

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD062318\
 Data File : VD058344.D
 Acq On : 23 Jun 2018 16:46
 Operator : VA/AP
 Sample : J3664-19
 Misc : 5.33g/5ml/MSVOA D/SOIL
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 DUP-0485-180619

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\82D061918S.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	1.228	94	102	104	rBV	2073268	4603596	9.45%	0.790%
2	1.297	104	109	130	rVB	3997508	14945068	30.68%	2.565%
3	5.570	536	543	547	rBV	333301	983254	2.02%	0.169%
4	5.658	547	552	568	rVB	637725	1983389	4.07%	0.340%
5	6.032	584	590	599	rBV	306514	804744	1.65%	0.138%
6	6.702	652	658	674	rBV	855471	2245278	4.61%	0.385%
7	7.293	707	718	727	rBV	571633	1893391	3.89%	0.325%
8	8.720	855	863	867	rBV	1084323	2704253	5.55%	0.464%
9	8.946	879	886	890	rBV4	115311	487584	1.00%	0.084%
10	9.222	907	914	921	rVB2	325573	912620	1.87%	0.157%
11	9.754	963	968	972	rBV	235447	519982	1.07%	0.089%
12	9.970	986	990	993	rVV3	185752	530554	1.09%	0.091%
13	10.039	993	997	1002	rVV	703530	1639337	3.37%	0.281%
14	10.128	1002	1006	1011	rVV	733496	1787165	3.67%	0.307%
15	10.443	1034	1038	1042	rVB	560896	1089850	2.24%	0.187%
16	10.649	1054	1059	1061	rBV	263838	601969	1.24%	0.103%
17	10.728	1061	1067	1071	rVV2	3543591	8436680	17.32%	1.448%
18	10.797	1071	1074	1079	rVB2	479796	1111301	2.28%	0.191%
19	10.896	1079	1084	1088	rBV2	479366	1066936	2.19%	0.183%
20	10.994	1088	1094	1097	rBV2	865065	2575037	5.29%	0.442%
21	11.053	1097	1100	1103	rVB	1645644	2912368	5.98%	0.500%
22	11.112	1103	1106	1113	rVB	1106654	2208701	4.53%	0.379%
23	11.211	1113	1116	1121	rBV	904603	1726913	3.55%	0.296%
24	11.368	1128	1132	1140	rBV4	1906313	7423332	15.24%	1.274%
25	11.467	1140	1142	1146	rVB	700267	1212396	2.49%	0.208%
26	11.604	1149	1156	1159	rBV3	3625141	8905882	18.28%	1.528%
27	11.683	1159	1164	1172	rBV5	3495825	10707792	21.98%	1.837%
28	11.782	1172	1174	1177	rVB2	940077	1566382	3.22%	0.269%
29	11.850	1177	1181	1188	rBV3	1772046	6298086	12.93%	1.081%
30	11.939	1188	1190	1192	rVB2	869649	997235	2.05%	0.171%
31	11.988	1192	1195	1197	rBV2	1794038	2589523	5.32%	0.444%
32	12.097	1203	1206	1211	rVB2	4763399	8267337	16.97%	1.419%
33	12.215	1214	1218	1222	rBV2	2484681	7091525	14.56%	1.217%
34	12.274	1222	1224	1226	rBV	2515374	4249119	8.72%	0.729%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	12.323	1226	1229	1231	rVV	2071560	3852440	7.91%	0.661%
36	12.362	1231	1233	1236	rVB2	2601883	3929395	8.07%	0.674%
37	12.412	1236	1238	1241	rVB3	1203641	2133473	4.38%	0.366%
38	12.461	1241	1243	1244	rBV	927939	1156765	2.37%	0.199%
39	12.510	1244	1248	1252	rBV3	2860945	5755498	11.82%	0.988%
40	12.579	1252	1255	1257	rBV2	993064	1483183	3.04%	0.255%
41	12.668	1262	1264	1269	rVV4	3227428	8518634	17.49%	1.462%
42	12.746	1269	1272	1279	rVV4	7099249	26926818	55.28%	4.621%
43	12.845	1279	1282	1284	rVV	5947760	11923908	24.48%	2.046%
44	12.933	1287	1291	1293	rVV2	3043354	7574003	15.55%	1.300%
45	13.032	1296	1301	1305	rVV2	5511102	13788260	28.31%	2.366%
46	13.111	1305	1309	1314	rVV	10511217	21216269	43.56%	3.641%
47	13.209	1314	1319	1322	rVV2	12050666	27459839	56.37%	4.712%
48	13.258	1322	1324	1326	rVV	3407950	6100785	12.52%	1.047%
49	13.307	1326	1329	1331	rVV	2966816	7313703	15.01%	1.255%
50	13.347	1331	1333	1336	rVV2	3492813	6905316	14.18%	1.185%
51	13.475	1336	1346	1351	rVV2	12981517	48710132	100.00%	8.359%
52	13.544	1351	1353	1358	rVV4	13197666	29177207	59.90%	5.007%
53	13.613	1358	1360	1361	rVV	3434351	5838867	11.99%	1.002%
54	13.672	1364	1366	1370	rVV3	8710164	19001367	39.01%	3.261%
55	13.760	1370	1375	1380	rVV4	12119453	31079120	63.80%	5.333%
56	13.839	1380	1383	1386	rVV2	9096724	14441398	29.65%	2.478%
57	13.908	1386	1390	1391	rVV	6241350	10582290	21.73%	1.816%
58	13.937	1391	1393	1395	rVV2	5518288	9110671	18.70%	1.563%
59	13.977	1395	1397	1401	rVV2	5897622	10986926	22.56%	1.885%
60	14.046	1401	1404	1405	rVV3	3920882	6100838	12.52%	1.047%
61	14.085	1405	1408	1412	rVV2	13307773	29832995	61.25%	5.119%
62	14.154	1412	1415	1417	rVV	5898902	9915227	20.36%	1.701%
63	14.193	1417	1419	1420	rVV2	2783038	4532467	9.30%	0.778%
64	14.302	1424	1430	1431	rVV4	4394289	12134759	24.91%	2.082%
65	14.341	1431	1434	1436	rVV2	6916411	14357218	29.47%	2.464%
66	14.400	1436	1440	1443	rVV3	6722568	17325310	35.57%	2.973%
67	14.449	1443	1445	1446	rVV2	1210640	1984269	4.07%	0.341%
68	14.479	1446	1448	1450	rVV	4507456	6476570	13.30%	1.111%
69	14.508	1450	1451	1454	rVV	2703542	3738298	7.67%	0.641%
70	14.567	1454	1457	1460	rVV3	2906490	5648169	11.60%	0.969%
71	14.636	1460	1464	1466	rVV3	3087951	5683062	11.67%	0.975%

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Stop Thrs : 0 Peak Location: TOP

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72	14.666	1466	1467	1469	rVV2	1639247	1731688	3.56%	0.297%
73	14.764	1473	1477	1479	rBV3	757639	1312640	2.69%	0.225%
74	14.804	1479	1481	1485	rVB5	1333877	1761623	3.62%	0.302%
75	14.873	1485	1488	1490	rBV	1397598	2310679	4.74%	0.397%
76	14.971	1495	1498	1500	rVV2	714981	1109571	2.28%	0.190%
77	15.109	1509	1512	1514	rVV	660025	1064493	2.19%	0.183%
78	15.286	1526	1530	1536	rVB5	562035	1284997	2.64%	0.221%
79	15.385	1537	1540	1545	rVB	1590570	2396681	4.92%	0.411%

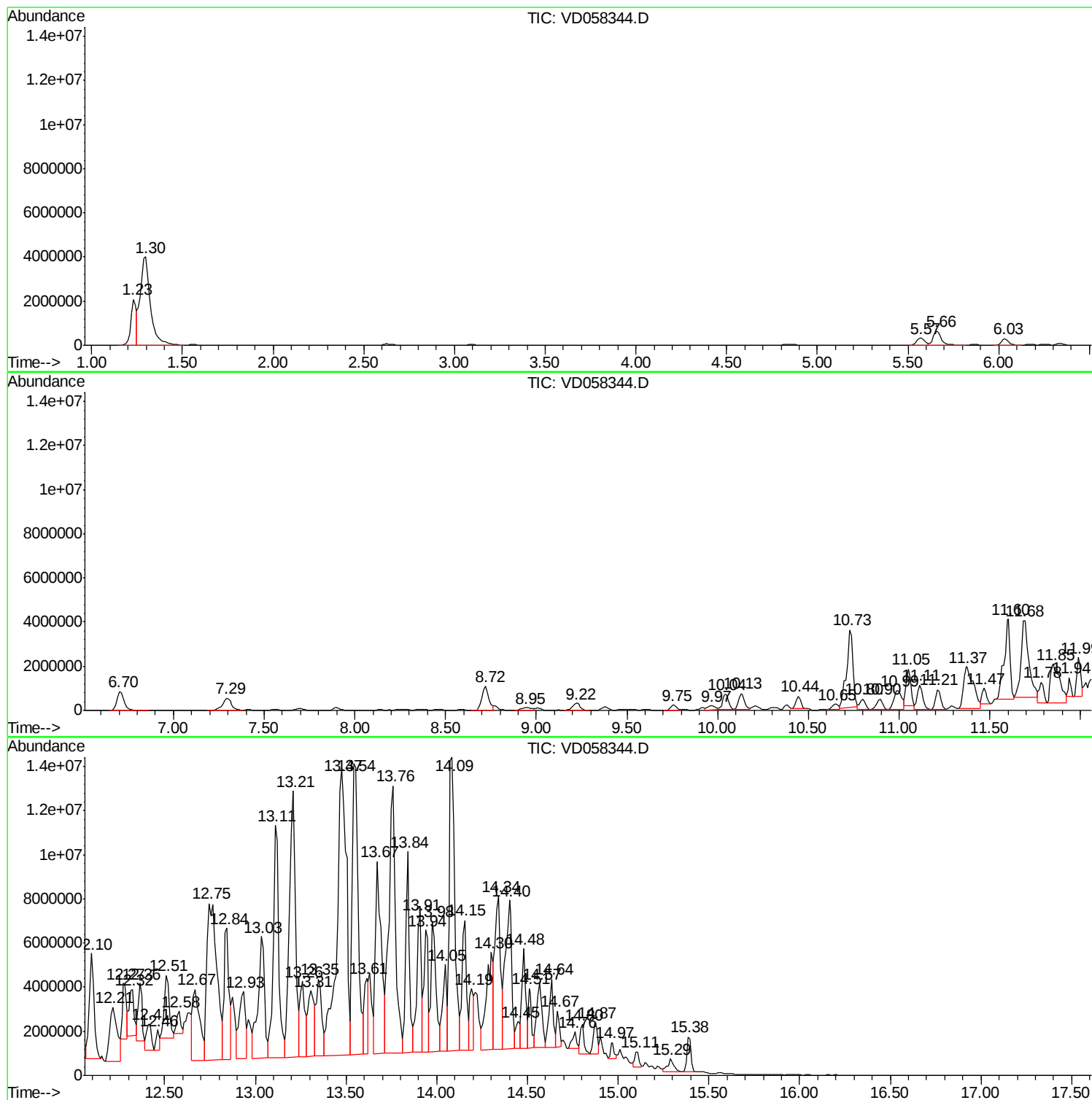
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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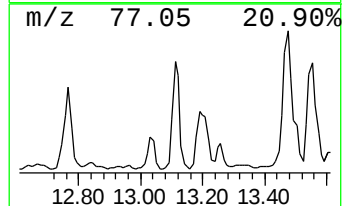
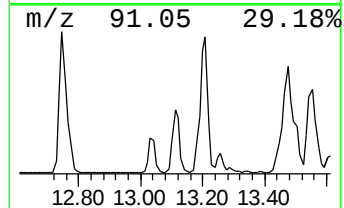
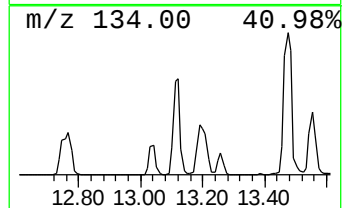
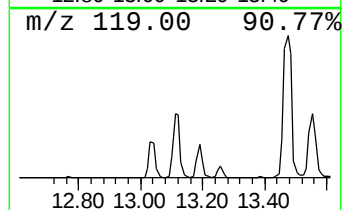
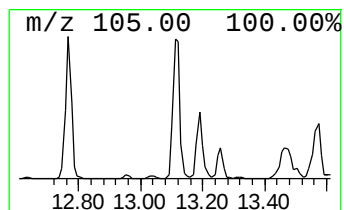
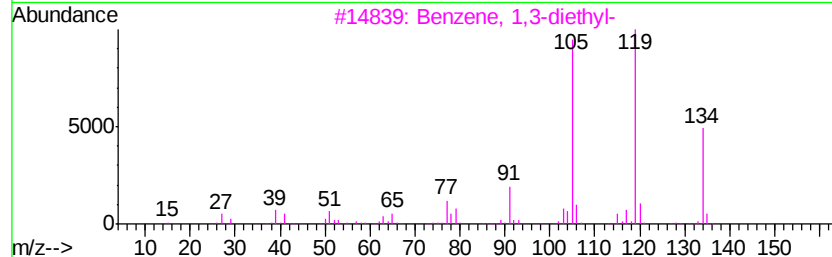
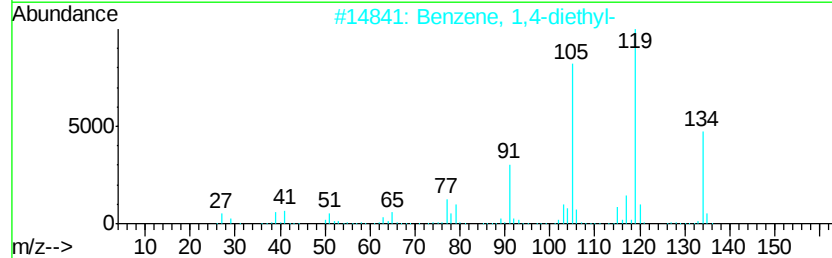
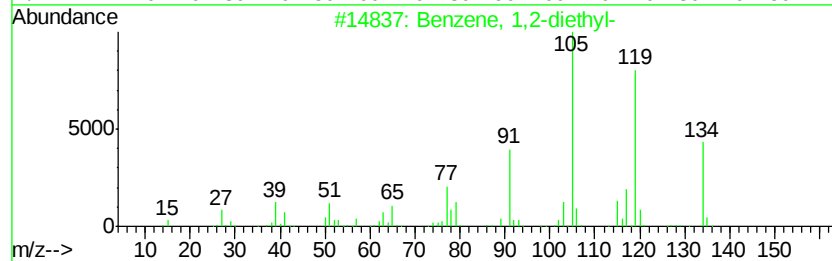
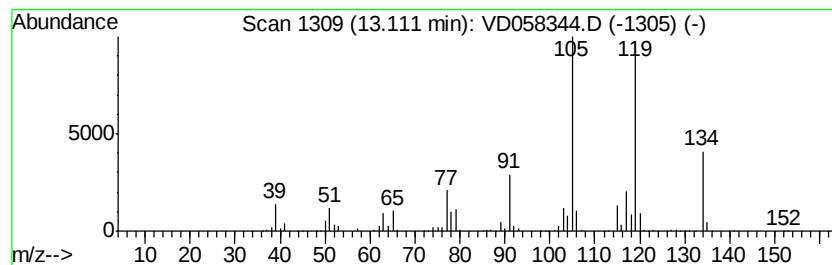
Instrument :
MSVOA_D
ClientSampleId :
DUP-0485-180619

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1,2-diethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.11	140.06 ug/l	21216300	1,4-Dichlorobenzene-d4	12.91
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2-diethyl-	134 C10H14	000135-01-3 96
2		Benzene, 1,4-diethyl-	134 C10H14	000105-05-5 90
3		Benzene, 1,3-diethyl-	134 C10H14	000141-93-5 90
4		Benzene, 1-ethyl-3,5-dimethyl-	134 C10H14	000934-74-7 76
5		Benzene, 1,2,3,4-tetramethyl-	134 C10H14	000488-23-3 62



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Instrument :
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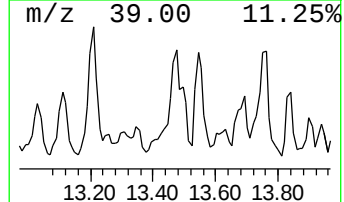
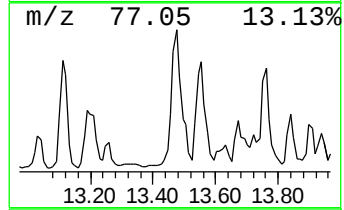
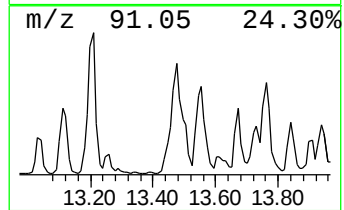
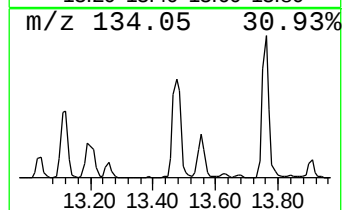
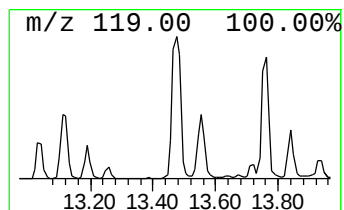
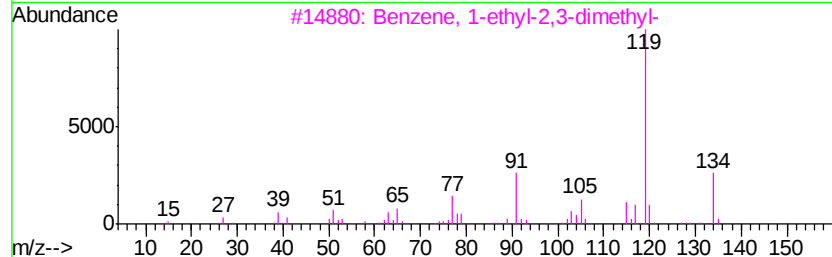
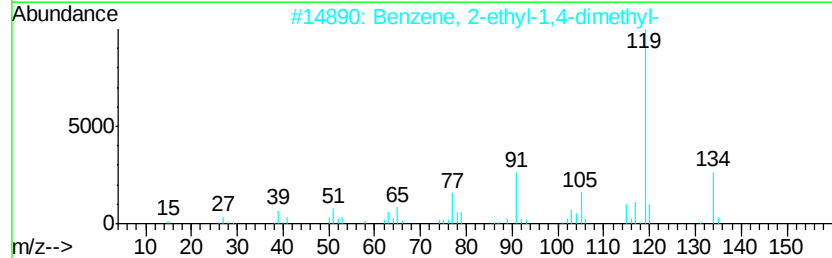
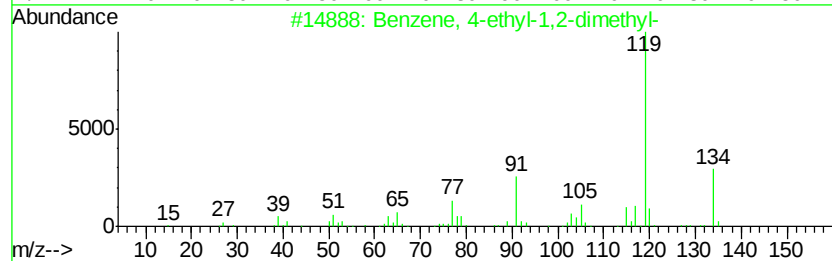
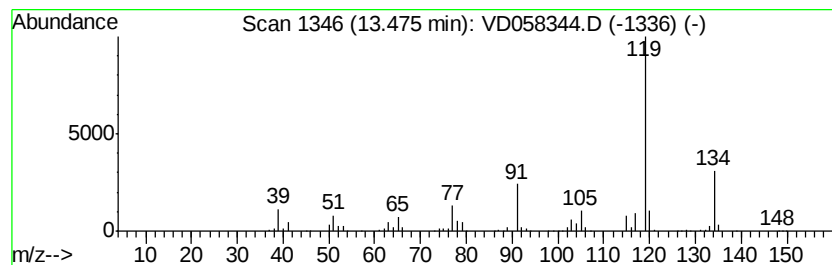
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	321.56 ug/l	48710100	1,4-Dichlorobenzene-d4	12.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4		o-Cymene	134	C10H14	000527-84-4	94
5		p-Cymene	134	C10H14	000099-87-6	94



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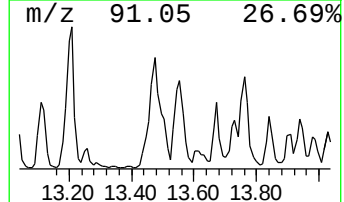
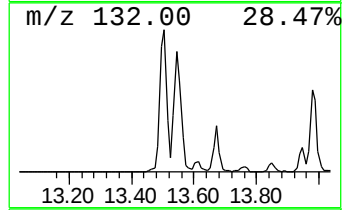
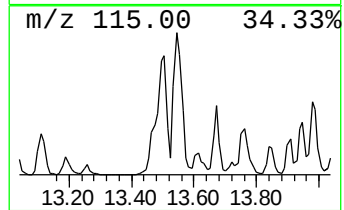
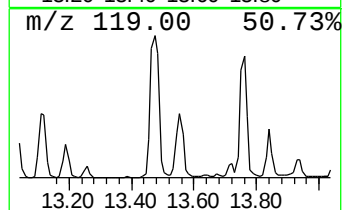
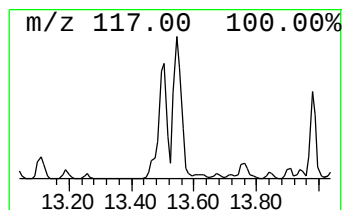
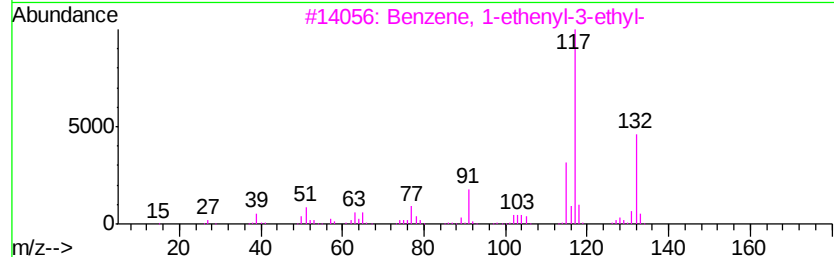
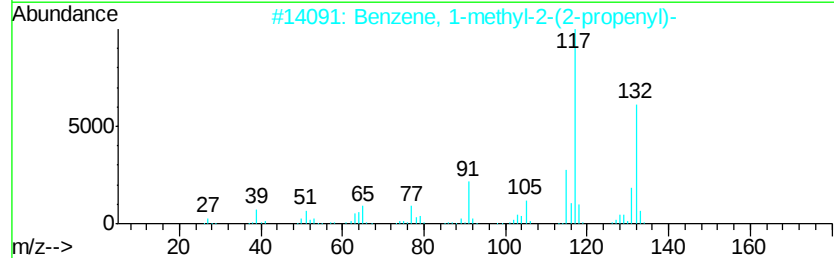
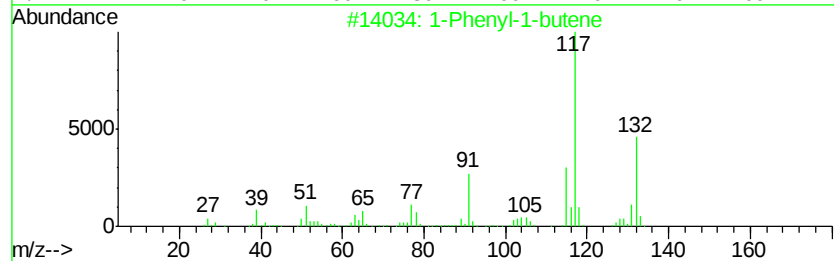
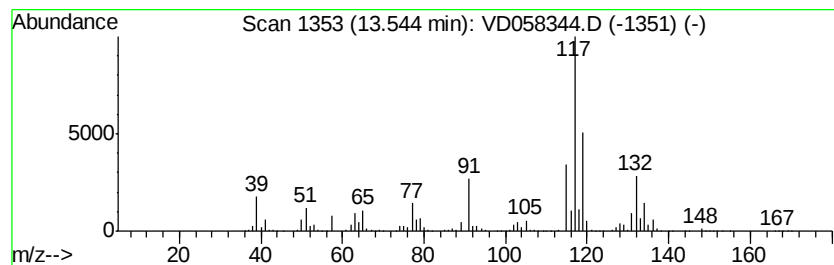
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.M
Quant Title : SW846 8260

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

Peak Number 5 1-Phenyl-1-butene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	192.61 ug/l	29177200	1,4-Dichlorobenzene-d4	12.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	70
2		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	70
4		Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	70
5		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	70



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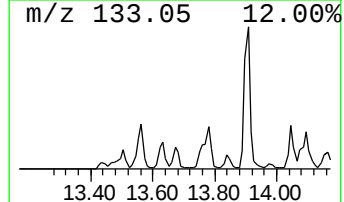
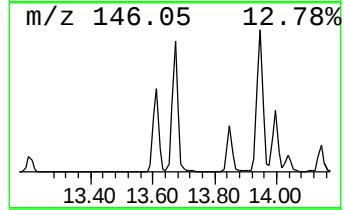
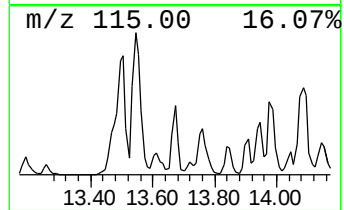
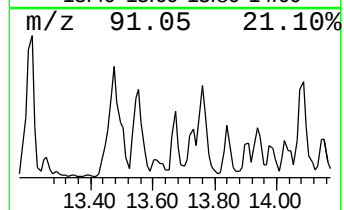
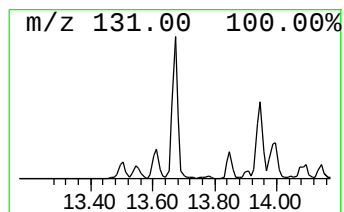
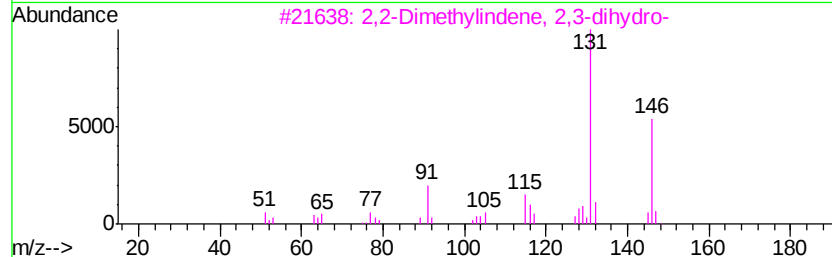
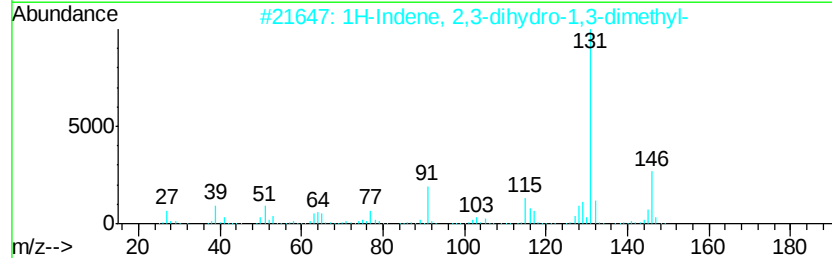
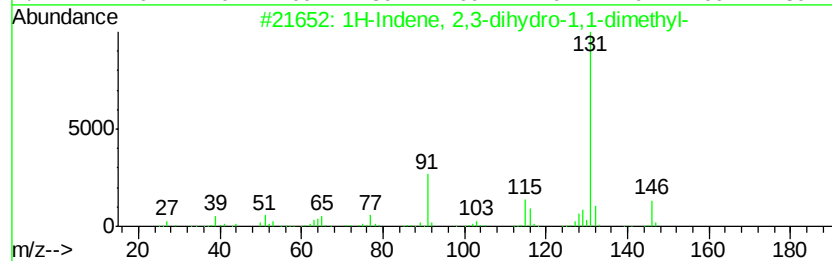
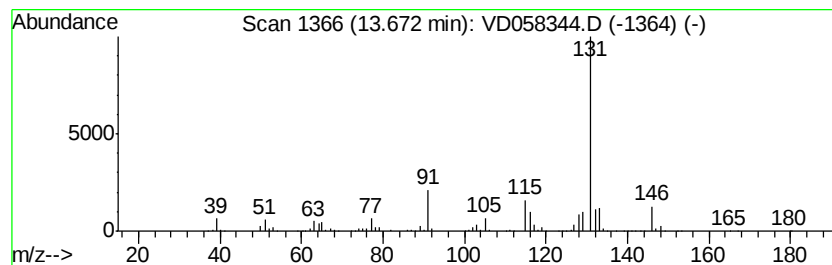
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.M
Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 2,3-dihydro-1,1-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.67	125.44 ug/l	19001400	1,4-Dichlorobenzene-d4	12.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	93
2		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90
3		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	87
4		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	86
5		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD062318\
Data File : VD058344.D
Acq On : 23 Jun 2018 16:46
Operator : VA/AP
Sample : J3664-19
Misc : 5.33g/5ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
DUP-0485-180619

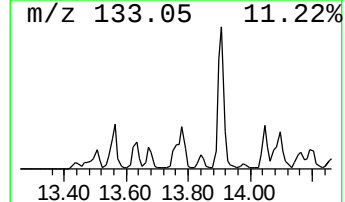
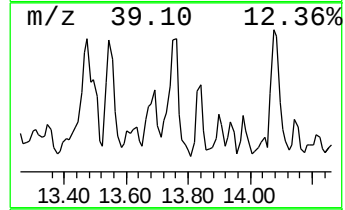
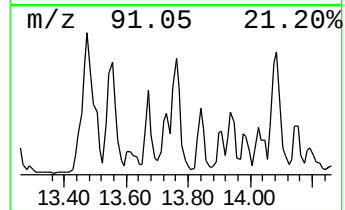
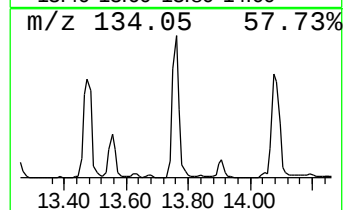
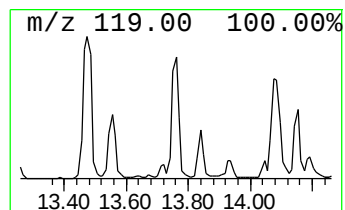
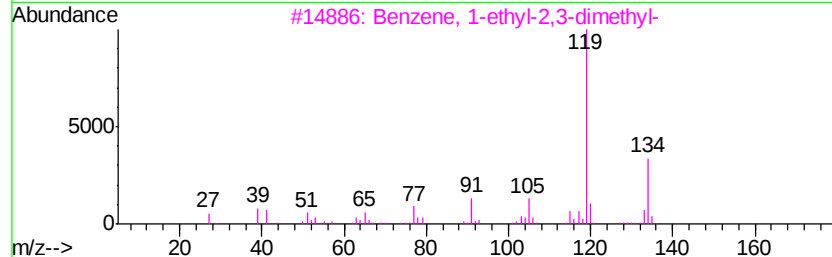
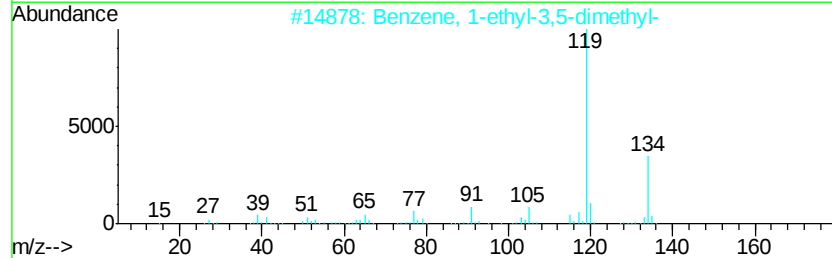
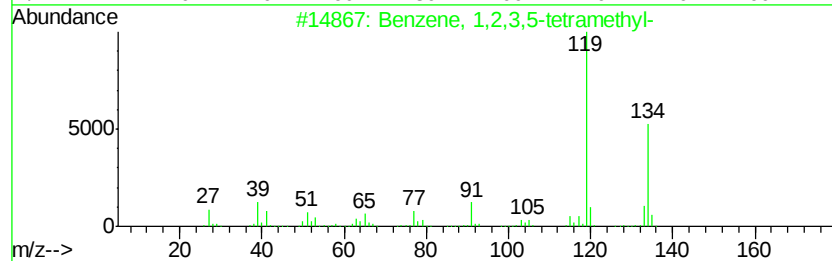
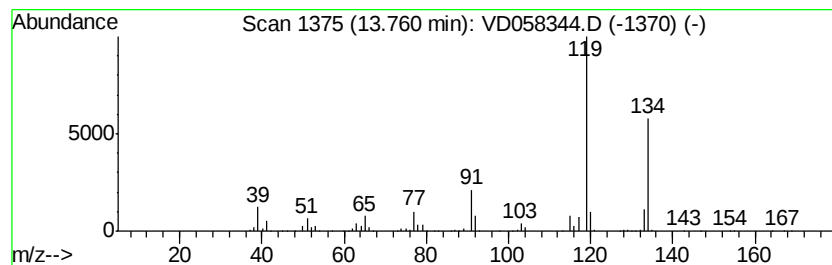
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	205.17 ug/l	31079100	1,4-Dichlorobenzene-d4	12.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
2		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
4		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91
5		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD062318\
Data File : VD058344.D
Acq On : 23 Jun 2018 16:46
Operator : VA/AP
Sample : J3664-19
Misc : 5.33g/5ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
DUP-0485-180619

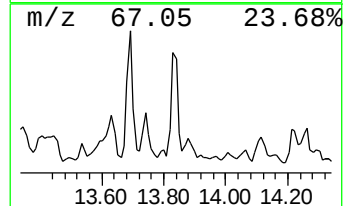
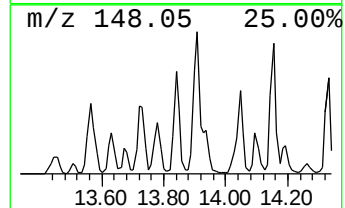
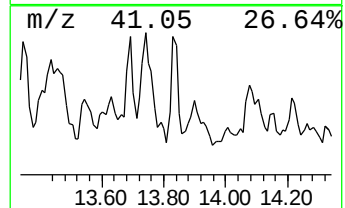
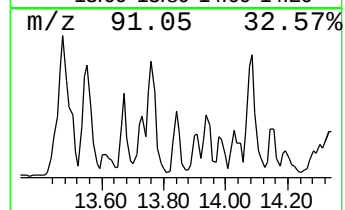
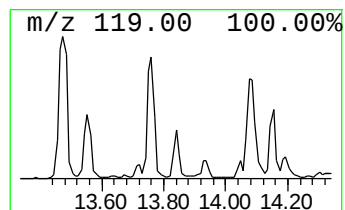
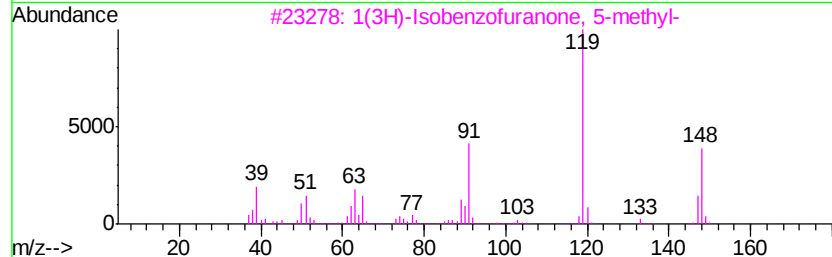
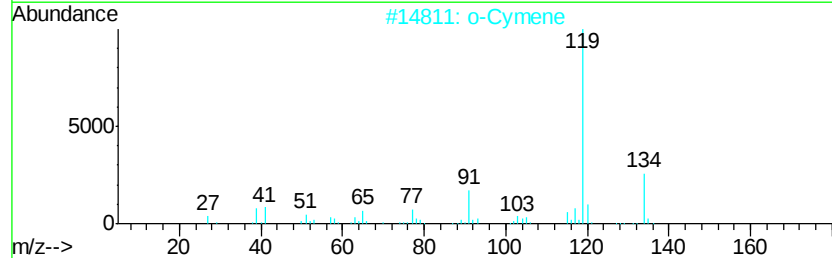
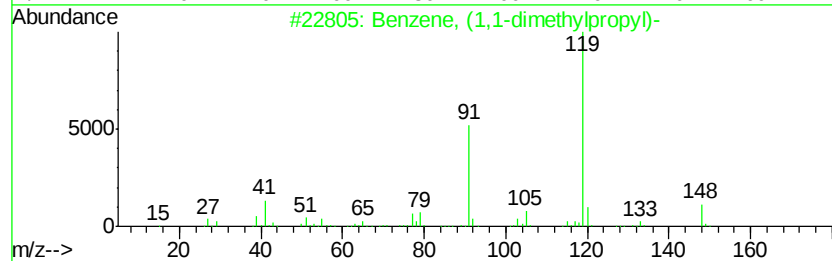
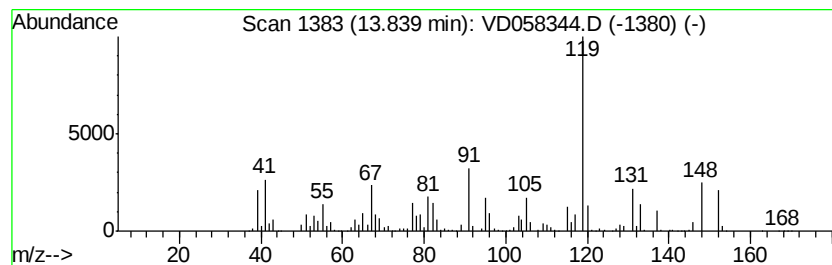
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Quant Title : SW846 8260

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

Peak Number 8 unknown13.84 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.84	95.34 ug/l	14441400	1,4-Dichlorobenzene-d4	12.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	46
2		o-Cymene	134	C10H14	000527-84-4	42
3		1(3H)-Isobenzofuranone, 5-methyl-	148	C9H8O2	054120-64-8	38
4		4'-Methylpropiophenone	148	C10H12O	005337-93-9	35
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	35



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD062318\
Data File : VD058344.D
Acq On : 23 Jun 2018 16:46
Operator : VA/AP
Sample : J3664-19
Misc : 5.33g/5ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
DUP-0485-180619

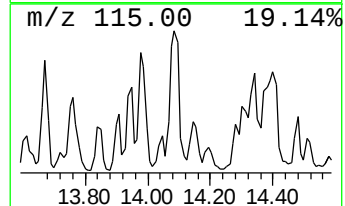
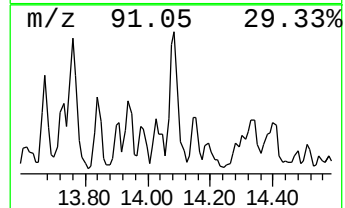
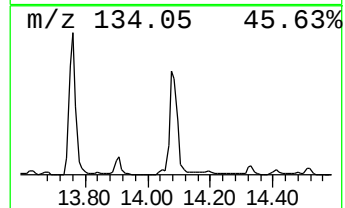
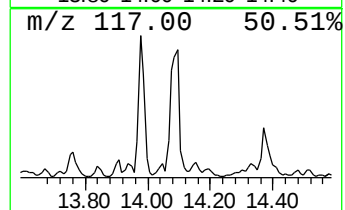
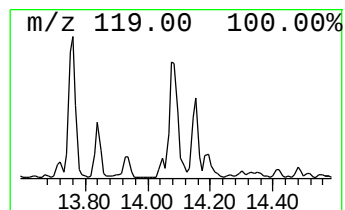
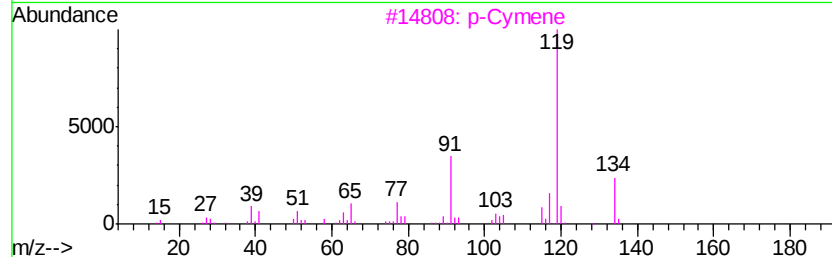
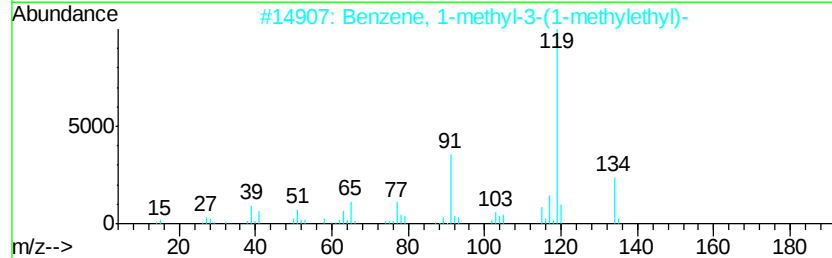
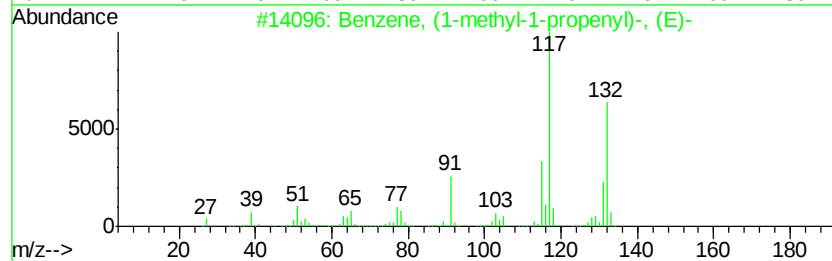
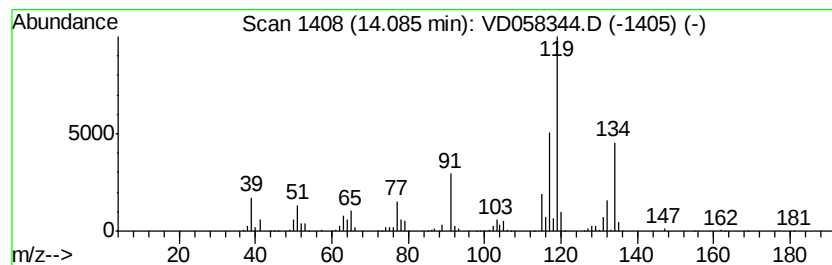
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.09	196.94 ug/l	29833000	1,4-Dichlorobenzene-d4	12.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	74
2		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	70
3		p-Cymene	134	C10H14	000099-87-6	70
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	64
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD062318\
Data File : VD058344.D
Acq On : 23 Jun 2018 16:46
Operator : VA/AP
Sample : J3664-19
Misc : 5.33g/5ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
DUP-0485-180619

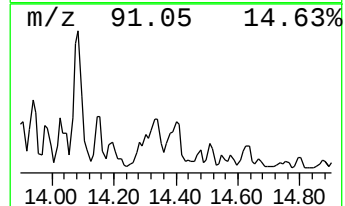
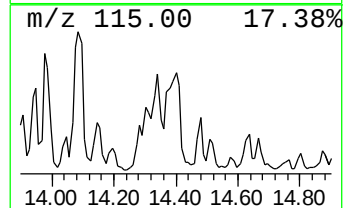
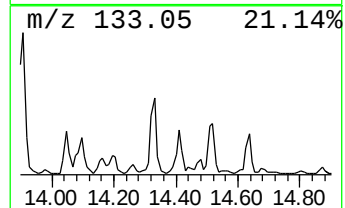
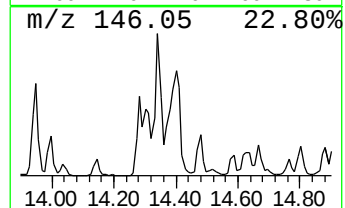
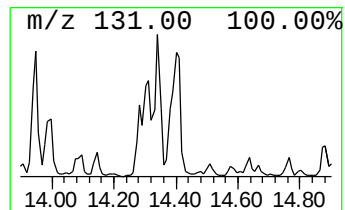
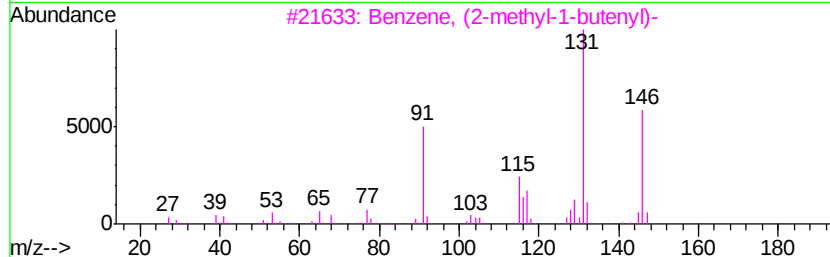
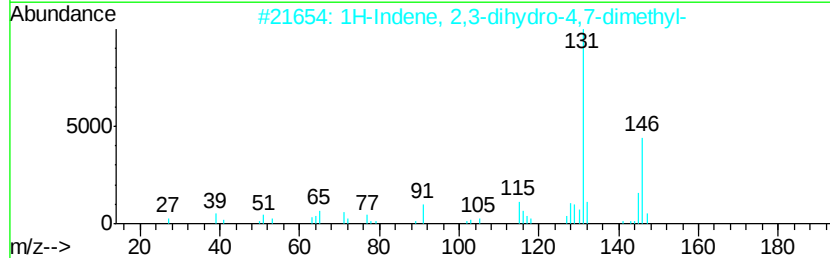
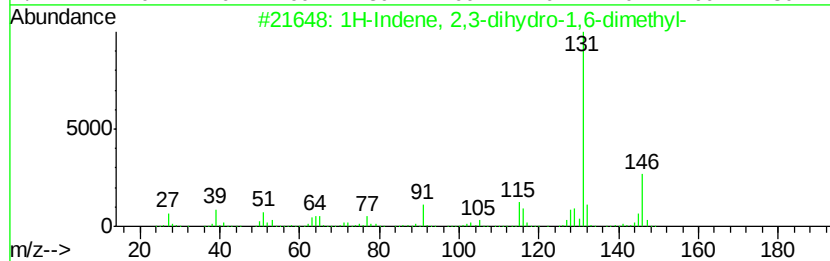
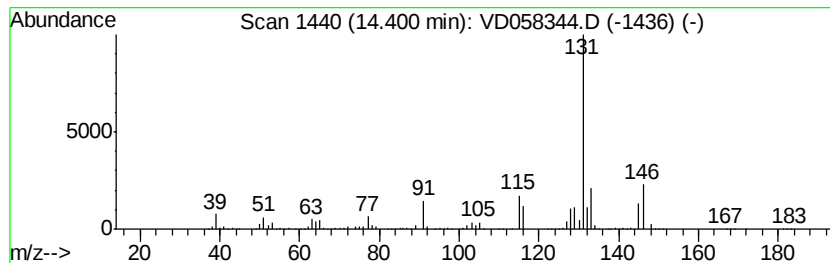
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.M
Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.40	114.37 ug/l	17325300	1,4-Dichlorobenzene-d4	12.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	87
2		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87
3		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	87
4		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	87
5		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	87



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD062318\
Data File : VD058344.D
Acq On : 23 Jun 2018 16:46
Operator : VA/AP
Sample : J3664-19
Misc : 5.33g/5ml/MSVOA_D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
DUP-0485-180619

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D061918S.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, 1,2-diet...	13.11	140.1	ug/l	21216300	4	12.91	7574000	50.0
Benzene, 4-ethyl-...	13.47	321.6	ug/l	48710100	4	12.91	7574000	50.0
1-Phenyl-1-butene	13.54	192.6	ug/l	29177200	4	12.91	7574000	50.0
1H-Indene, 2,3-di...	13.67	125.4	ug/l	19001400	4	12.91	7574000	50.0
Benzene, 1,2,3,5-...	13.76	205.2	ug/l	31079100	4	12.91	7574000	50.0
unknown13.84	13.84	95.3	ug/l	14441400	4	12.91	7574000	50.0
Benzene, (1-methy...	14.09	196.9	ug/l	29833000	4	12.91	7574000	50.0
1H-Indene, 2,3-di...	14.40	114.4	ug/l	17325300	4	12.91	7574000	50.0