304%
80	16.019	1469	1473	1477	rVV2	40092	84327	5.36%	0.481%
81	16.068	1477	1478	1483	rVB4	7102	16506	1.05%	0.094%
82	16.167	1483	1488	1493	rVB6	10910	26698	1.70%	0.152%
83	16.314	1498	1503	1505	rBV2	27464	51303	3.26%	0.293%
84	16.344	1505	1506	1512	rVB2	13683	21365	1.36%	0.122%

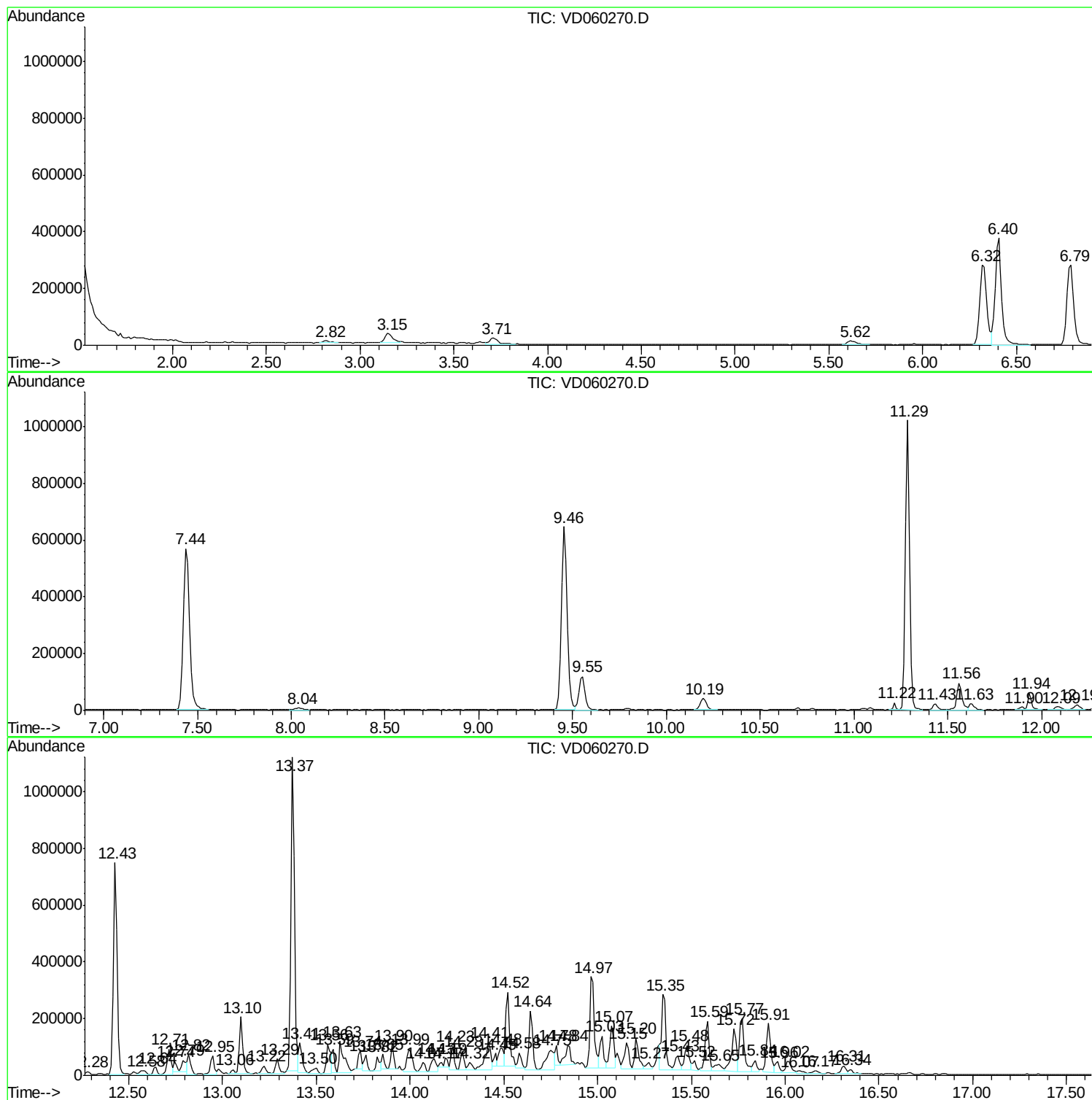
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D100118S.M
Quant Title : SW846 8260

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



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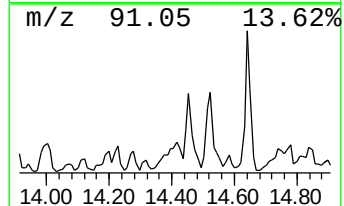
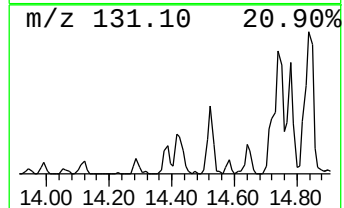
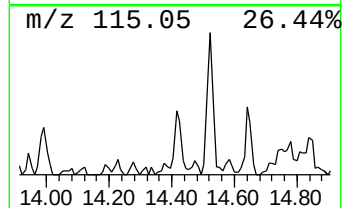
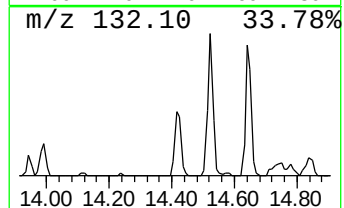
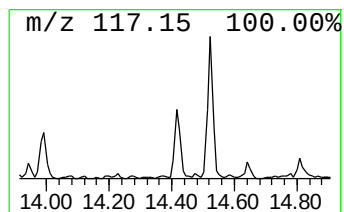
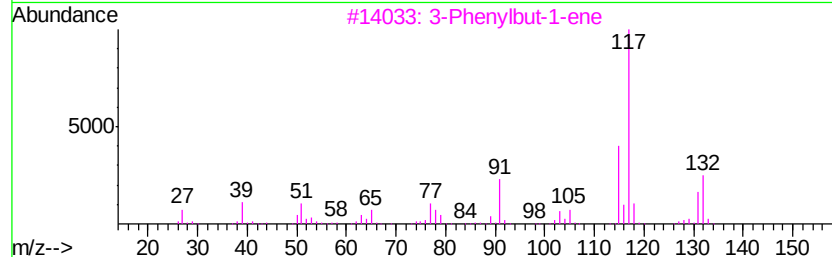
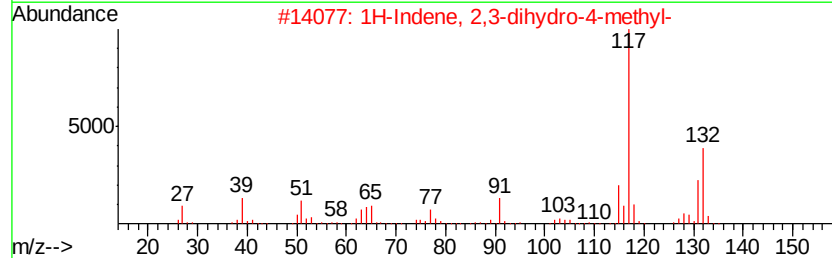
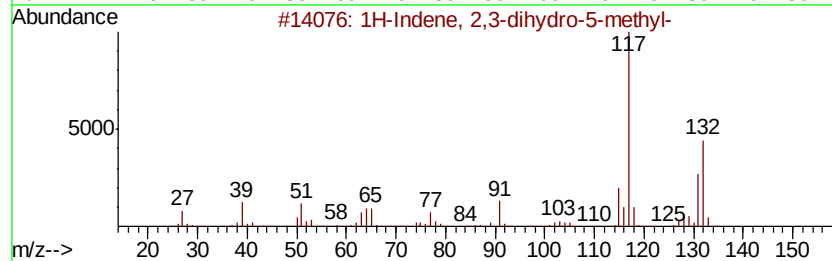
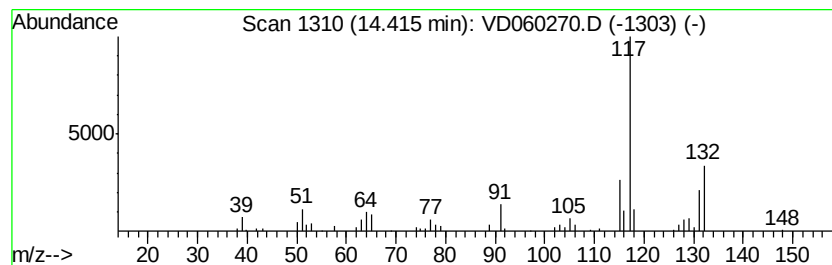
Instrument :
MSVOA_D
ClientSampleId :
OR-03-103118-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D100118S.M
Quant Title : SW846 8260

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

Peak Number 1 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.41	7.53 ug/l	213807	1,4-Dichlorobenzene-d4	13.37	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	83
4	Indan, 1-methyl-	132	C10H12	000767-58-8	81
5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	74



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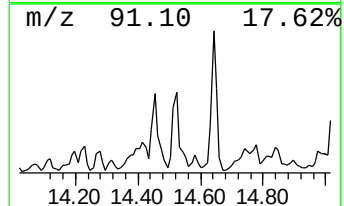
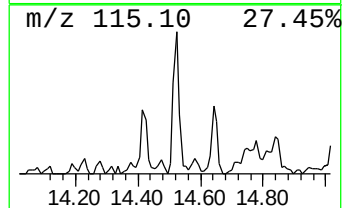
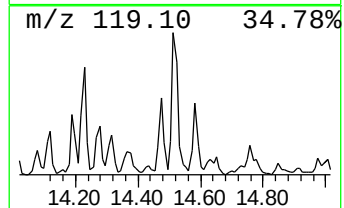
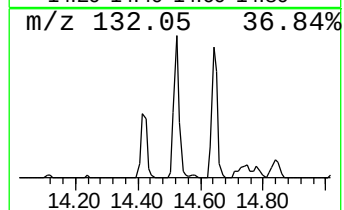
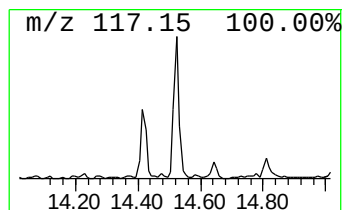
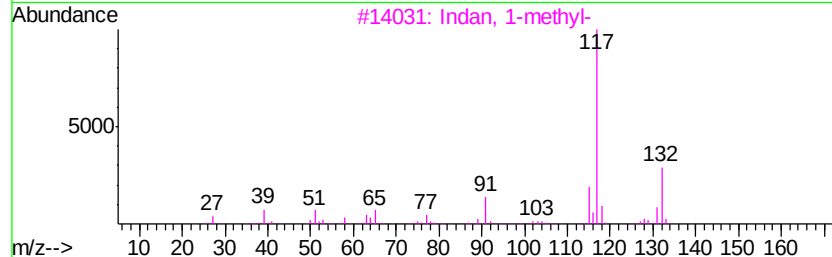
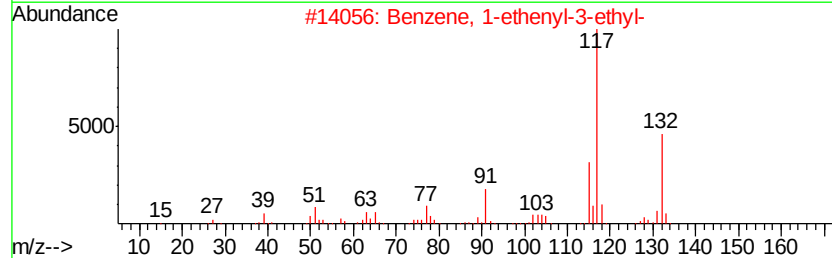
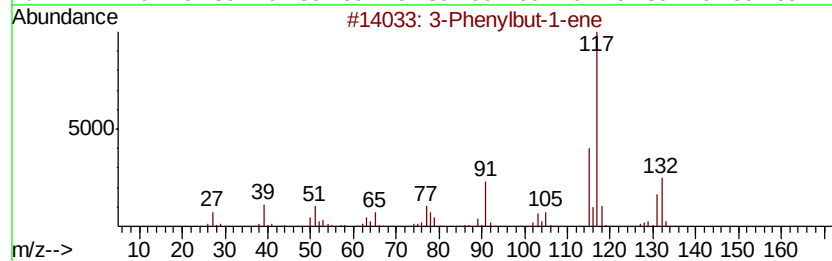
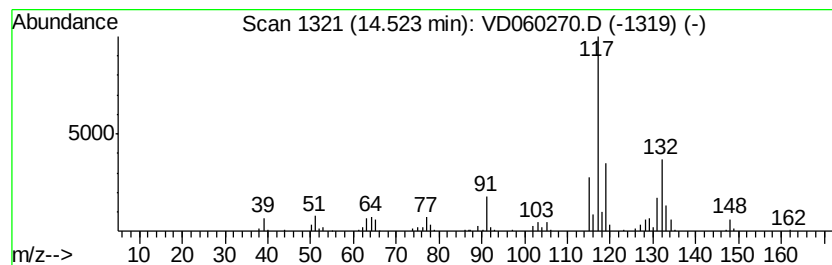
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D100118S.M
Quant Title : SW846 8260

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

Peak Number 2 3-Phenylbut-1-ene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.52	12.81 ug/l	363689	1,4-Dichlorobenzene-d4	13.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	76
3		Indan, 1-methyl-	132	C10H12	000767-58-8	76
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	72



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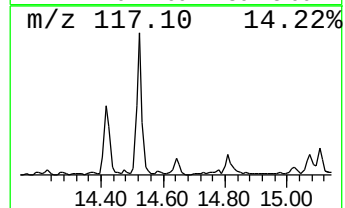
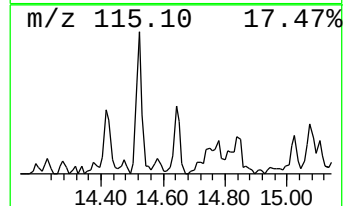
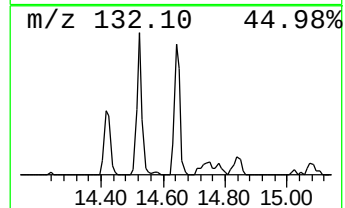
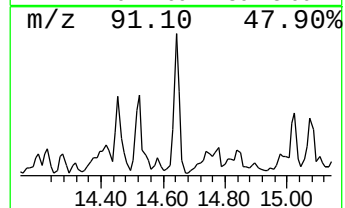
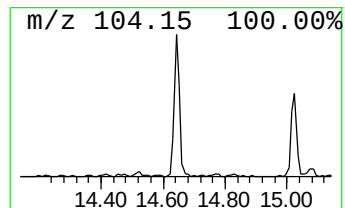
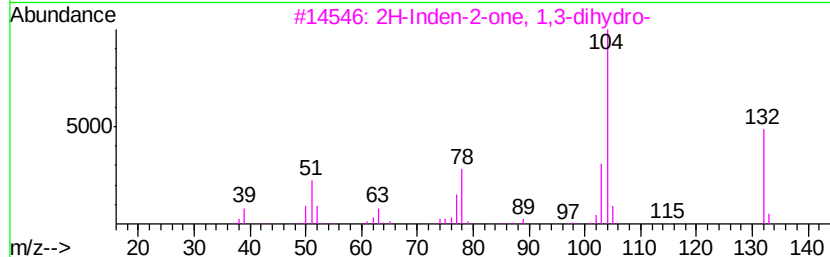
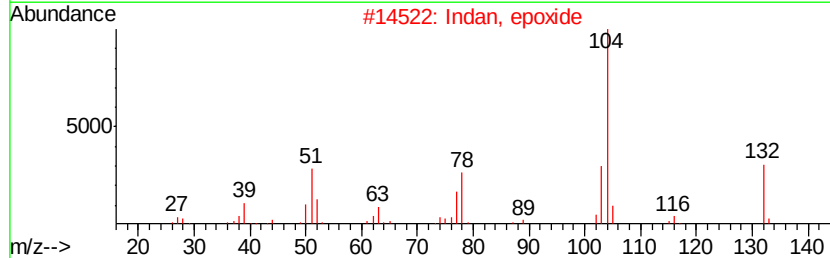
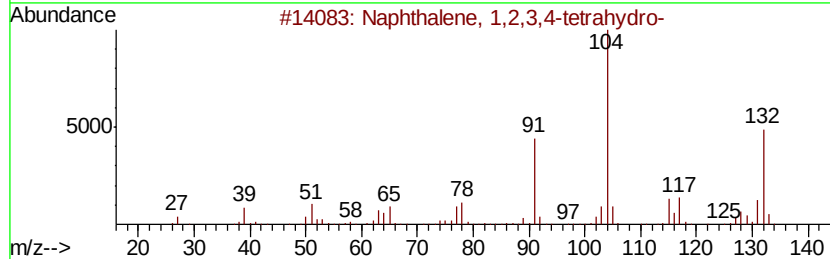
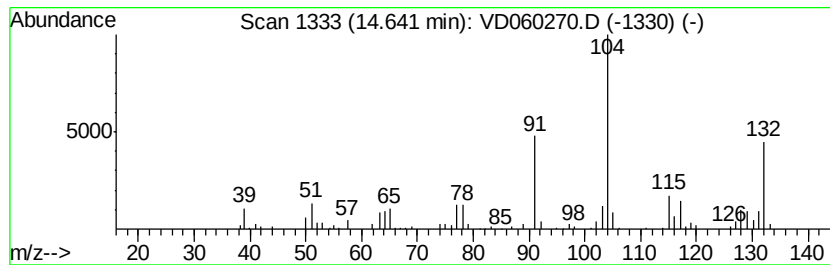
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.64	9.92 ug/l	281623	1,4-Dichlorobenzene-d4	13.37		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90
2		Indan, epoxide	132	C9H8O	000768-22-9	58
3		2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	50
4		Acetic acid, trichloro-, 2-pheny...	266	C10H9Cl3O2	071965-07-6	43
5		N-Ethylbenzimidoyl bromide	211	C9H10BrN	041182-87-0	40



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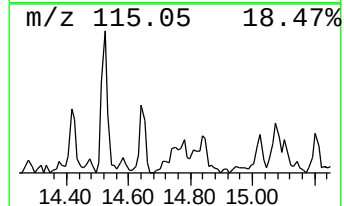
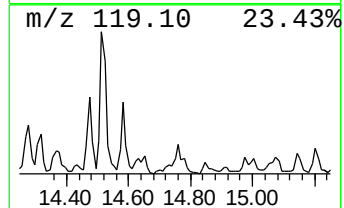
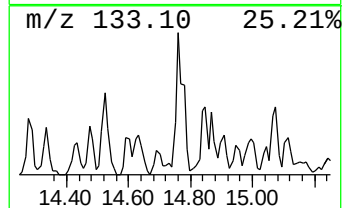
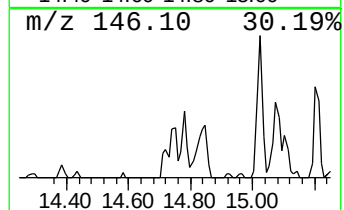
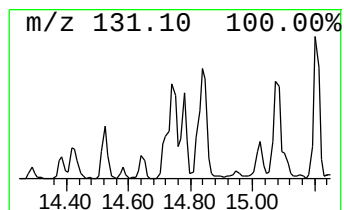
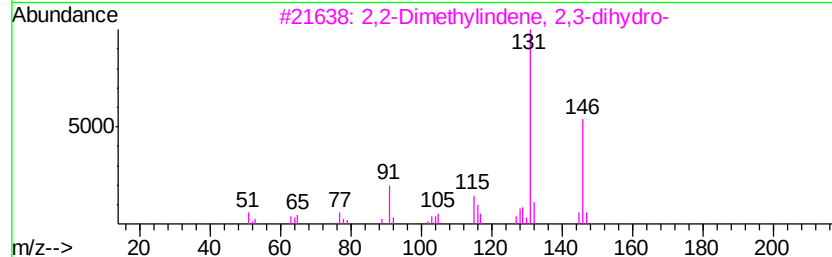
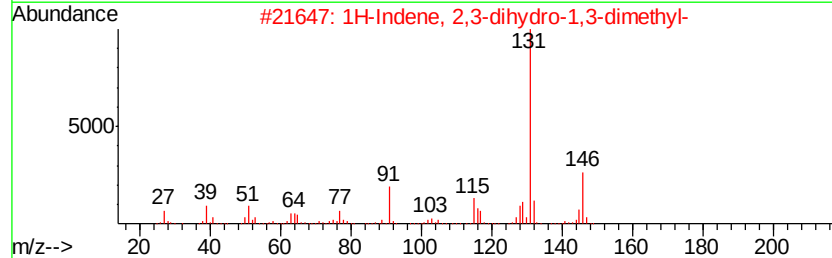
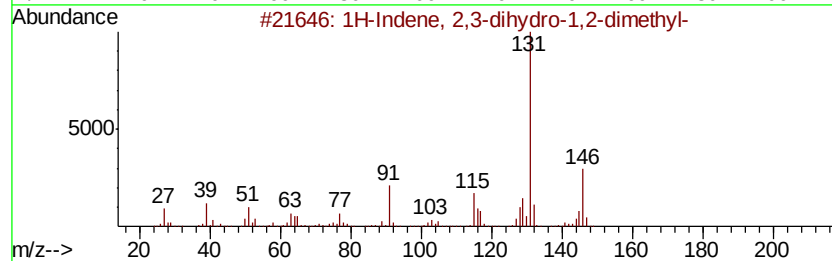
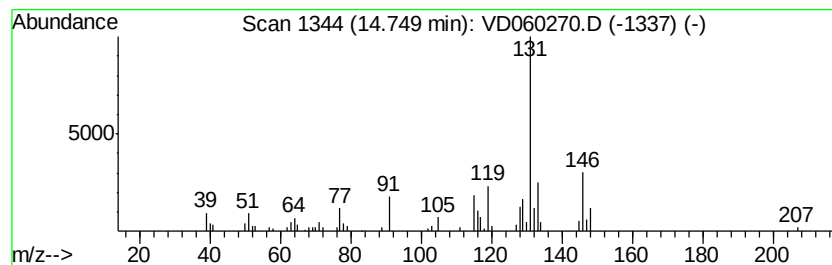
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MSVOA_D
ClientSampleId :
OR-03-103118-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D100118S.M
Quant Title : SW846 8260

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

Peak Number 4 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.75	7.30 ug/l	207285	1,4-Dichlorobenzene-d4	13.37
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8	93
2	1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5	92
3	2,2-Dimethylindene, 2,3-dihydro-	146 C11H14	020836-11-7	89
4	Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3	76
5	Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD110118\
Data File : VD060270.D
Acq On : 1 Nov 2018 18:57
Operator : VA/AP
Sample : J5200-19
Misc : 6.83g/5ml/MSVOA D/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
OR-03-103118-B

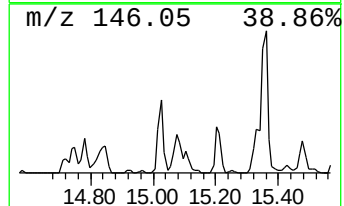
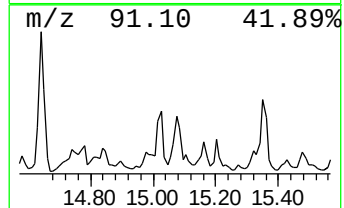
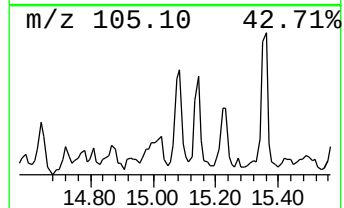
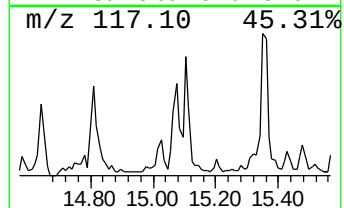
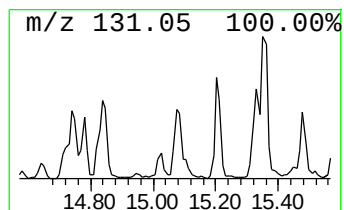
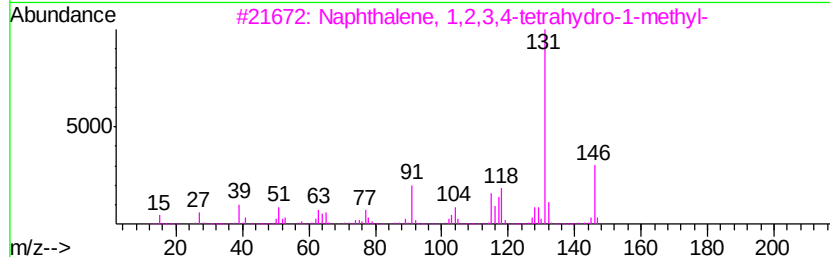
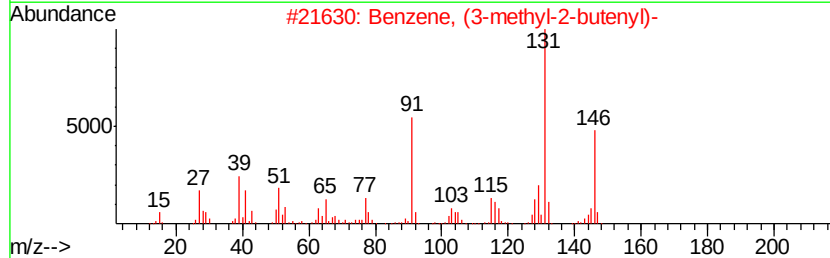
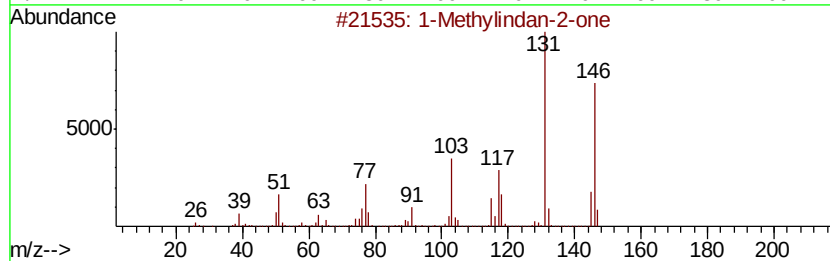
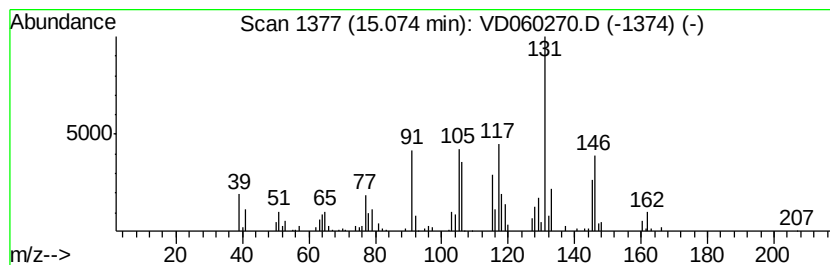
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Quant Title : SW846 8260

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

Peak Number 5 1-Methylindan-2-one Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	8.28 ug/l	235139	1,4-Dichlorobenzene-d4	13.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Methylindan-2-one	146	C10H10O	035587-60-1	70
2		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70
3		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	60
4		2-Propenal, 3-(4-methylphenyl)-	146	C10H10O	001504-75-2	53
5		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD110118\
Data File : VD060270.D
Acq On : 1 Nov 2018 18:57
Operator : VA/AP
Sample : J5200-19
Misc : 6.83g/5ml/MSVOA D/SOIL
ALS Vial : 13 Sample Multiplier: 1

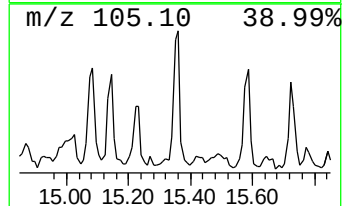
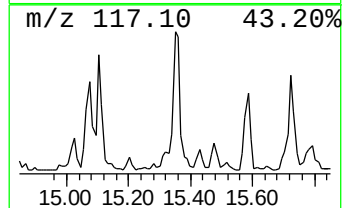
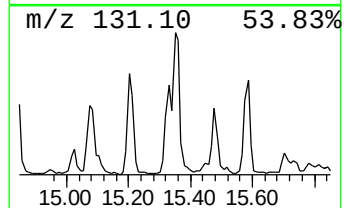
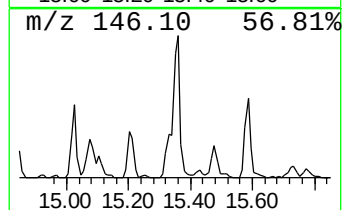
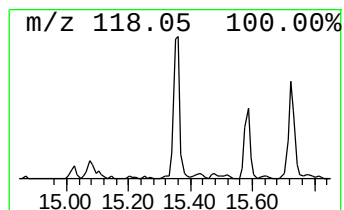
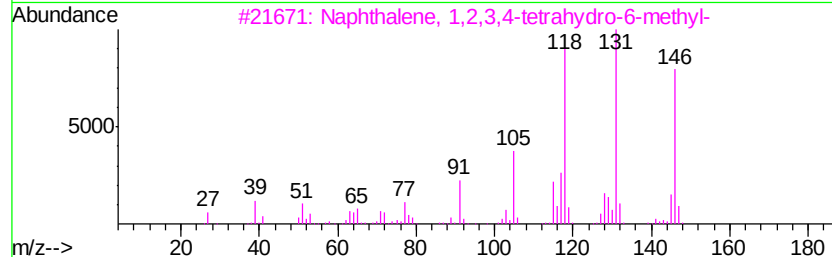
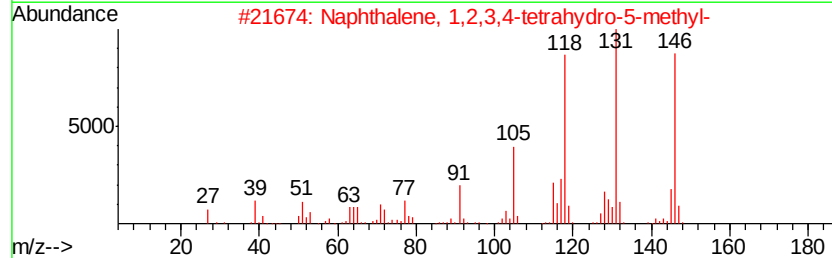
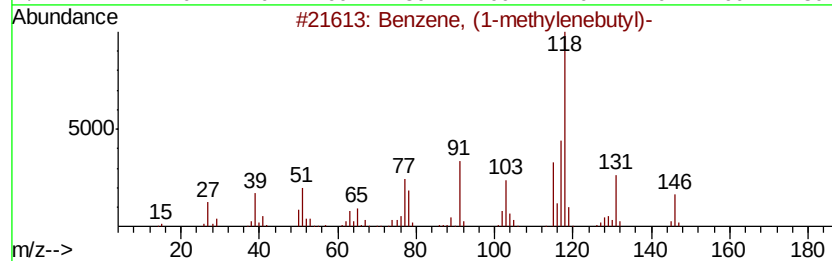
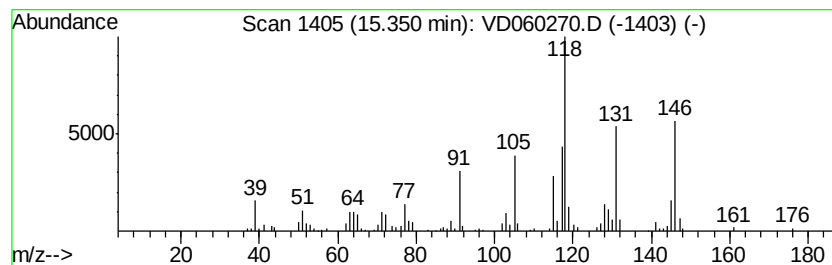
Instrument :
MSVOA_D
ClientSampleId :
OR-03-103118-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D100118S.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, (1-methylenebutyl)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.35	14.37 ug/l	408080	1,4-Dichlorobenzene-d4	13.37
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1-methylenebutyl)-	146 C11H14	005676-32-4 91
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 91
3		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 81
4		Benzene, (1-ethyl-2-propenyl)-	146 C11H14	019947-22-9 80
5		2,3,4,5,6,7-Hexahydro-1H-cyclope...	146 C11H14	1000189-31-0 68



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD110118\
Data File : VD060270.D
Acq On : 1 Nov 2018 18:57
Operator : VA/AP
Sample : J5200-19
Misc : 6.83g/5ml/MSVOA D/SOIL
ALS Vial : 13 Sample Multiplier: 1

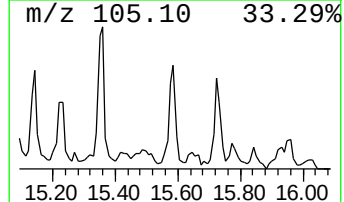
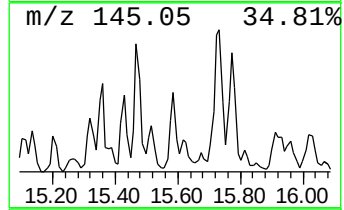
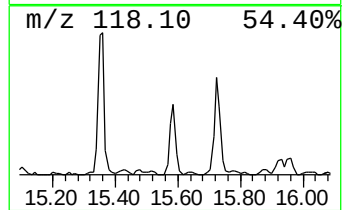
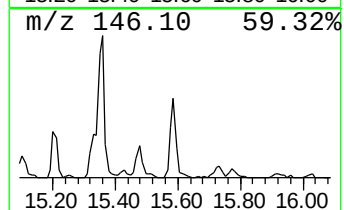
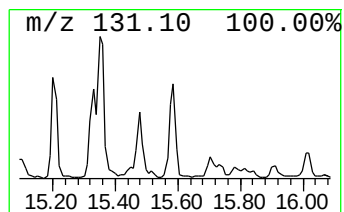
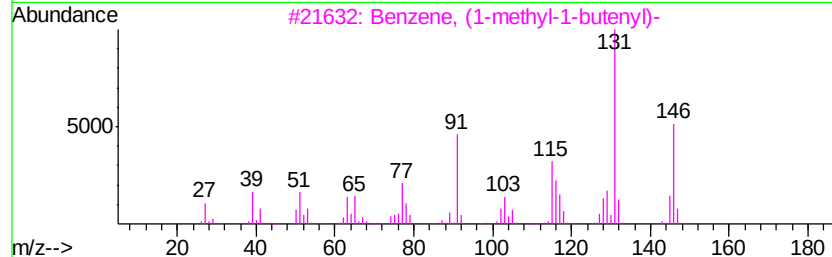
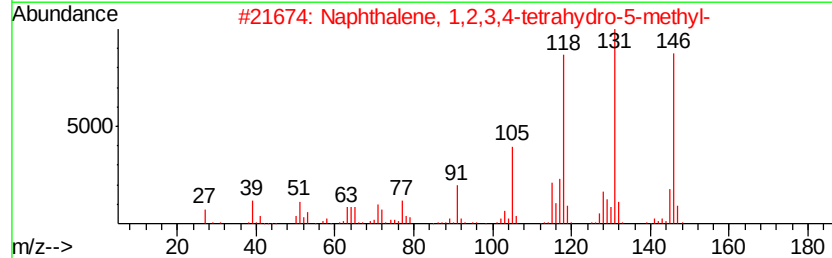
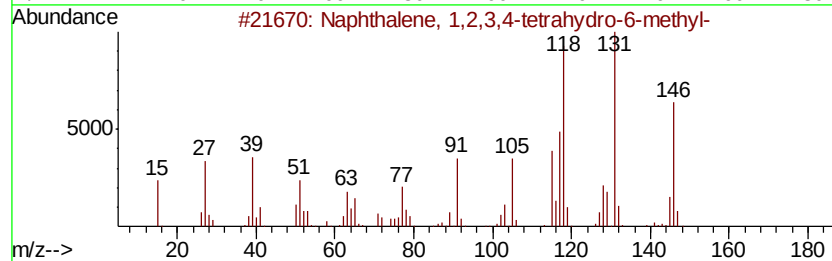
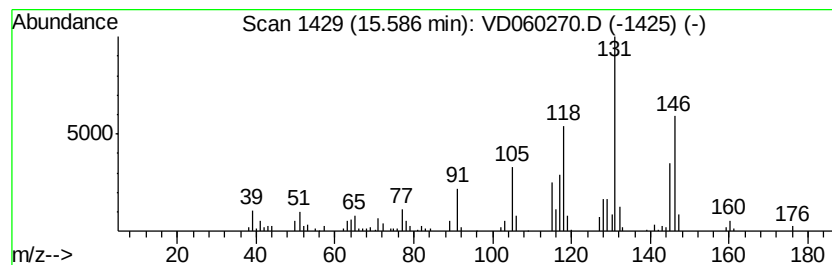
Instrument :
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ClientSampleId :
OR-03-103118-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D100118S.M
Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetra... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.59	8.97 ug/l	254652	1,4-Dichlorobenzene-d4	13.37
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 94
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 93
3		Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2 76
4		2-Propenal, 3-(4-methylphenyl)-	146 C10H10O	001504-75-2 76
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD110118\
Data File : VD060270.D
Acq On : 1 Nov 2018 18:57
Operator : VA/AP
Sample : J5200-19
Misc : 6.83g/5ml/MSVOA D/SOIL
ALS Vial : 13 Sample Multiplier: 1

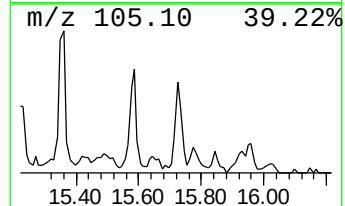
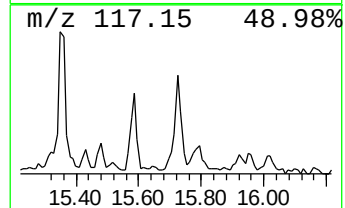
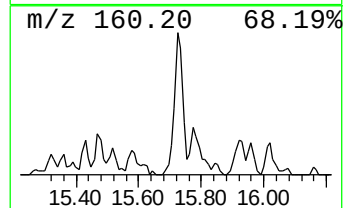
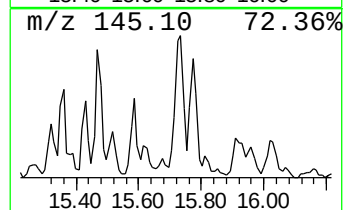
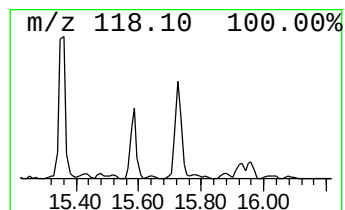
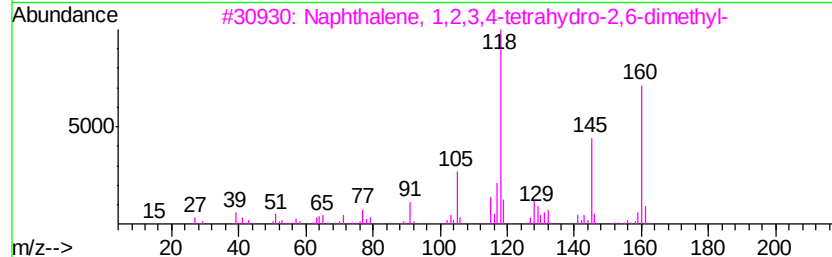
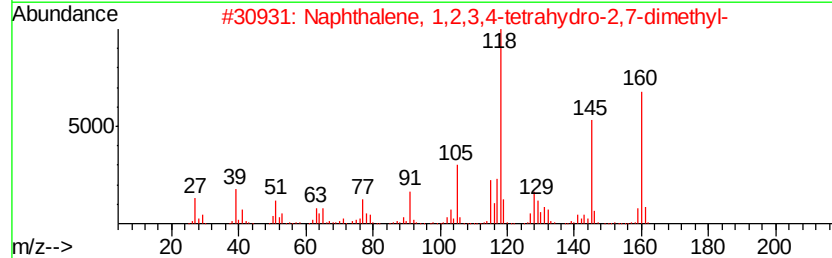
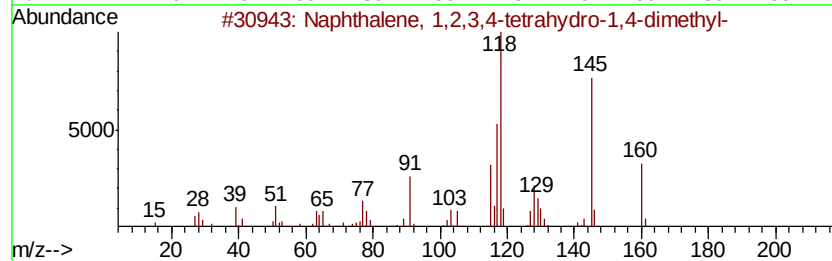
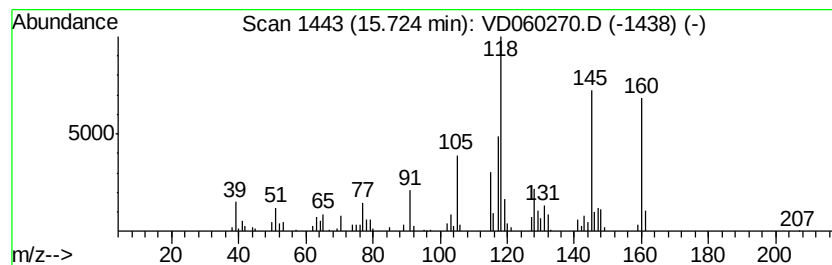
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D100118S.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetra... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.72	9.73 ug/l	276351	1,4-Dichlorobenzene-d4	13.37
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	004175-54-6 93
2		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	013065-07-1 81
3		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	007524-63-2 81
4		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	021564-91-0 64
5		Benzene, 1,3,5-trimethyl-2-(1-me...	160 C12H16	014679-13-1 50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD110118\
Data File : VD060270.D
Acq On : 1 Nov 2018 18:57
Operator : VA/AP
Sample : J5200-19
Misc : 6.83g/5ml/MSVOA D/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
OR-03-103118-B

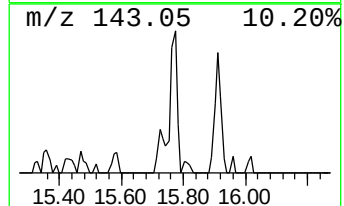
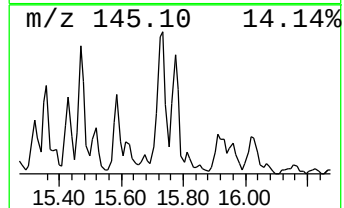
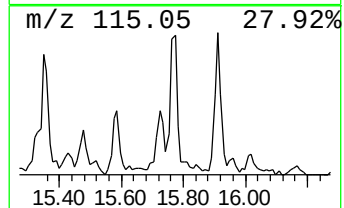
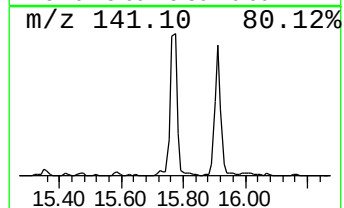
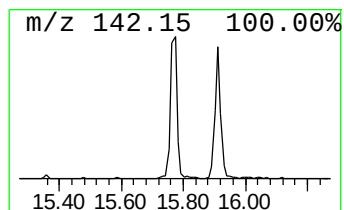
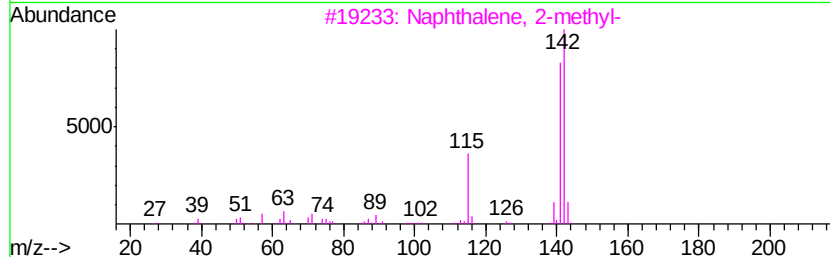
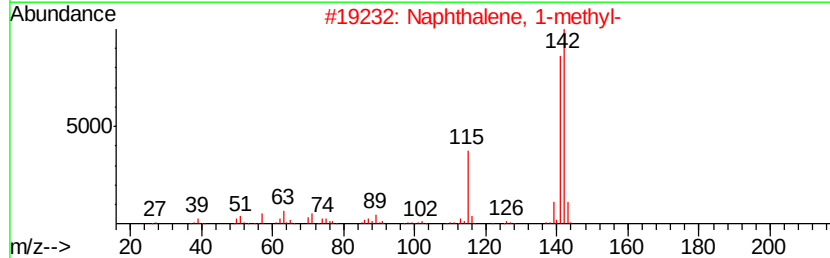
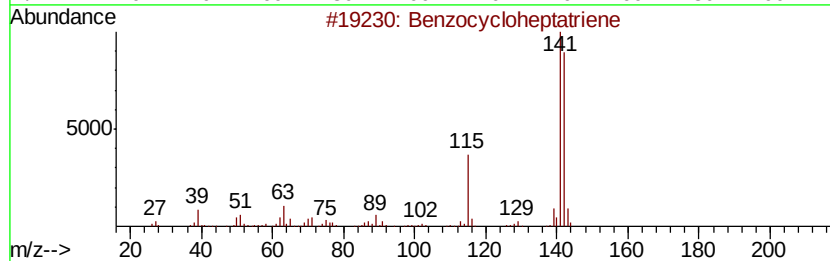
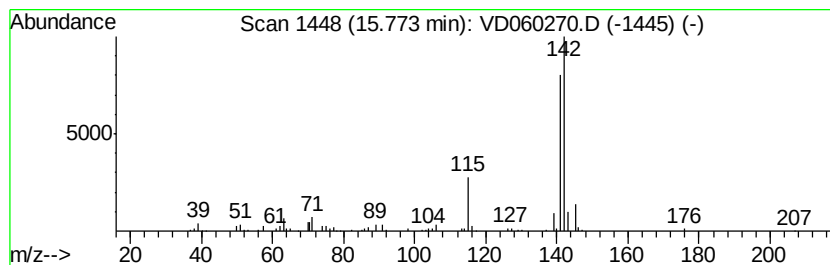
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D100118S.M
Quant Title : SW846 8260

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

Peak Number 9 Benzocycloheptatriene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.77	12.26 ug/l	348282	1,4-Dichlorobenzene-d4	13.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzocycloheptatriene	142	C11H10	000264-09-5	90
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	90
3		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	90
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	80
5		1-Naphthaleneethanol	172	C12H12O	000773-99-9	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD110118\
Data File : VD060270.D
Acq On : 1 Nov 2018 18:57
Operator : VA/AP
Sample : J5200-19
Misc : 6.83g/5ml/MSVOA D/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
OR-03-103118-B

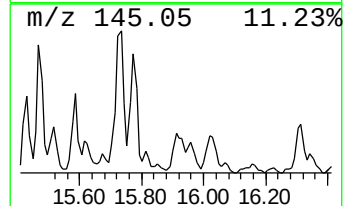
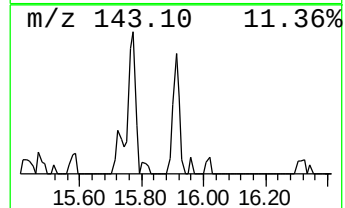
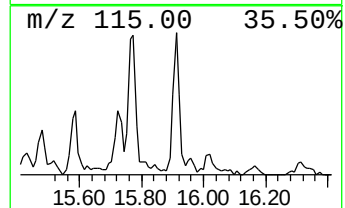
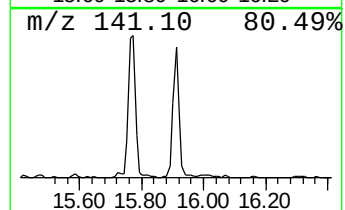
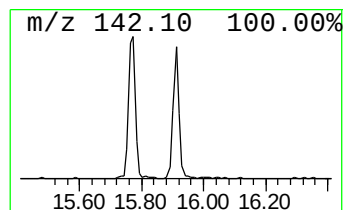
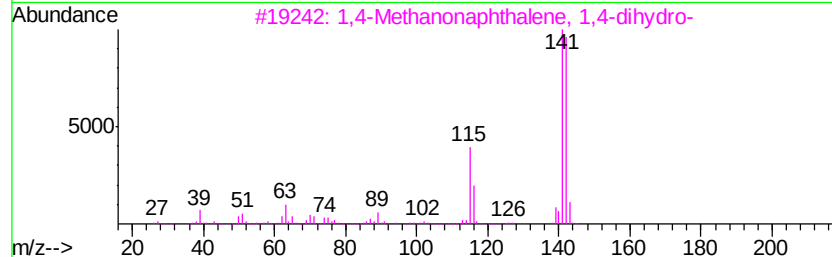
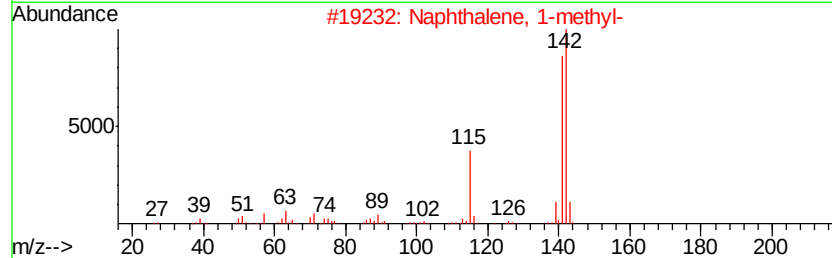
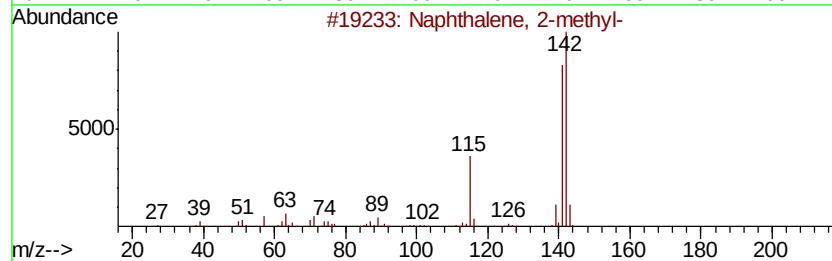
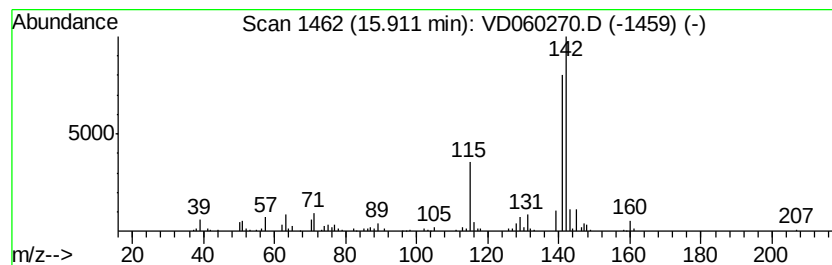
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.91	9.65 ug/l	274174	1,4-Dichlorobenzene-d4	13.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	95
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
3		1,4-Methanonaphthalene, 1,4-dihydro...	142	C11H10	004453-90-1	87
4		Benzocycloheptatriene	142	C11H10	000264-09-5	81
5		Spiro(9-methylenetricyclo[6.2.1]nona-2,5,8-trien-9-ylidene)-2,3-dihydro-1H-benzofuran	184	C13H12O	033929-47-4	78



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD110118\
Data File : VD060270.D
Acq On : 1 Nov 2018 18:57
Operator : VA/AP
Sample : J5200-19
Misc : 6.83g/5ml/MSVOA_D/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
OR-03-103118-B

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D100118S.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
1H-Indene, 2,3-di...	14.41	7.5	ug/l	213807	4	13.37	1419950	50.0
3-Phenylbut-1-ene	14.52	12.8	ug/l	363689	4	13.37	1419950	50.0
Naphthalene, 1,2,...	14.64	9.9	ug/l	281623	4	13.37	1419950	50.0
1H-Indene, 2,3-di...	14.75	7.3	ug/l	207285	4	13.37	1419950	50.0
1-Methylindan-2-one	15.07	8.3	ug/l	235139	4	13.37	1419950	50.0
Benzene, (1-methy...	15.35	14.4	ug/l	408080	4	13.37	1419950	50.0
Naphthalene, 1,2,...	15.59	9.0	ug/l	254652	4	13.37	1419950	50.0
Naphthalene, 1,2,...	15.72	9.7	ug/l	276351	4	13.37	1419950	50.0
Benzocycloheptatr...	15.77	12.3	ug/l	348282	4	13.37	1419950	50.0
Naphthalene, 1-me...	15.91	9.7	ug/l	274174	4	13.37	1419950	50.0