

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD040919\
 Data File : VD061689.D
 Acq On : 9 Apr 2019 21:52
 Operator : VA/AP
 Sample : K2316-06
 Misc : 5.10a/5ml/MSVOA D/SOIL
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 20190174-AMND-COMP-CS-1

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_D\METHOD\82D040619S.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	1.982	53	57	64	rVB2	7678	19355	1.08%	0.070%
2	2.120	69	71	81	rVB3	8424	28608	1.60%	0.104%
3	3.291	180	190	197	rBV2	239647	686348	38.28%	2.499%
4	3.410	198	202	211	rVB	83416	204734	11.42%	0.745%
5	3.843	241	246	255	rBV	28772	74787	4.17%	0.272%
6	6.097	468	475	481	rBV2	6626	21736	1.21%	0.079%
7	6.983	558	565	569	rBV3	8412	31736	1.77%	0.116%
8	7.082	569	575	591	rVB	442433	1268866	70.76%	4.620%
9	7.485	609	616	629	rBV	187370	522581	29.14%	1.903%
10	7.731	635	641	646	rVB	15432	43930	2.45%	0.160%
11	8.253	687	694	706	rBV	588368	1532228	85.45%	5.579%
12	8.893	751	759	764	rBV3	23970	76214	4.25%	0.278%
13	9.582	823	829	835	rBV3	16421	49438	2.76%	0.180%
14	9.759	842	847	851	rBV2	15700	45049	2.51%	0.164%
15	10.438	906	916	930	rBV	759041	1793082	100.00%	6.529%
16	10.655	934	938	943	rVV5	6773	22031	1.23%	0.080%
17	10.950	964	968	976	rVB3	26882	64359	3.59%	0.234%
18	11.128	981	986	992	rVV4	10301	24676	1.38%	0.090%
19	11.246	992	998	1005	rVB6	7007	32422	1.81%	0.118%
20	11.512	1020	1025	1029	rBV2	22355	42230	2.36%	0.154%
21	11.649	1034	1039	1041	rVV2	15598	32757	1.83%	0.119%
22	11.699	1041	1044	1047	rVV3	22070	51408	2.87%	0.187%
23	11.767	1047	1051	1056	rVB	34349	75708	4.22%	0.276%
24	11.856	1056	1060	1064	rBV	50248	114322	6.38%	0.416%
25	11.915	1064	1066	1071	rVV3	17615	38044	2.12%	0.139%
26	12.023	1071	1077	1080	rVV2	9886	24737	1.38%	0.090%
27	12.082	1080	1083	1085	rVV	18320	30739	1.71%	0.112%
28	12.142	1085	1089	1092	rVV	49701	92727	5.17%	0.338%
29	12.319	1104	1107	1110	rBV3	9198	21839	1.22%	0.080%
30	12.388	1110	1114	1118	rVV	787843	1307879	72.94%	4.762%
31	12.476	1121	1123	1127	rVB3	14201	29491	1.64%	0.107%
32	12.565	1127	1132	1136	rVB2	27693	61094	3.41%	0.222%
33	12.673	1136	1143	1145	rBV3	42899	123508	6.89%	0.450%
34	12.713	1145	1147	1150	rVB	61980	97827	5.46%	0.356%

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35	12.772	1150	1153	1158	rVB	49612	83673	4.67%	0.305%
36	12.870	1158	1163	1167	rBV3	24966	50624	2.82%	0.184%
37	12.939	1167	1170	1174	rVB4	9549	18765	1.05%	0.068%
38	13.018	1174	1178	1181	rBV2	87876	216350	12.07%	0.788%
39	13.067	1181	1183	1185	rVV2	64068	95397	5.32%	0.347%
40	13.116	1185	1188	1191	rVV	37170	63995	3.57%	0.233%
41	13.234	1194	1200	1203	rVV2	136322	280384	15.64%	1.021%
42	13.323	1203	1209	1216	rVV4	122754	405812	22.63%	1.478%
43	13.411	1216	1218	1222	rVB	57066	95041	5.30%	0.346%
44	13.490	1222	1226	1231	rBV3	72114	206942	11.54%	0.754%
45	13.569	1231	1234	1239	rVV2	676592	1027702	57.31%	3.742%
46	13.628	1239	1240	1242	rVV2	39258	38959	2.17%	0.142%
47	13.687	1242	1246	1247	rVV4	27687	62203	3.47%	0.226%
48	13.726	1247	1250	1254	rVV2	180008	314316	17.53%	1.145%
49	13.845	1257	1262	1265	rVV5	63700	184606	10.30%	0.672%
50	13.904	1265	1268	1270	rVV2	117713	240612	13.42%	0.876%
51	13.943	1270	1272	1274	rVV2	159234	258776	14.43%	0.942%
52	13.982	1274	1276	1279	rVV2	134607	232556	12.97%	0.847%
53	14.041	1279	1282	1283	rVV2	58190	102210	5.70%	0.372%
54	14.071	1283	1285	1286	rVV	79950	108663	6.06%	0.396%
55	14.101	1286	1288	1290	rVV2	91427	179128	9.99%	0.652%
56	14.140	1290	1292	1295	rVV2	131388	260766	14.54%	0.950%
57	14.209	1297	1299	1300	rVV2	66098	111481	6.22%	0.406%
58	14.248	1300	1303	1310	rVV2	202888	499146	27.84%	1.818%
59	14.337	1310	1312	1314	rVV	89456	171981	9.59%	0.626%
60	14.376	1314	1316	1318	rVV3	151227	272069	15.17%	0.991%
61	14.455	1318	1324	1326	rVV4	200939	596955	33.29%	2.174%
62	14.534	1329	1332	1334	rVV	661263	971955	54.21%	3.539%
63	14.573	1334	1336	1338	rVV	143684	233837	13.04%	0.851%
64	14.612	1338	1340	1341	rVV	108837	153235	8.55%	0.558%
65	14.642	1341	1343	1347	rVV3	124541	303857	16.95%	1.106%
66	14.721	1349	1351	1352	rVV2	74519	106676	5.95%	0.388%
67	14.740	1352	1353	1355	rVV	74425	103050	5.75%	0.375%
68	14.790	1355	1358	1359	rVV2	207583	282839	15.77%	1.030%
69	14.819	1359	1361	1364	rVV	291163	495493	27.63%	1.804%
70	14.927	1368	1372	1374	rVV2	152363	390348	21.77%	1.421%
71	14.987	1377	1378	1379	rVV	124431	135661	7.57%	0.494%

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72	15.016	1379	1381	1384	rVV3	188126	408025	22.76%	1.486%
73	15.065	1384	1386	1388	rVV	245992	328321	18.31%	1.195%
74	15.105	1388	1390	1392	rVV2	56057	102101	5.69%	0.372%
75	15.154	1392	1395	1399	rVV2	275297	627929	35.02%	2.286%
76	15.233	1399	1403	1406	rVV4	181365	455525	25.40%	1.659%
77	15.292	1406	1409	1411	rVV3	252167	553591	30.87%	2.016%
78	15.331	1411	1413	1414	rVV2	177673	257088	14.34%	0.936%
79	15.361	1414	1416	1417	rVV	277175	356706	19.89%	1.299%
80	15.390	1417	1419	1422	rVV2	344131	473075	26.38%	1.723%
81	15.439	1422	1424	1426	rVV2	350727	512724	28.59%	1.867%
82	15.479	1426	1428	1432	rVV3	203519	466170	26.00%	1.697%
83	15.548	1432	1435	1438	rVV3	150596	358106	19.97%	1.304%
84	15.597	1438	1440	1443	rVV3	103664	216939	12.10%	0.790%
85	15.646	1443	1445	1446	rVV	170973	233304	13.01%	0.850%
86	15.685	1446	1449	1451	rVV	459984	778576	43.42%	2.835%
87	15.745	1454	1455	1458	rVV3	169047	304422	16.98%	1.108%
88	15.813	1458	1462	1465	rVV4	182834	563306	31.42%	2.051%
89	15.863	1465	1467	1468	rVV	101514	145655	8.12%	0.530%
90	15.922	1468	1473	1474	rVV3	216037	549169	30.63%	2.000%
91	16.010	1477	1482	1484	rVV2	245833	632790	35.29%	2.304%
92	16.079	1488	1489	1491	rVV2	120834	125995	7.03%	0.459%
93	16.158	1491	1497	1499	rVV5	118857	351070	19.58%	1.278%
94	16.237	1502	1505	1507	rVV3	128595	228045	12.72%	0.830%
95	16.276	1507	1509	1511	rVB	52859	53637	2.99%	0.195%
96	16.375	1516	1519	1520	rVB	35959	42560	2.37%	0.155%
97	16.493	1527	1531	1533	rBV4	63159	140571	7.84%	0.512%
98	16.591	1539	1541	1544	rVB3	24226	31893	1.78%	0.116%
99	16.640	1544	1546	1548	rVV2	26208	34430	1.92%	0.125%
100	16.729	1552	1555	1559	rVB2	54227	88887	4.96%	0.324%

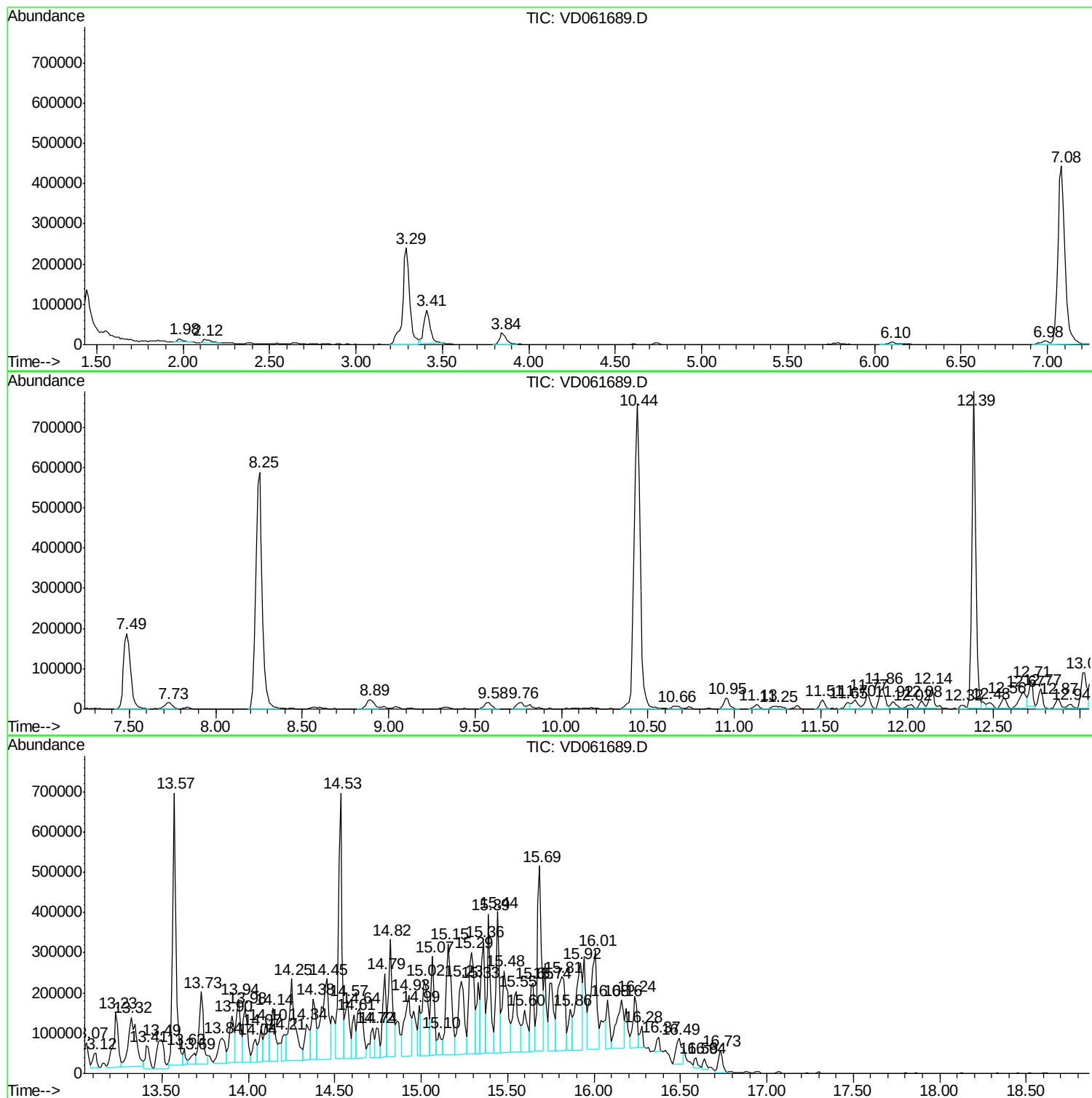
Sum of corrected areas: 27463163

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Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D040619S.M
Quant Title : SW846 8260

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



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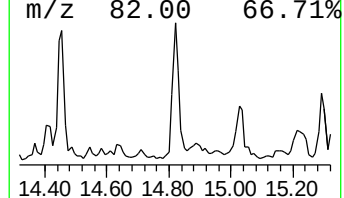
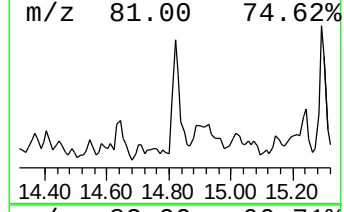
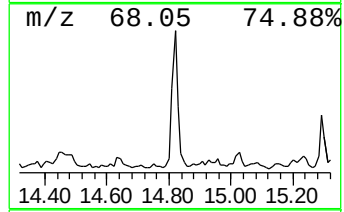
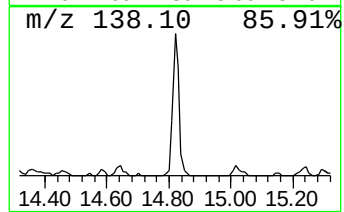
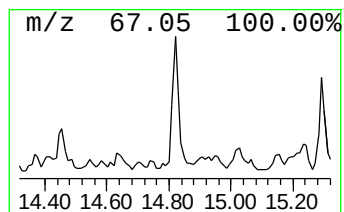
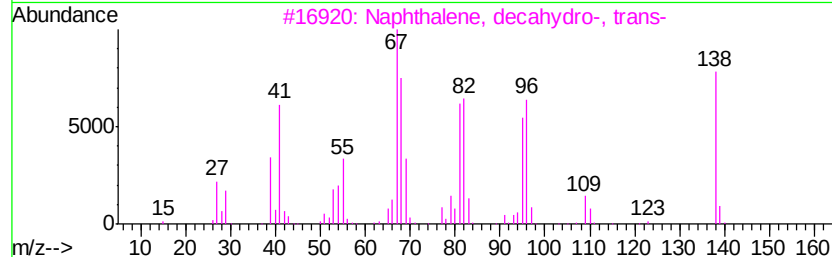
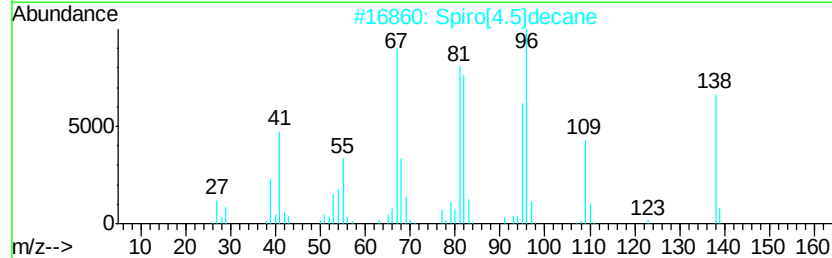
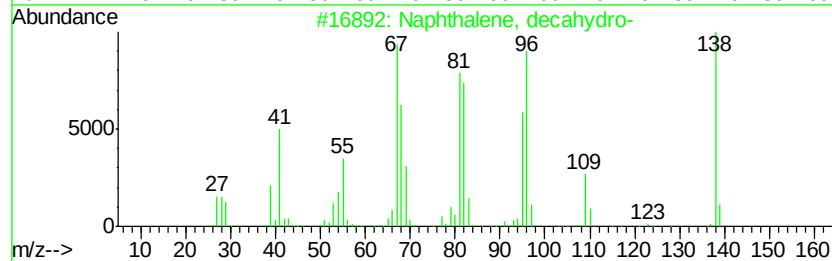
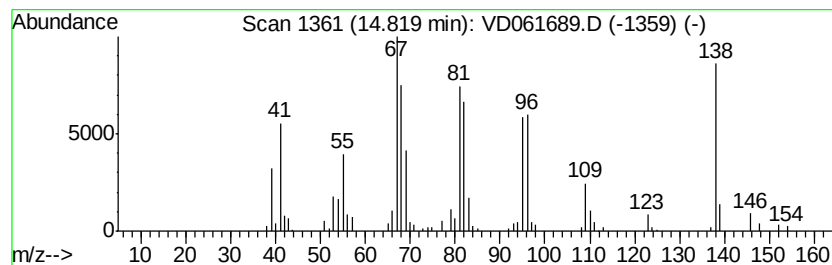
Instrument :
MSVOA_D
ClientSampleId :
20190174-AMND-COMP-CS-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D040619S.M
Quant Title : SW846 8260

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

Peak Number 1 Naphthalene, decahydro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	25.49 ug/l	495493	1,4-Dichlorobenzene-d4	14.53
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, decahydro-	138 C10H18	000091-17-8 93
2		Spiro[4.5]decane	138 C10H18	000176-63-6 90
3		Naphthalene, decahydro-, trans-	138 C10H18	000493-02-7 90
4		Naphthalene, decahydro-, cis-	138 C10H18	000493-01-6 89
5		1,1'-Bicyclopentyl	138 C10H18	001636-39-1 83



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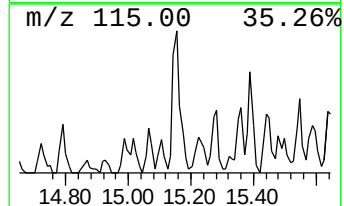
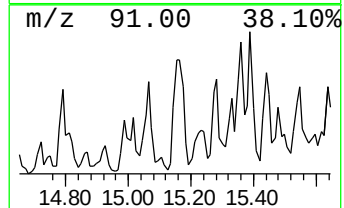
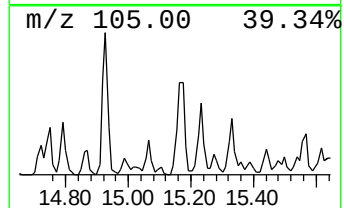
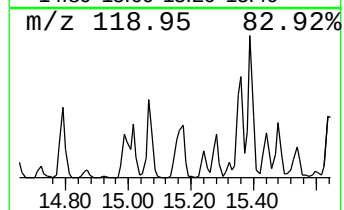
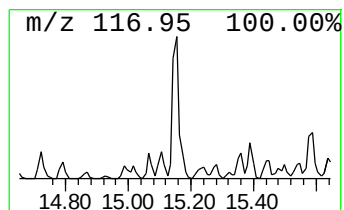
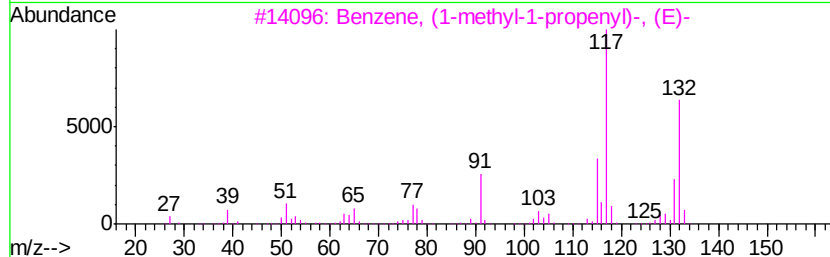
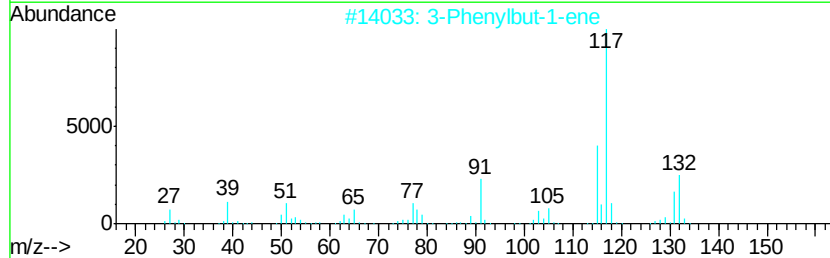
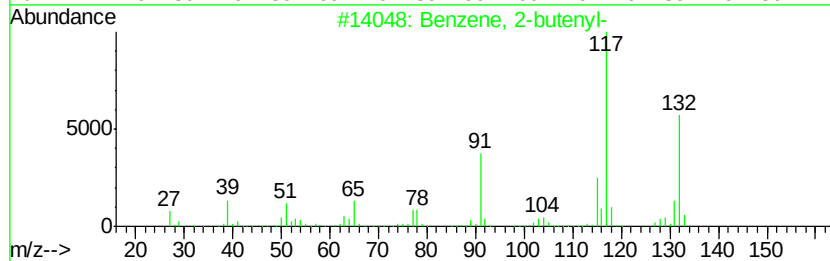
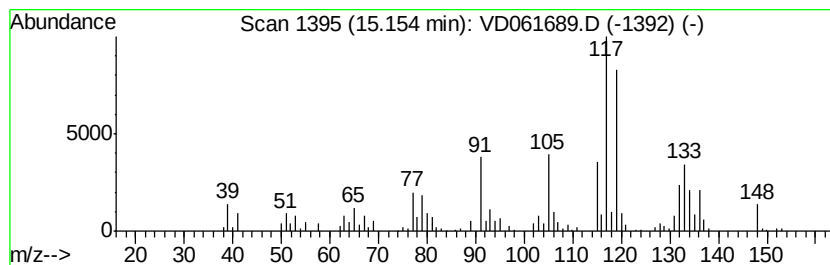
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown15.15 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.15	32.30 ug/l	627929	1,4-Dichlorobenzene-d4	14.53
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 2-butenyl-	132 C10H12	001560-06-1 46
2		3-Phenylbut-1-ene	132 C10H12	000934-10-1 46
3		Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000768-00-3 46
4		1-Phenyl-1-butene	132 C10H12	000824-90-8 46
5		Benzene, 1-methyl-4-(2-propenyl)-	132 C10H12	003333-13-9 46



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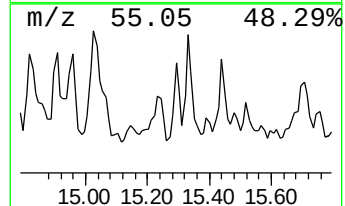
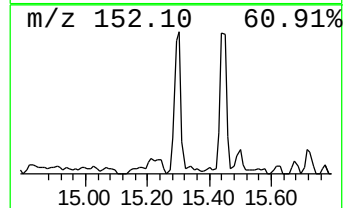
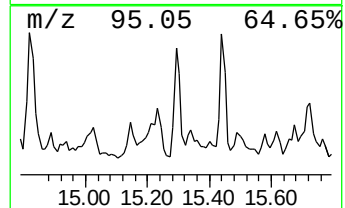
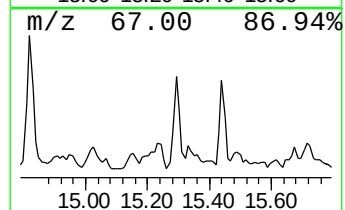
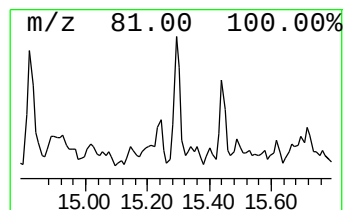
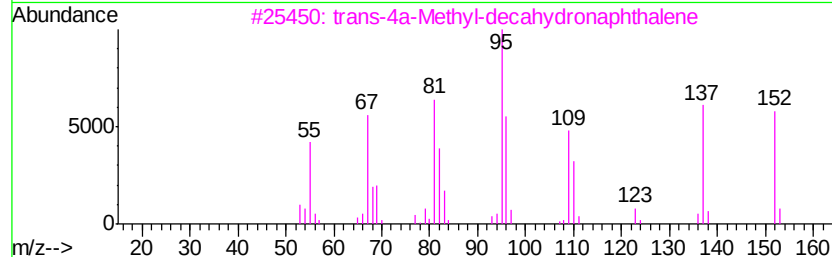
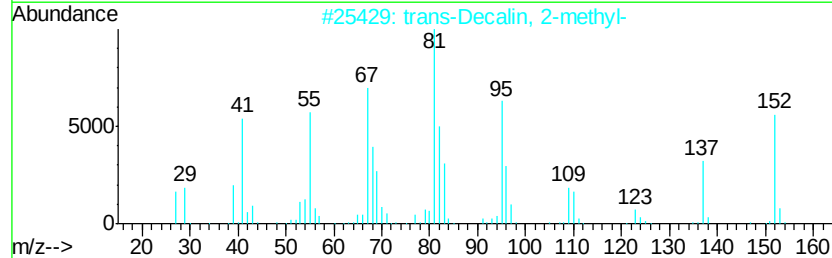
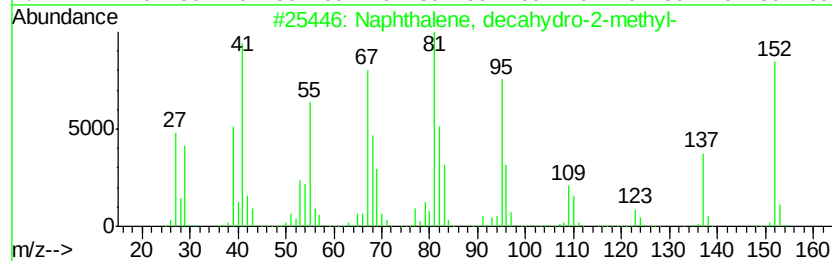
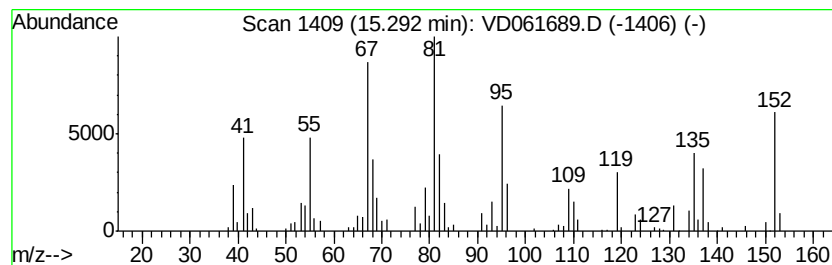
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Naphthalene, decahydro-2-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.29	28.48 ug/l	553591	1,4-Dichlorobenzene-d4	14.53
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, decahydro-2-methyl-	152 C11H20	002958-76-1 89
2		trans-Decalin, 2-methyl-	152 C11H20	1000152-47-3 89
3		trans-4a-Methyl-decahydronaphtha...	152 C11H20	002547-27-5 64
4		Cyclohexanone, 5-methyl-2-(1-met...	152 C10H16O	015932-80-6 64
5		cis-Decalin, 2-syn-methyl-	152 C11H20	1000155-85-6 62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD040919\
Data File : VD061689.D
Acq On : 9 Apr 2019 21:52
Operator : VA/AP
Sample : K2316-06
Misc : 5.10a/5ml/MSVOA D/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
20190174-AMND-COMP-CS-1

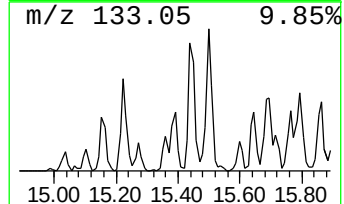
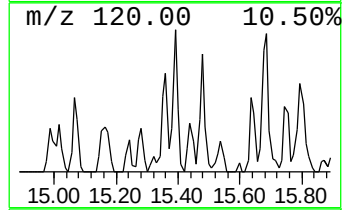
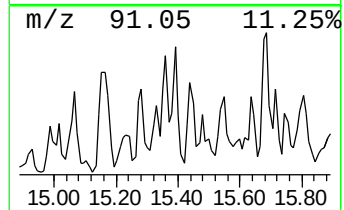
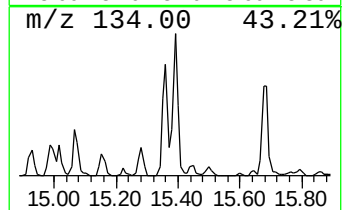
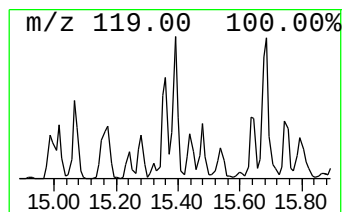
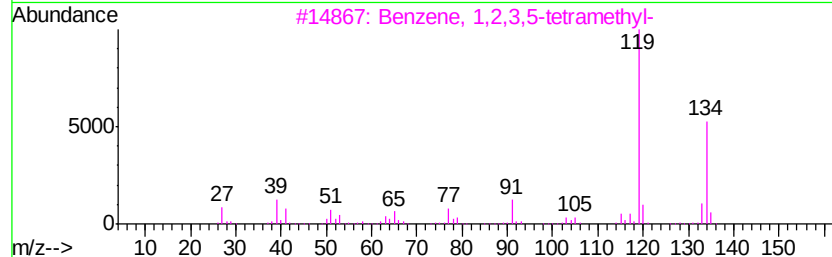
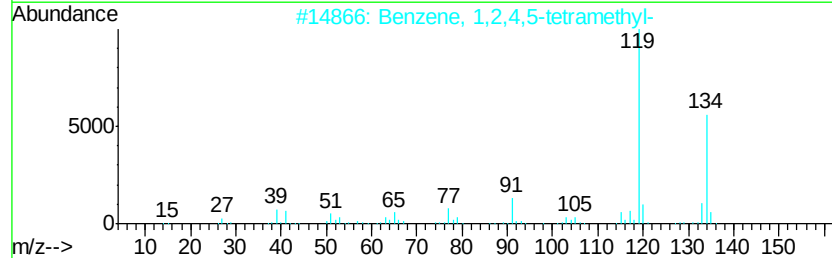
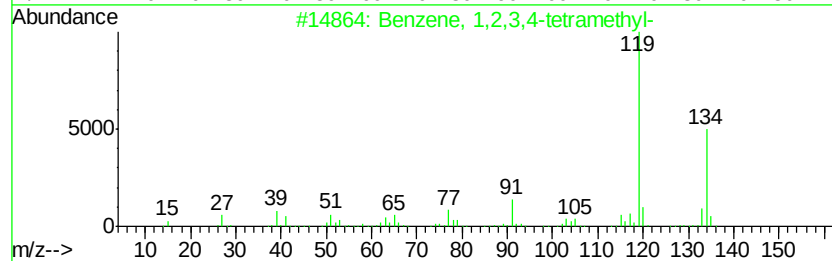
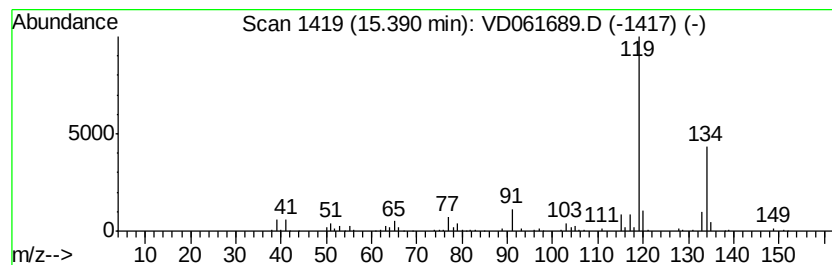
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.39	24.34 ug/l	473075	1,4-Dichlorobenzene-d4	14.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4		o-Cymene	134	C10H14	000527-84-4	94
5		p-Cymene	134	C10H14	000099-87-6	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD040919\
Data File : VD061689.D
Acq On : 9 Apr 2019 21:52
Operator : VA/AP
Sample : K2316-06
Misc : 5.10a/5ml/MSVOA D/SOIL
ALS Vial : 21 Sample Multiplier: 1

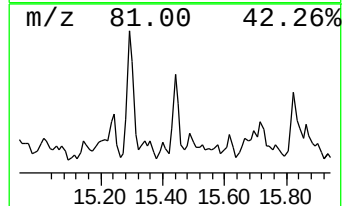
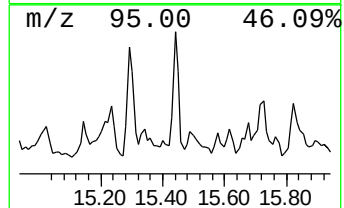
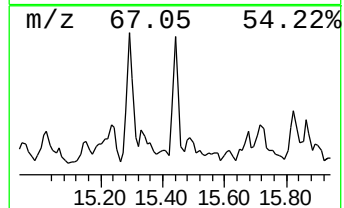
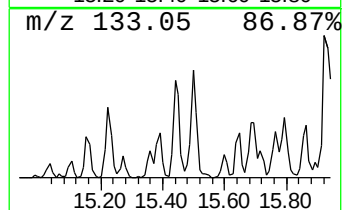
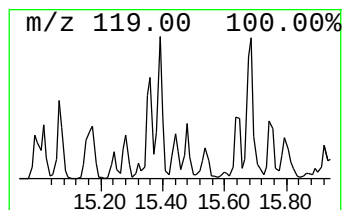
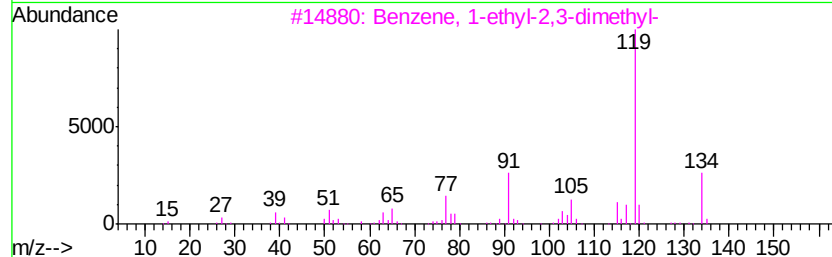
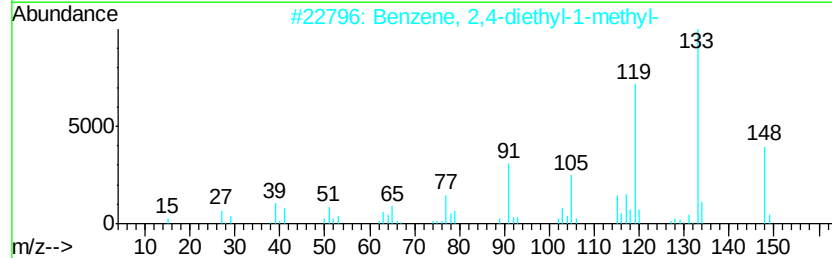
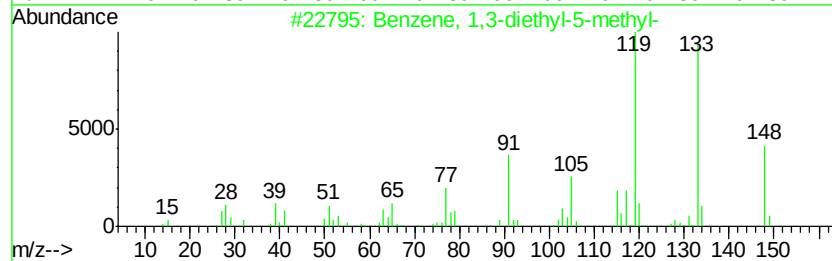
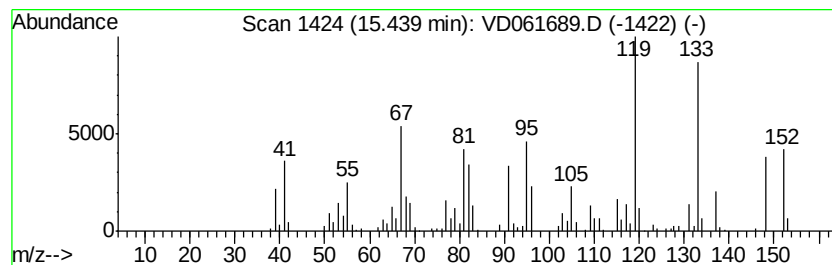
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ClientSampleId :
20190174-AMND-COMP-CS-1

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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,3-diethyl-5-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.44	26.38 ug/l	512724	1,4-Dichlorobenzene-d4	14.53	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	89
2	Benzene, 2,4-diethyl-1-methyl-	148	C11H16	001758-85-6	46
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	43
4	Ethanone, 1-(2,4-dimethylphenyl)-	148	C10H12O	000089-74-7	35
5	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	30



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD040919\
Data File : VD061689.D
Acq On : 9 Apr 2019 21:52
Operator : VA/AP
Sample : K2316-06
Misc : 5.10a/5ml/MSVOA D/SOIL
ALS Vial : 21 Sample Multiplier: 1

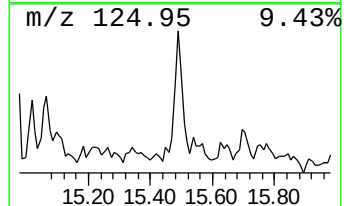
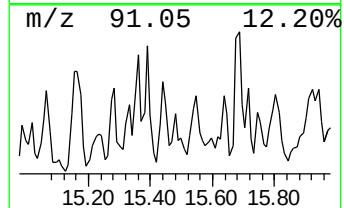
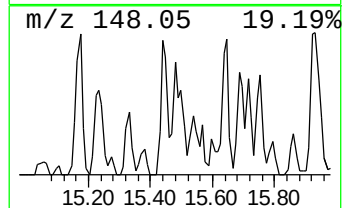
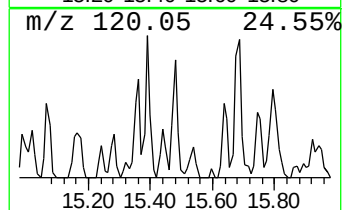
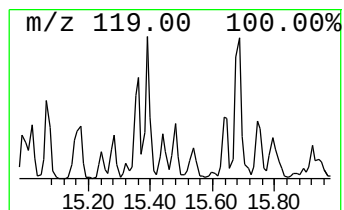
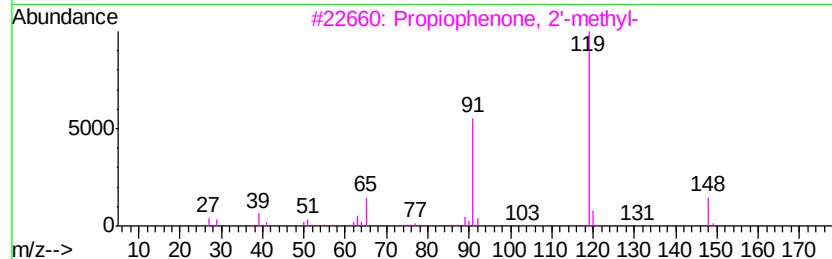
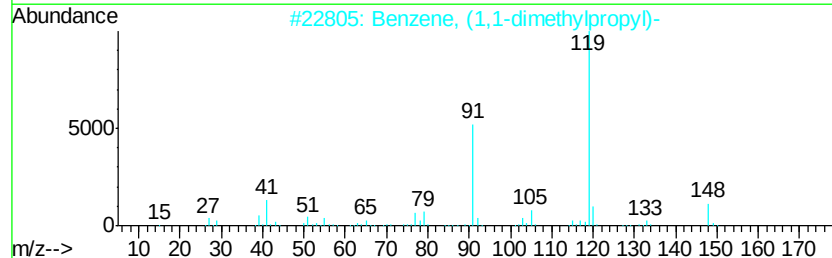
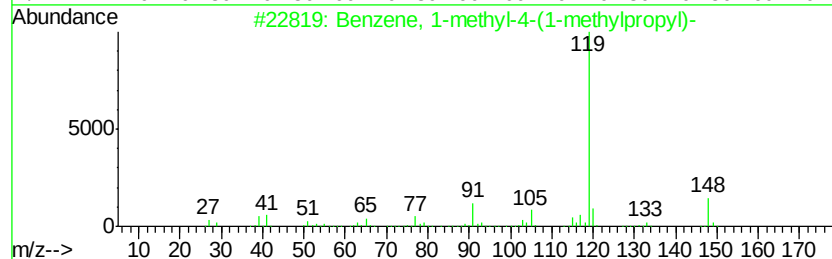
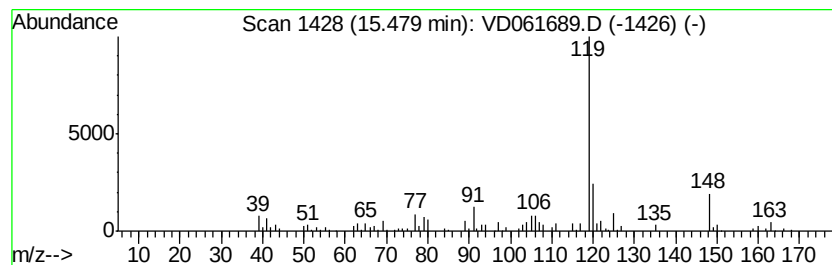
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D040619S.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-methyl-4-(1-meth... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.48	23.98 ug/l	466170	1,4-Dichlorobenzene-d4	14.53
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-methyl-4-(1-methylpro...	148 C11H16	001595-16-0 50
2		Benzene, (1,1-dimethylpropyl)-	148 C11H16	002049-95-8 50
3		Propiophenone, 2'-methyl-	148 C10H12O	002040-14-4 50
4		Chrysanthemic acid 2,4-dimethy...	286 C19H26O2	000070-38-2 47
5		Benzene, 1,4-dimethyl-2-(2-methy...	162 C12H18	055669-88-0 47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD040919\
Data File : VD061689.D
Acq On : 9 Apr 2019 21:52
Operator : VA/AP
Sample : K2316-06
Misc : 5.10g/5ml/MSVOA D/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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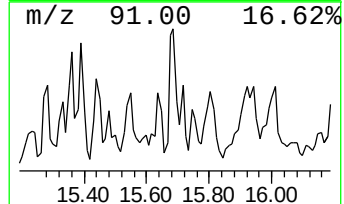
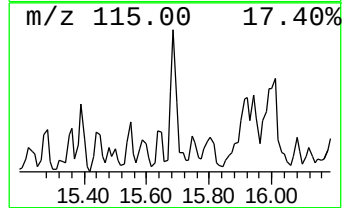
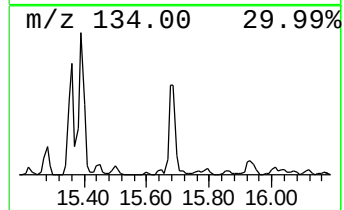
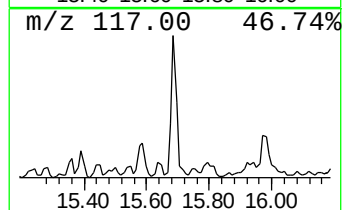
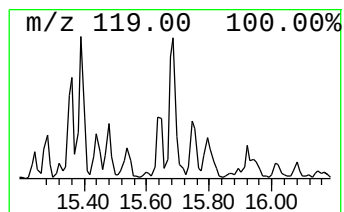
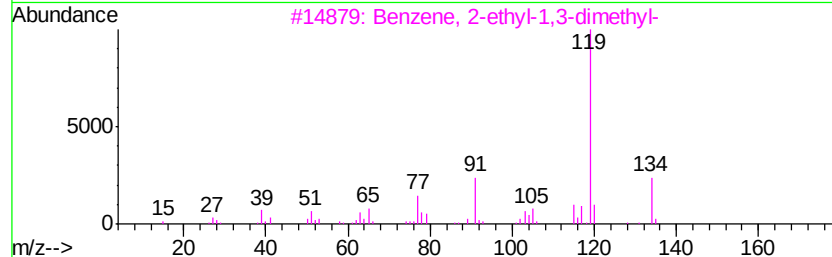
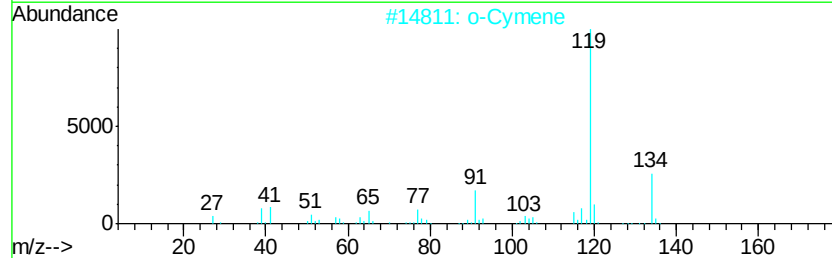
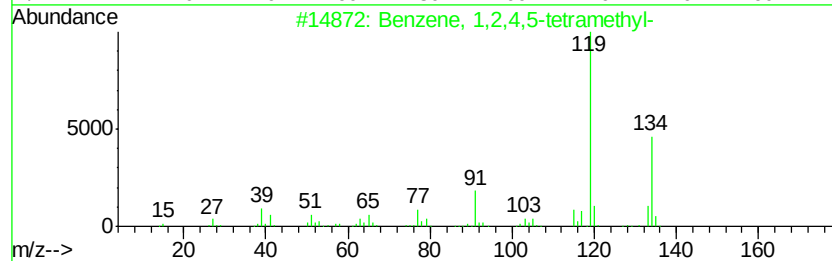
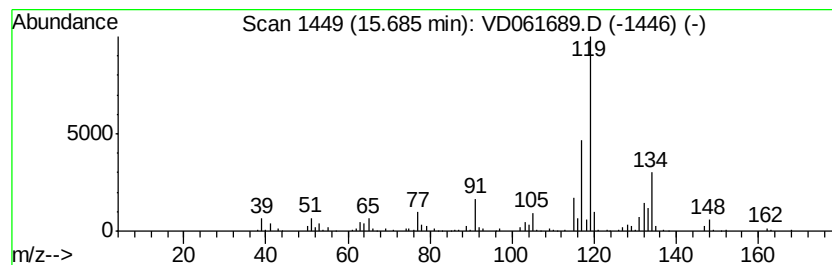
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.69	40.05 ug/l	778576	1,4-Dichlorobenzene-d4	14.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
2		o-Cymene	134	C10H14	000527-84-4	60
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	60
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	60
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD040919\
Data File : VD061689.D
Acq On : 9 Apr 2019 21:52
Operator : VA/AP
Sample : K2316-06
Misc : 5.10g/5ml/MSVOA D/SOIL
ALS Vial : 21 Sample Multiplier: 1

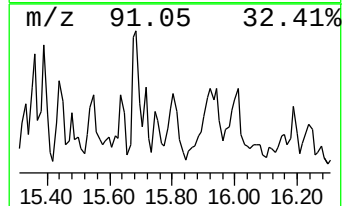
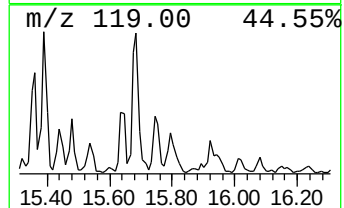
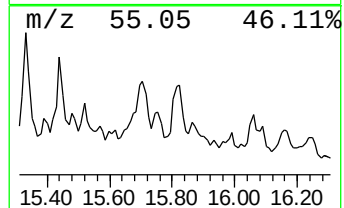
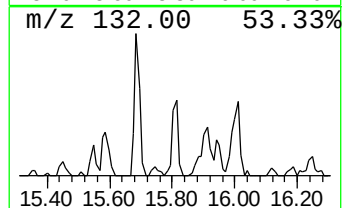
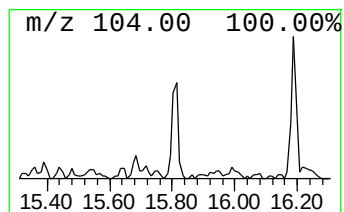
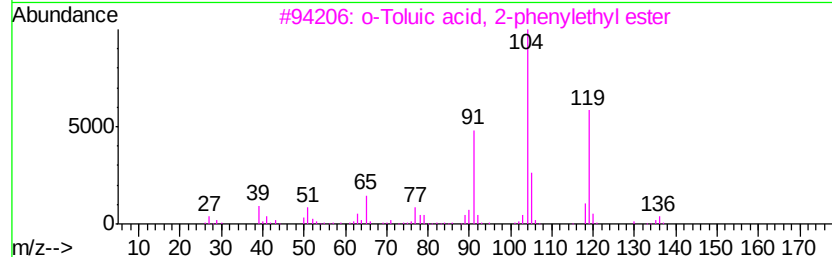
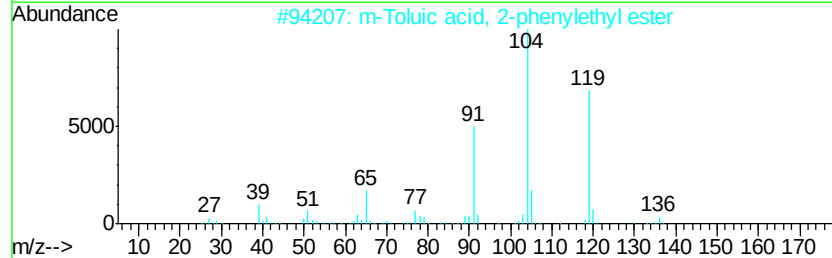
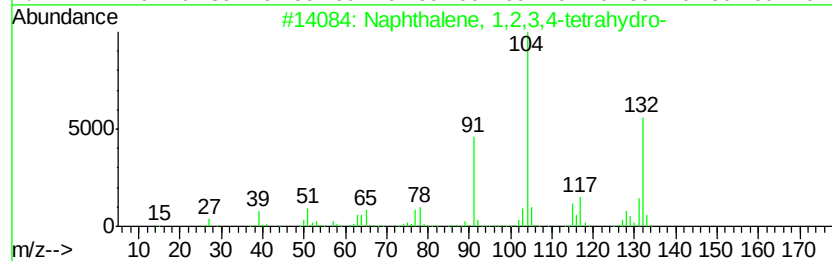
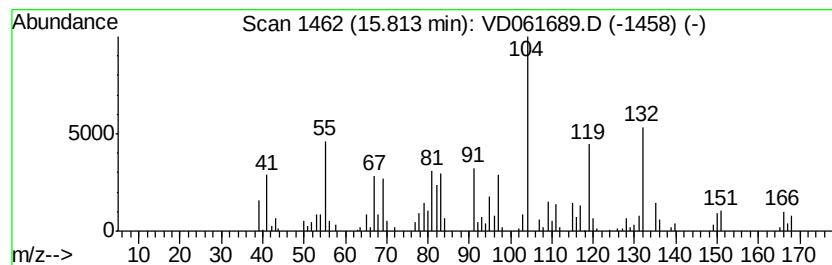
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D040619S.M
Quant Title : SW846 8260

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

Peak Number 8 unknown15.81 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.81	28.98 ug/l	563306	1,4-Dichlorobenzene-d4	14.53
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-	132 C10H12	000119-64-2 49
2		m-Toluic acid, 2-phenylethyl ester	240 C16H16O2	006641-74-3 38
3		o-Toluic acid, 2-phenylethyl ester	240 C16H16O2	1000293-42-6 38
4		p-Toluic acid, 2-phenylethyl ester	240 C16H16O2	203587-50-2 38
5		5-Chloropentanoic acid, 2-phenyl...	240 C13H17ClO2	1000293-49-1 22



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD040919\
Data File : VD061689.D
Acq On : 9 Apr 2019 21:52
Operator : VA/AP
Sample : K2316-06
Misc : 5.10g/5ml/MSVOA D/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
20190174-AMND-COMP-CS-1

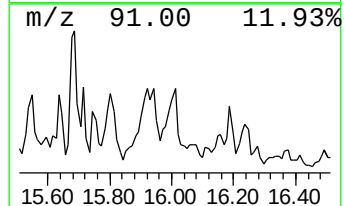
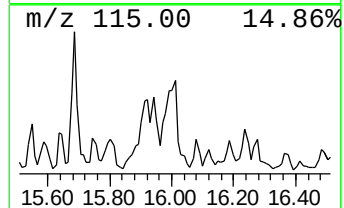
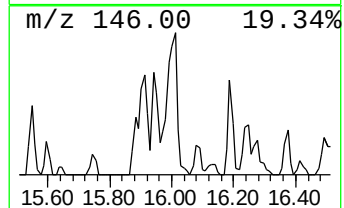
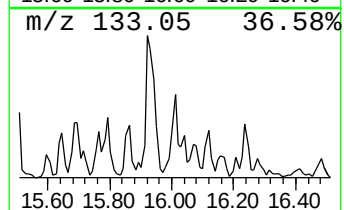
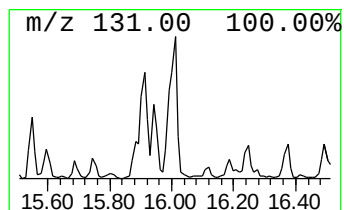
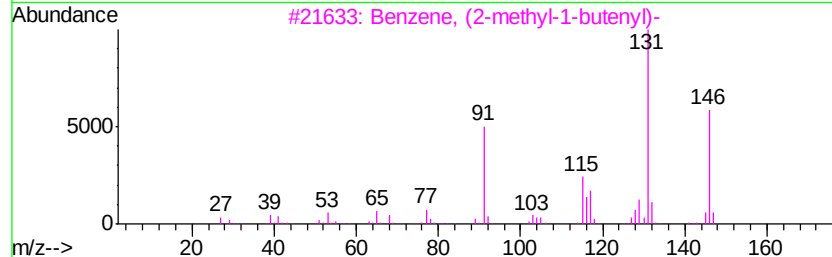
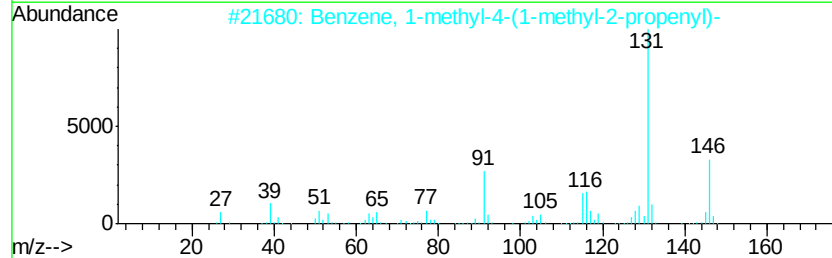
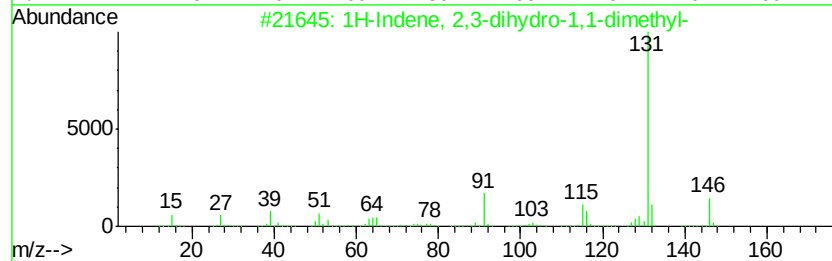
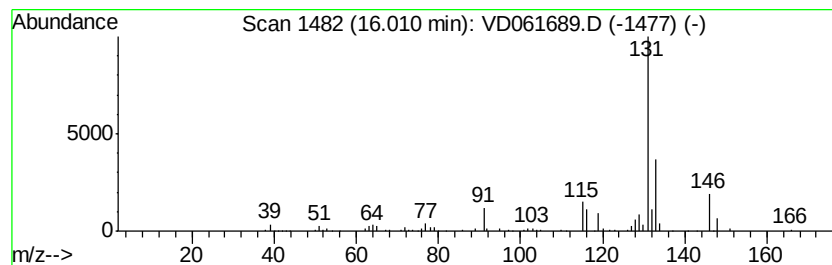
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.01	32.55 ug/l	632790	1,4-Dichlorobenzene-d4	14.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	81
2		Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	81
3		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	74
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	74
5		Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	74



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD040919\
Data File : VD061689.D
Acq On : 9 Apr 2019 21:52
Operator : VA/AP
Sample : K2316-06
Misc : 5.10g/5ml/MSVOA_D/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
20190174-AMND-COMP-CS-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D040619S.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
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unknown15.15	15.15	32.3	ug/l	627929	4	14.53	971955	50.0
Naphthalene, deca...	15.29	28.5	ug/l	553591	4	14.53	971955	50.0
Benzene, 1,2,3,4-...	15.39	24.3	ug/l	473075	4	14.53	971955	50.0
Benzene, 1,3-diet...	15.44	26.4	ug/l	512724	4	14.53	971955	50.0
Benzene, 1-methyl...	15.48	24.0	ug/l	466170	4	14.53	971955	50.0
Benzene, 1,2,4,5-...	15.69	40.0	ug/l	778576	4	14.53	971955	50.0
unknown15.81	15.81	29.0	ug/l	563306	4	14.53	971955	50.0
1H-Indene, 2,3-di...	16.01	32.5	ug/l	632790	4	14.53	971955	50.0