

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD041819\
 Data File : VD061897.D
 Acq On : 19 Apr 2019 5:29
 Operator : VA/AP
 Sample : K2197-07
 Misc : 5.72a/5ml/MSVOA D/SOIL
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 CL-01-041619-B

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D041519S.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	1.495	4	7	21	rVB	188358	592021	26.56%	1.601%
2	3.829	239	244	252	rBV	51284	131514	5.90%	0.356%
3	6.162	475	481	485	rBV2	9877	28220	1.27%	0.076%
4	6.939	551	560	567	rBV2	293293	847790	38.04%	2.293%
5	7.057	567	572	584	rVB	333973	927311	41.60%	2.508%
6	7.471	607	614	629	rVB	205195	566537	25.42%	1.532%
7	8.229	686	691	708	rBV	450066	1113875	49.97%	3.012%
8	10.424	908	914	921	rBV	472408	1100039	49.35%	2.975%
9	12.373	1108	1112	1119	rVB	545695	858520	38.52%	2.322%
10	12.531	1124	1128	1132	rVB	30916	43435	1.95%	0.117%
11	12.669	1138	1142	1147	rBV	78771	141846	6.36%	0.384%
12	12.748	1147	1150	1158	rVB	28518	46428	2.08%	0.126%
13	13.053	1178	1181	1185	rVB	166211	235727	10.58%	0.637%
14	13.230	1193	1199	1202	rBV	22910	39392	1.77%	0.107%
15	13.318	1202	1208	1214	rVV3	26523	68519	3.07%	0.185%
16	13.407	1214	1217	1220	rVV2	26907	45582	2.04%	0.123%
17	13.496	1220	1226	1228	rVV3	11715	29281	1.31%	0.079%
18	13.565	1228	1233	1238	rVV	417490	682010	30.60%	1.844%
19	13.673	1241	1244	1246	rVV	35588	57008	2.56%	0.154%
20	13.722	1246	1249	1252	rVV	33185	55790	2.50%	0.151%
21	13.781	1252	1255	1258	rVV	86906	118865	5.33%	0.321%
22	13.850	1258	1262	1264	rVV	388160	596386	26.76%	1.613%
23	13.889	1264	1266	1268	rVV	215712	312118	14.00%	0.844%
24	13.939	1268	1271	1273	rVV	245563	436719	19.59%	1.181%
25	13.968	1273	1274	1278	rVV	248924	298684	13.40%	0.808%
26	14.086	1282	1286	1292	rVV	386024	664354	29.81%	1.797%
27	14.175	1292	1295	1296	rVV2	30876	59274	2.66%	0.160%
28	14.204	1296	1298	1300	rVV	100660	173224	7.77%	0.468%
29	14.244	1300	1302	1309	rVV	773658	1080417	48.47%	2.922%
30	14.332	1309	1311	1312	rVV	45324	67674	3.04%	0.183%
31	14.372	1312	1315	1319	rVV2	97985	241091	10.82%	0.652%
32	14.441	1319	1322	1325	rVV2	190745	329208	14.77%	0.890%
33	14.520	1325	1330	1332	rVV2	388739	709141	31.81%	1.918%
34	14.569	1332	1335	1339	rVV	561214	1001433	44.93%	2.708%

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Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
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35	14.657	1339	1344	1346	rVV3	145015	321447	14.42%	0.869%
36	14.716	1346	1350	1351	rVV	759003	1016092	45.59%	2.748%
37	14.736	1351	1352	1355	rVV	350667	465881	20.90%	1.260%
38	14.785	1355	1357	1362	rVV2	624807	1058955	47.51%	2.864%
39	14.854	1362	1364	1365	rVV	121381	178173	7.99%	0.482%
40	14.894	1365	1368	1369	rVV	527352	738100	33.11%	1.996%
41	14.923	1369	1371	1375	rVV	277280	528909	23.73%	1.430%
42	14.982	1375	1377	1378	rVV	336750	417565	18.73%	1.129%
43	15.012	1378	1380	1383	rVV	396421	693750	31.12%	1.876%
44	15.061	1383	1385	1387	rVV	604818	781418	35.06%	2.113%
45	15.100	1387	1389	1391	rVV	233287	357216	16.03%	0.966%
46	15.140	1391	1393	1399	rVV2	582941	1261746	56.61%	3.412%
47	15.228	1399	1402	1404	rVV4	156545	368160	16.52%	0.996%
48	15.278	1404	1407	1410	rVV2	296381	649439	29.14%	1.756%
49	15.327	1410	1412	1413	rVV	234673	364916	16.37%	0.987%
50	15.356	1413	1415	1416	rVV	348174	521031	23.38%	1.409%
51	15.386	1416	1418	1421	rVV	539738	810938	36.38%	2.193%
52	15.435	1421	1423	1426	rVV3	331561	660928	29.65%	1.787%
53	15.474	1426	1427	1430	rVV3	205379	357106	16.02%	0.966%
54	15.573	1430	1437	1440	rVV2	566537	1437121	64.47%	3.886%
55	15.612	1440	1441	1442	rVV	248591	243722	10.93%	0.659%
56	15.652	1442	1445	1446	rVV2	417699	765894	34.36%	2.071%
57	15.681	1446	1448	1453	rVV	1429673	2228973	100.00%	6.028%
58	15.750	1453	1455	1458	rVV3	271852	523454	23.48%	1.416%
59	15.799	1458	1460	1464	rVV	935435	1426705	64.01%	3.858%
60	15.908	1465	1471	1473	rVV2	289763	1001406	44.93%	2.708%
61	15.937	1473	1474	1476	rVV	361616	394999	17.72%	1.068%
62	16.006	1476	1481	1485	rVV2	383548	1072782	48.13%	2.901%
63	16.055	1485	1486	1491	rVV5	123941	333834	14.98%	0.903%
64	16.144	1491	1495	1497	rVV5	124341	363631	16.31%	0.983%
65	16.183	1497	1499	1501	rVV	294853	386879	17.36%	1.046%
66	16.242	1501	1505	1507	rVV	245943	535188	24.01%	1.447%
67	16.321	1509	1513	1515	rVV4	71281	195879	8.79%	0.530%
68	16.360	1515	1517	1522	rVV	97977	208383	9.35%	0.564%
69	16.429	1522	1524	1526	rVV2	33450	65460	2.94%	0.177%
70	16.488	1526	1530	1531	rVV2	78986	155640	6.98%	0.421%
71	16.508	1531	1532	1538	rVV	165505	244279	10.96%	0.661%

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Instrument :
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Integration Parameters: RTEINT.P

Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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72	16.587	1538	1540	1542	rVV3	20609	35484	1.59%	0.096%
73	16.636	1542	1545	1548	rVV	36067	56736	2.55%	0.153%
74	16.734	1552	1555	1558	rBV	36059	52148	2.34%	0.141%

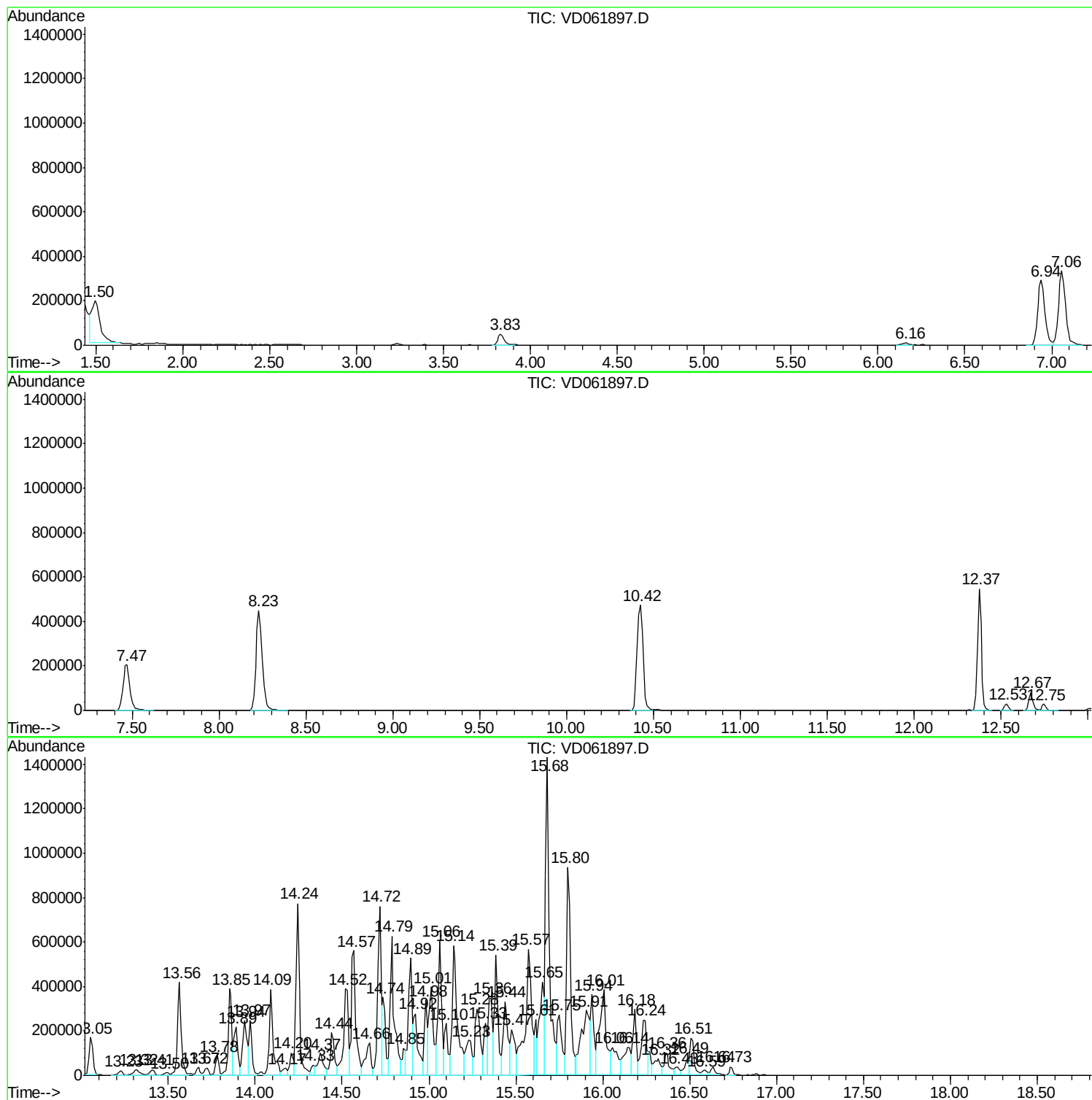
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Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041819\
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ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D041519S.M
Quant Title : SW846 8260

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



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Data File : VD061897.D
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Operator : VA/AP
Sample : K2197-07
Misc : 5.72a/5ml/MSVOA D/SOIL
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Instrument :
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CL-01-041619-B

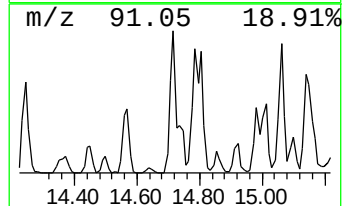
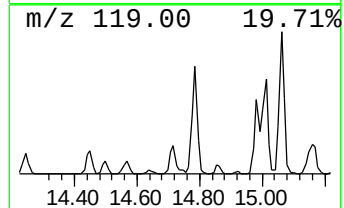
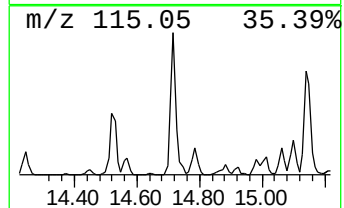
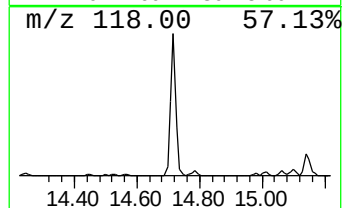
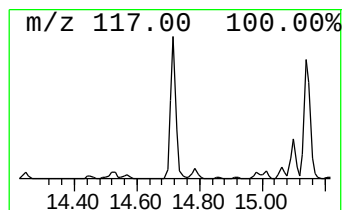
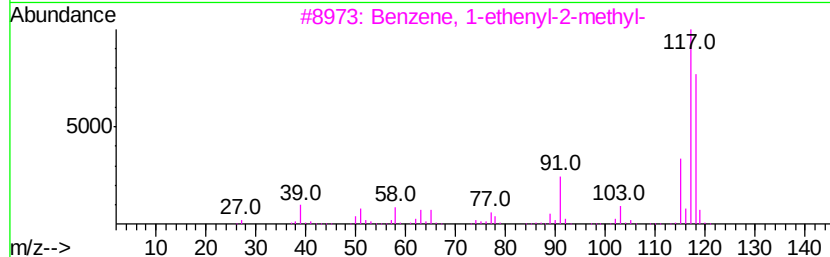
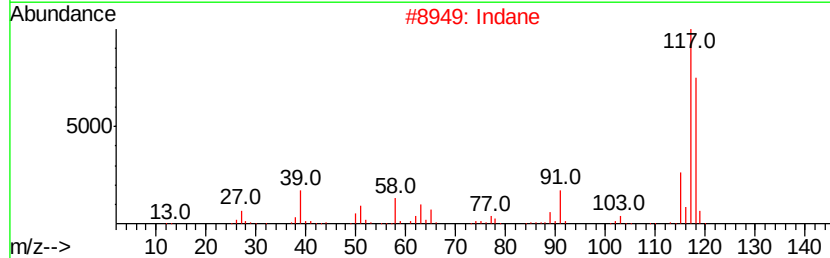
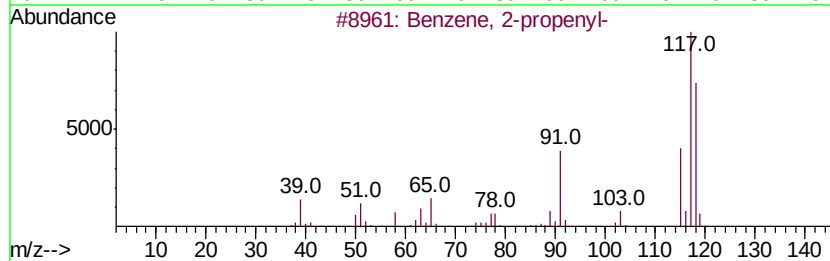
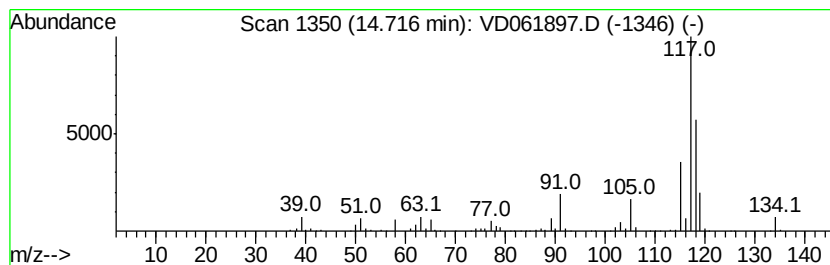
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D041519S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 2-propenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.72	71.64 ug/l	1016090	1,4-Dichlorobenzene-d4	14.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
2		Indane	118	C9H10	000496-11-7	76
3		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	70
4		Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
5		1,4:5,8-Dimethanobiphenylene, 1,...	184	C14H16	001624-13-1	59



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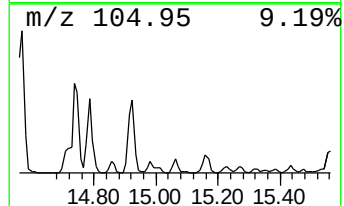
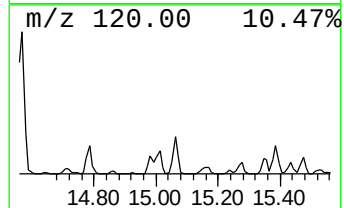
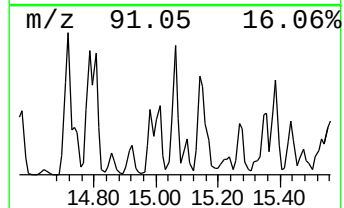
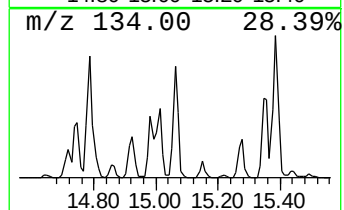
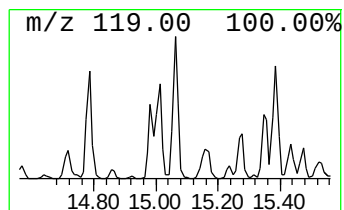
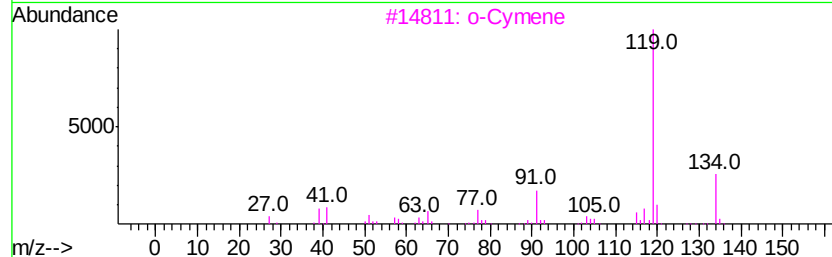
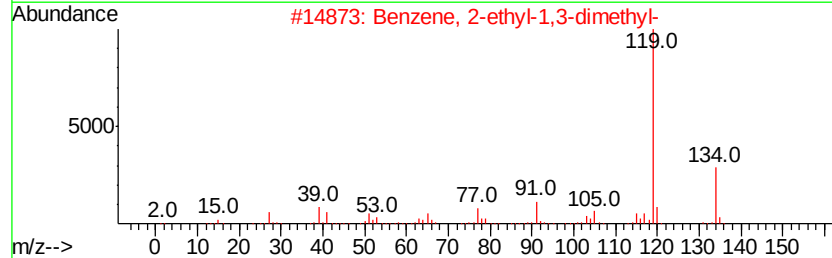
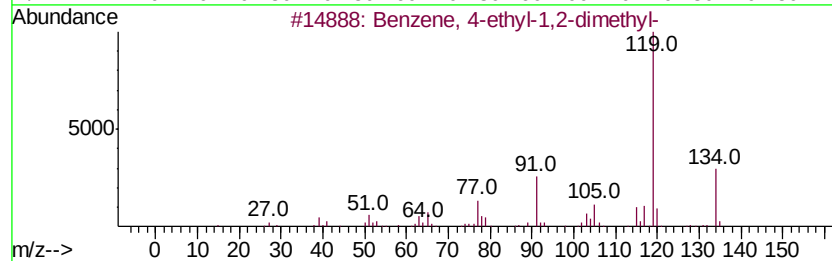
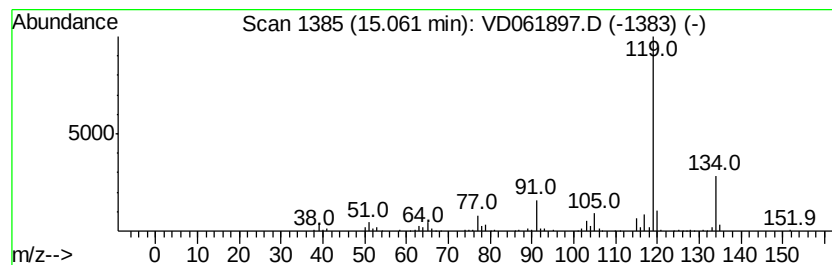
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D041519S.M
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 2 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.06	55.10 ug/l	781418	1,4-Dichlorobenzene-d4	14.53

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-1,3-dimethyl-	134	C10H14	002870-04-4	95
3		o-Cymene	134	C10H14	000527-84-4	95
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



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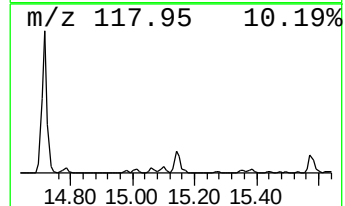
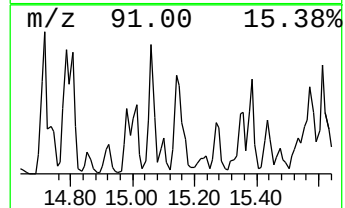
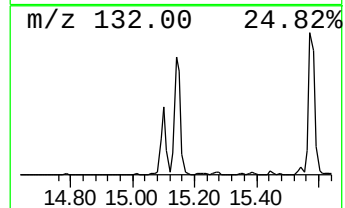
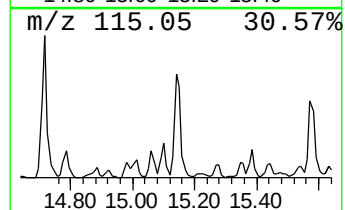
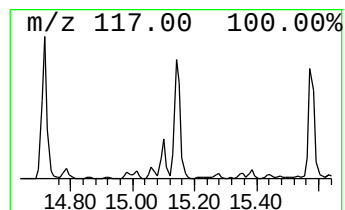
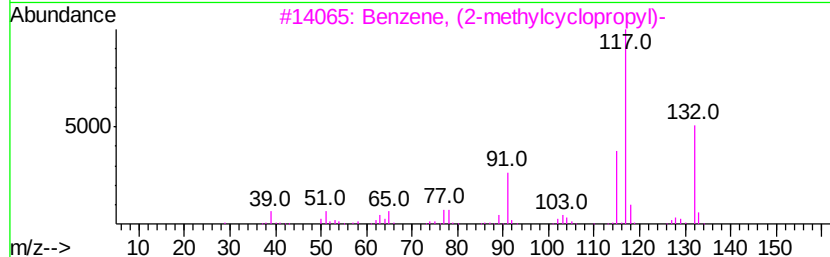
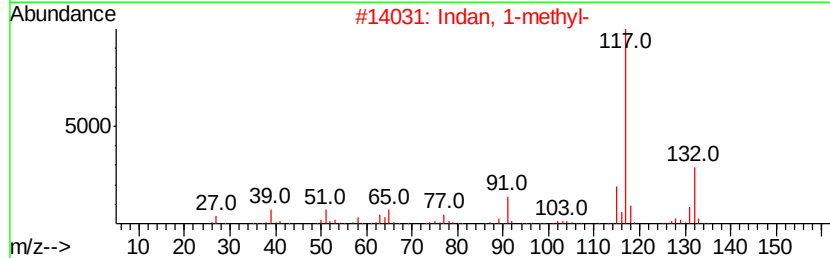
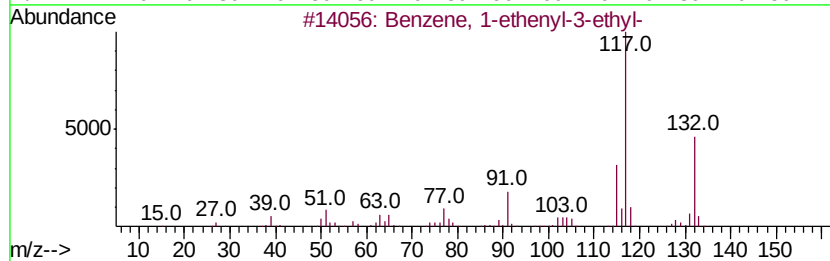
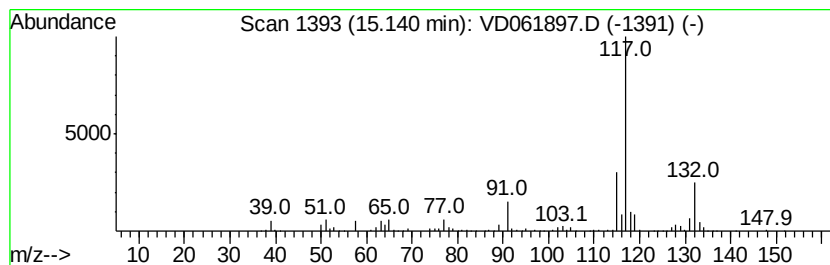
Instrument :
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ClientSampleId :
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.14	88.96 ug/l	1261750	1,4-Dichlorobenzene-d4	14.53
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4 90
2		Indan, 1-methyl-	132 C10H12	000767-58-8 87
3		Benzene, (2-methylcyclopropyl)-	132 C10H12	1000327-39-0 83
4		Benzene, 1-methyl-2-(2-propenyl)-	132 C10H12	001587-04-8 83
5		Benzene, 2-butenyl-	132 C10H12	001560-06-1 80



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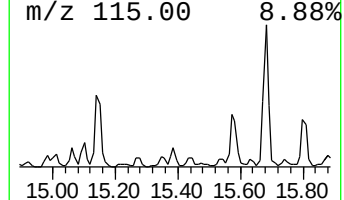
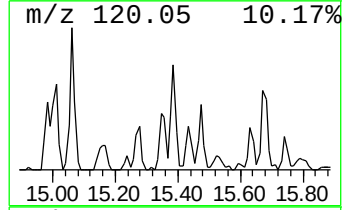
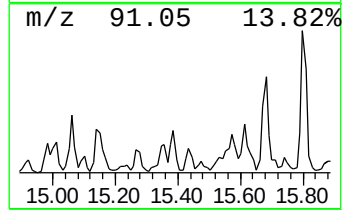
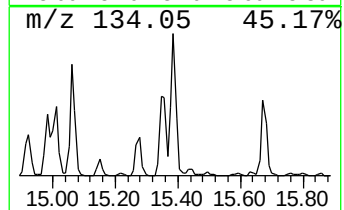
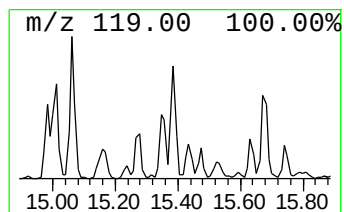
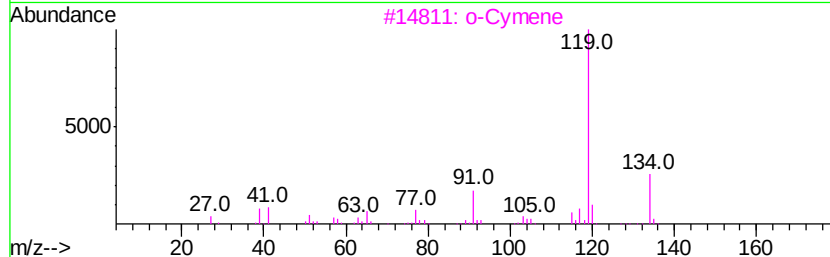
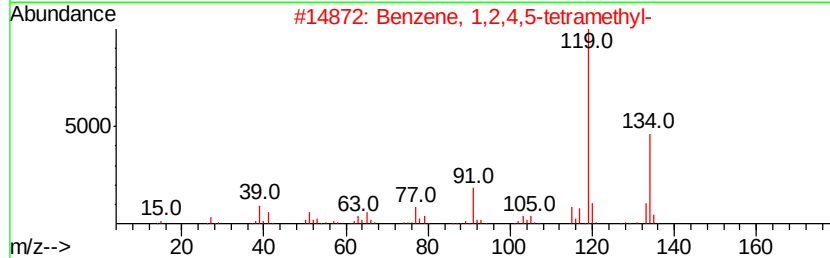
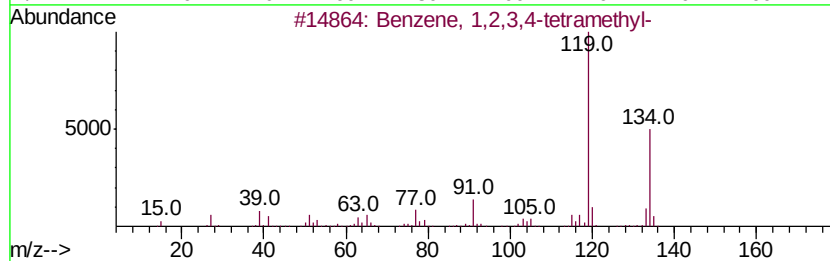
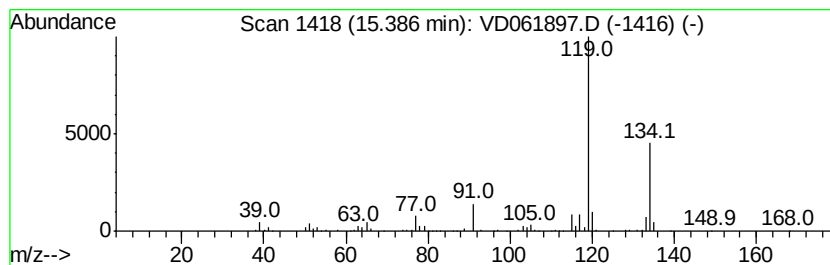
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.39	57.18 ug/l	810938	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		o-Cymene	134	C10H14	000527-84-4	95
4		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041819\
Data File : VD061897.D
Acq On : 19 Apr 2019 5:29
Operator : VA/AP
Sample : K2197-07
Misc : 5.72a/5ml/MSVOA D/SOIL
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
CL-01-041619-B

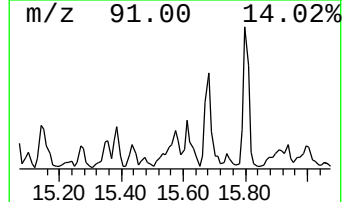
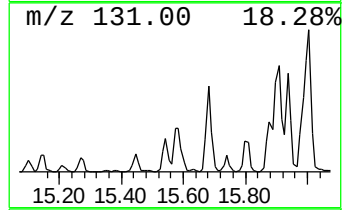
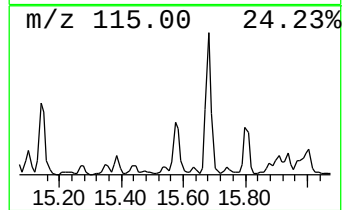
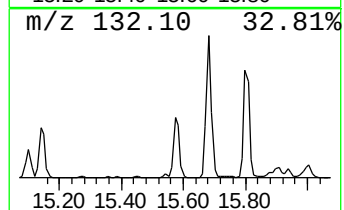
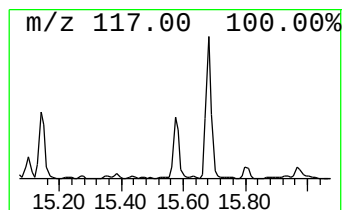
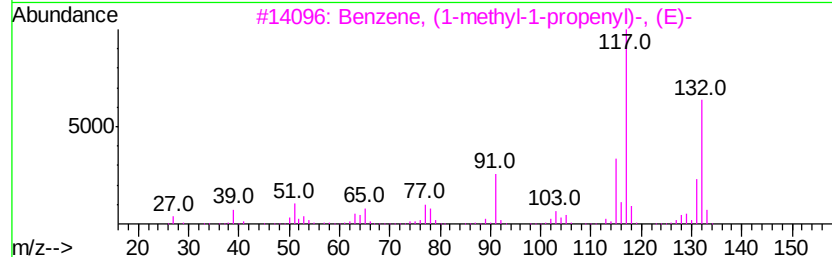
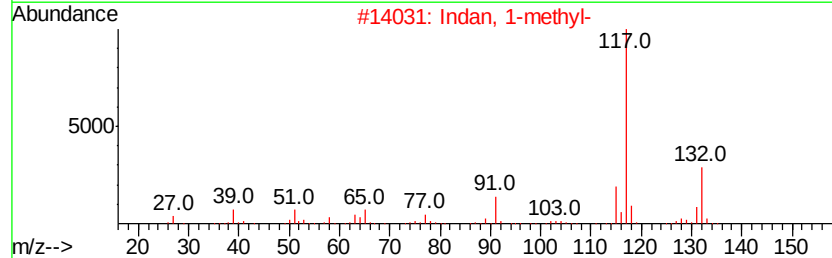
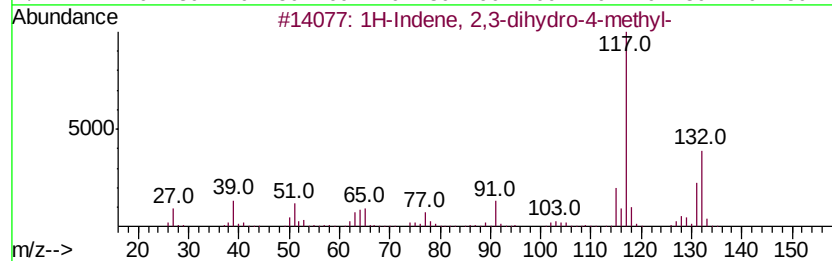
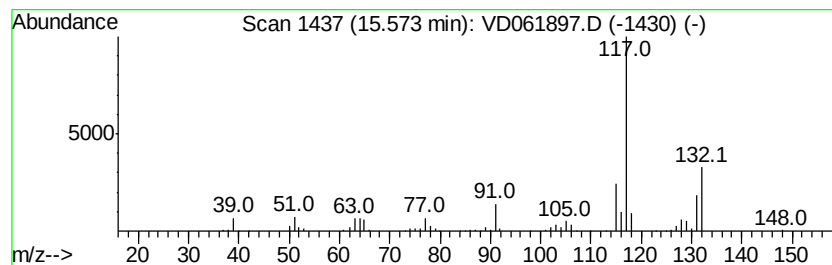
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Quant Title : SW846 8260

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

Peak Number 5 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.57	101.33 ug/l	1437120	1,4-Dichlorobenzene-d4	14.53

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	91
2	Indan, 1-methyl-	132	C10H12	000767-58-8	90
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	80
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041819\
Data File : VD061897.D
Acq On : 19 Apr 2019 5:29
Operator : VA/AP
Sample : K2197-07
Misc : 5.72a/5ml/MSVOA D/SOIL
ALS Vial : 41 Sample Multiplier: 1

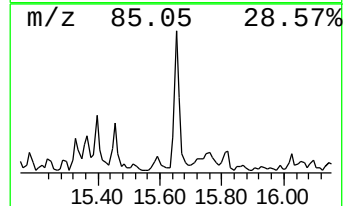
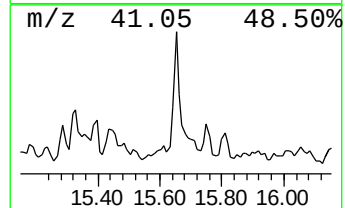
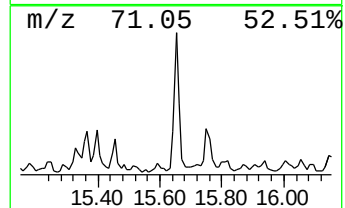
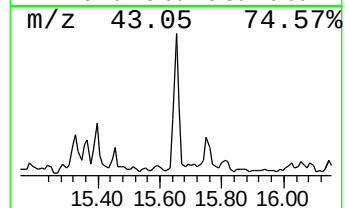
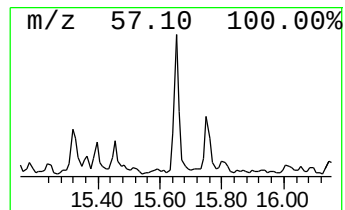
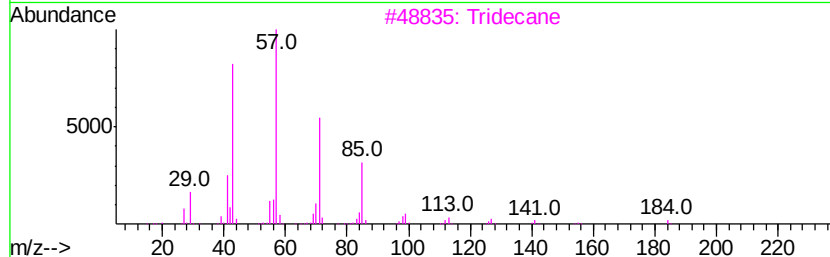
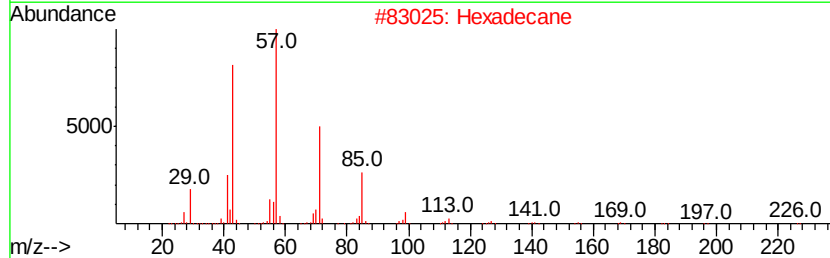
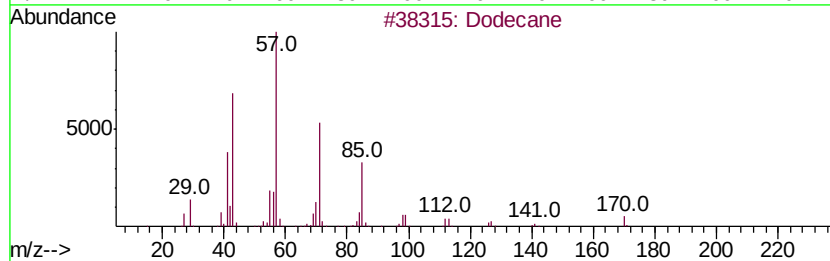
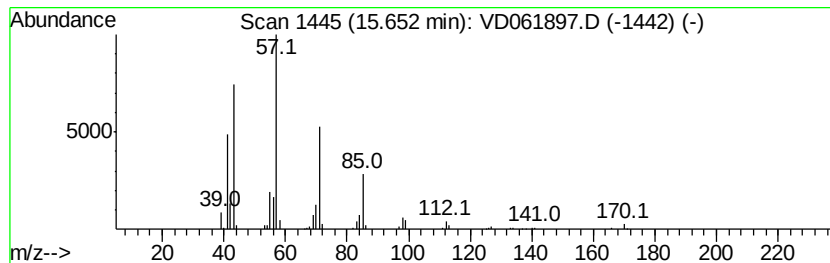
Instrument :
MSVOA_D
ClientSampleId :
CL-01-041619-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D041519S.M
Quant Title : SW846 8260

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

Peak Number 6 Dodecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.65	54.00 ug/l	765894	1,4-Dichlorobenzene-d4	14.53
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Dodecane	170 C12H26	000112-40-3	91
2	Hexadecane	226 C16H34	000544-76-3	86
3	Tridecane	184 C13H28	000629-50-5	83
4	Undecane	156 C11H24	001120-21-4	83
5	Decane, 6-ethyl-2-methyl-	184 C13H28	062108-21-8	78



Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD041819\
Data File : VD061897.D
Acq On : 19 Apr 2019 5:29
Operator : VA/AP
Sample : K2197-07
Misc : 5.72a/5ml/MSVOA D/SOIL
ALS Vial : 41 Sample Multiplier: 1

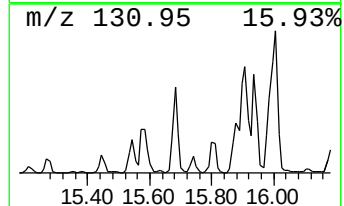
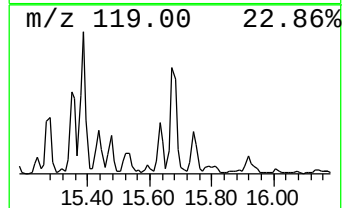
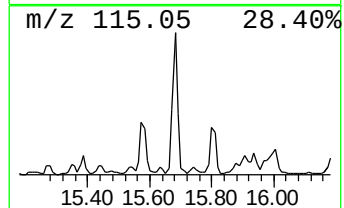
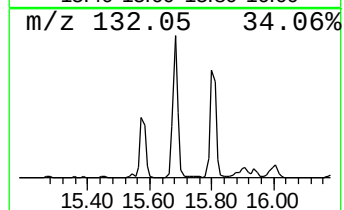
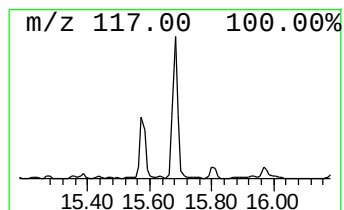
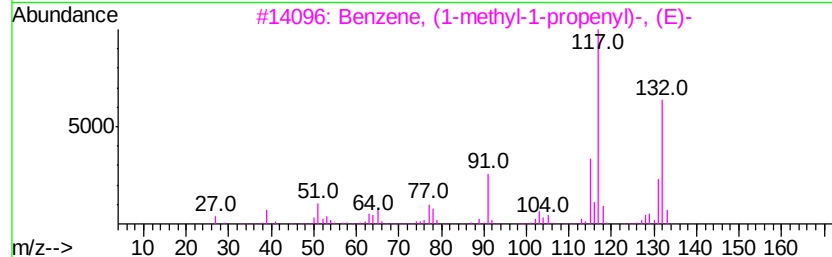
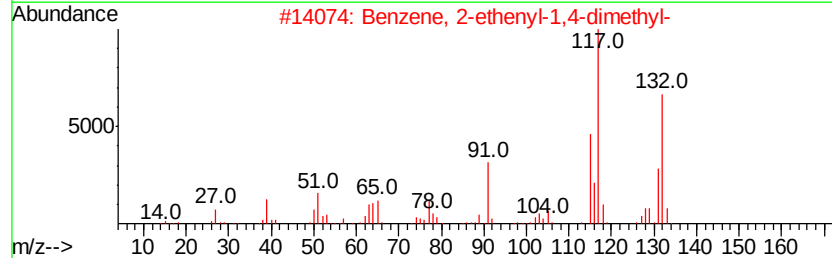
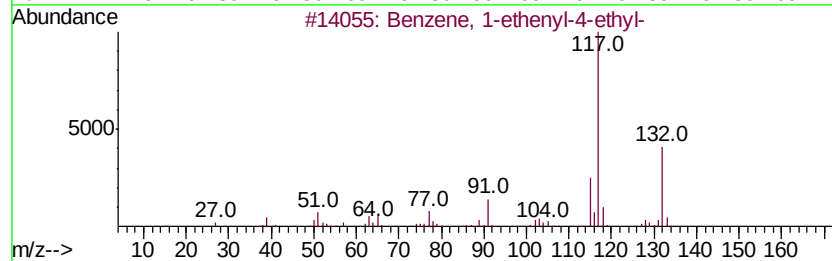
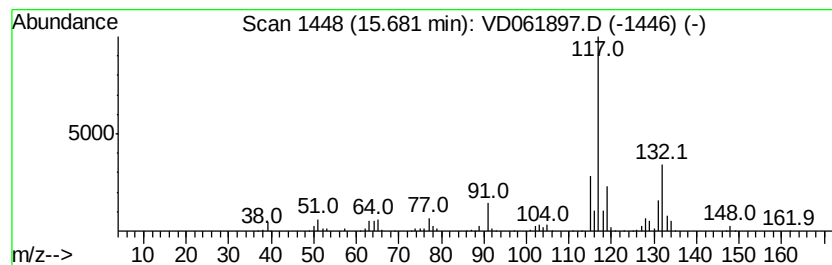
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CL-01-041619-B

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D041519S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.68	157.16 ug/l	2228970	1,4-Dichlorobenzene-d4	14.53
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-4-ethyl-	132 C10H12	003454-07-7 93
2		Benzene, 2-ethenyl-1,4-dimethyl-	132 C10H12	002039-89-6 91
3		Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000768-00-3 87
4		Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0 83
5		Benzene, 2-butenyl-	132 C10H12	001560-06-1 83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041819\
Data File : VD061897.D
Acq On : 19 Apr 2019 5:29
Operator : VA/AP
Sample : K2197-07
Misc : 5.72g/5ml/MSVOA D/SOIL
ALS Vial : 41 Sample Multiplier: 1

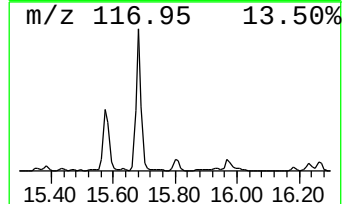
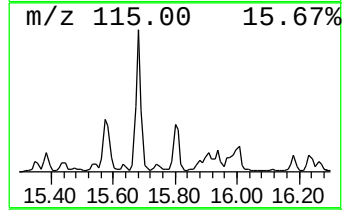
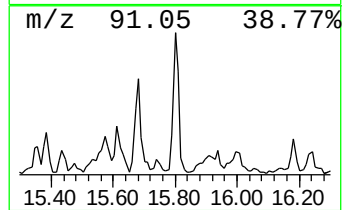
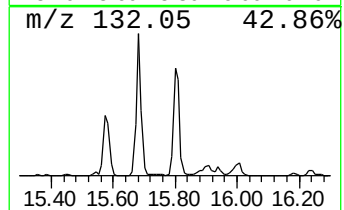
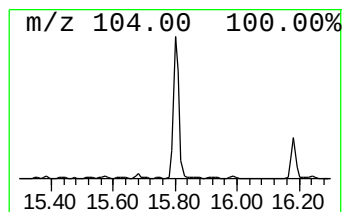
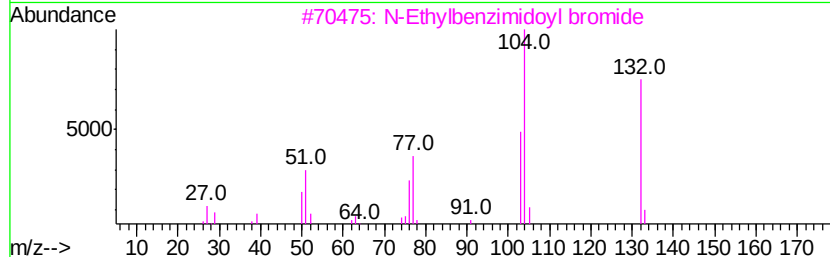
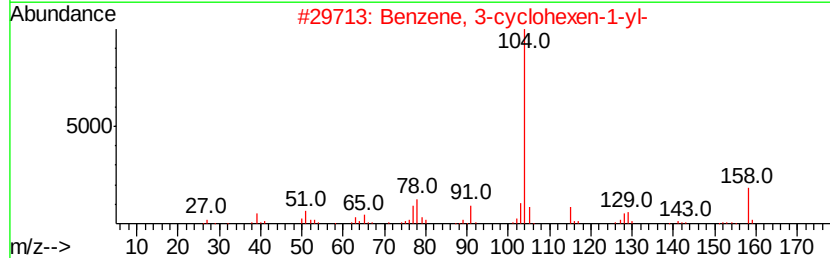
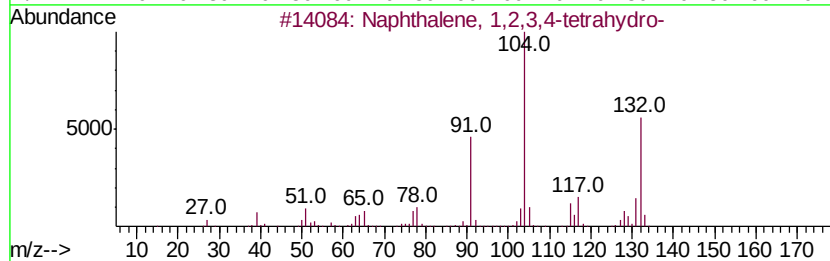
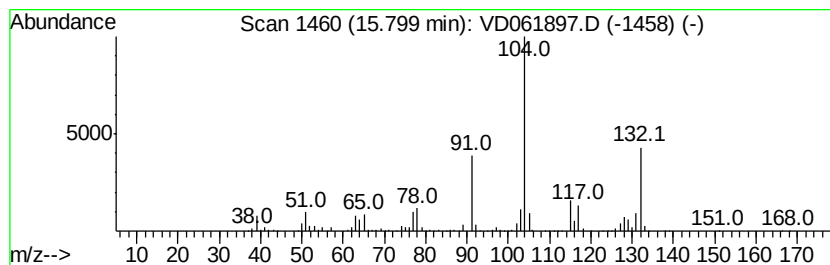
Instrument :
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ClientSampleId :
CL-01-041619-B

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D041519S.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 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.80	100.59 ug/l	1426710	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 91
2		Benzene, 3-cyclohexen-1-yl-	158 C12H14	004994-16-5 50
3		N-Ethylbenzimidoyl bromide	211 C9H10BrN	041182-87-0 50
4		Benzene, cyclobutyl-	132 C10H12	004392-30-7 42
5		3(2H)-Furanone, dihydro-2,2-dime...	190 C12H14O2	063678-00-2 38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041819\
Data File : VD061897.D
Acq On : 19 Apr 2019 5:29
Operator : VA/AP
Sample : K2197-07
Misc : 5.72a/5ml/MSVOA D/SOIL
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
CL-01-041619-B

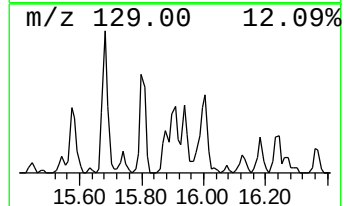
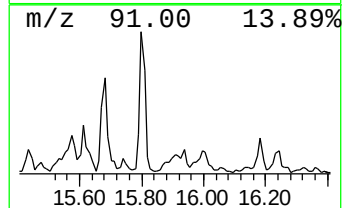
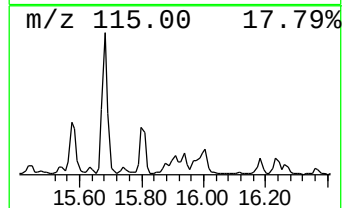
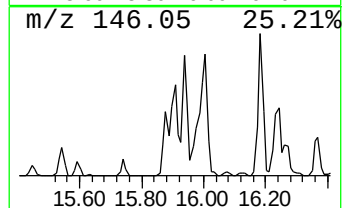
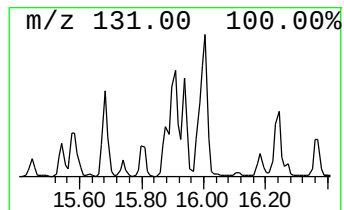
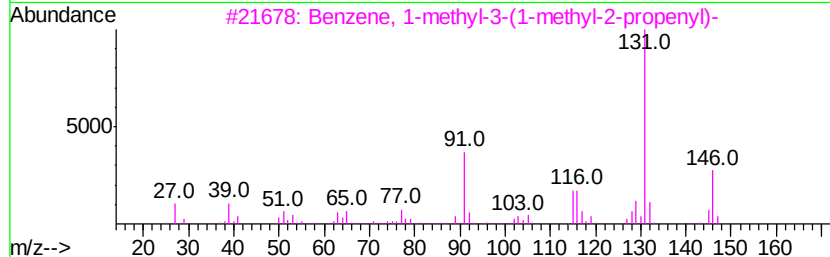
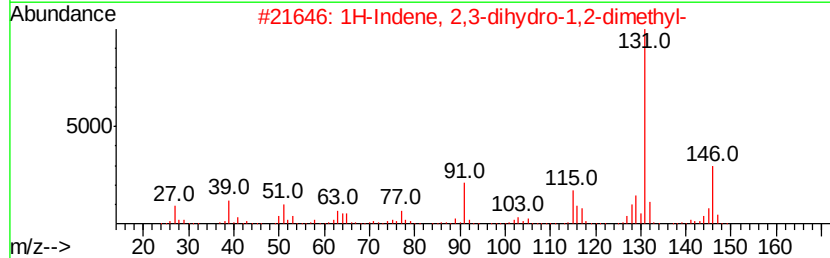
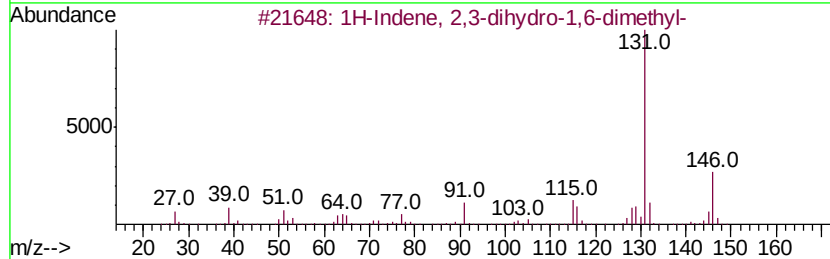
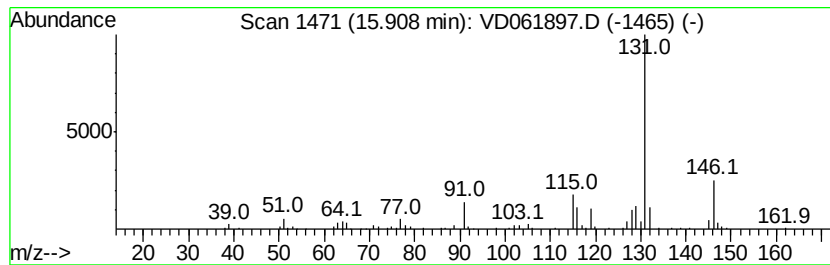
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Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.91	70.61 ug/l	1001410	1,4-Dichlorobenzene-d4	14.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	95
2		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	93
3		Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	90
4		Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	90
5		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041819\
Data File : VD061897.D
Acq On : 19 Apr 2019 5:29
Operator : VA/AP
Sample : K2197-07
Misc : 5.72a/5ml/MSVOA D/SOIL
ALS Vial : 41 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CL-01-041619-B

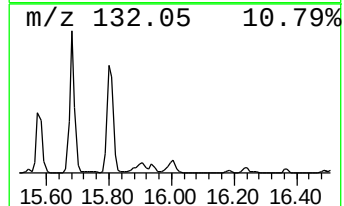
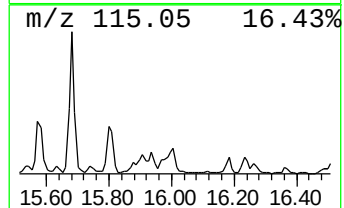
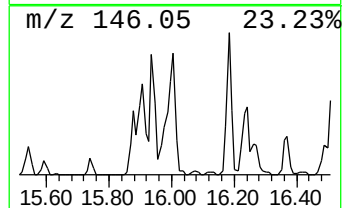
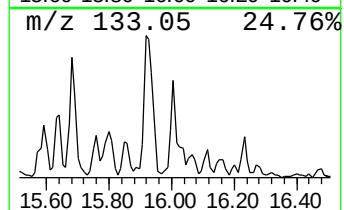
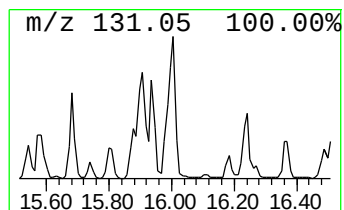
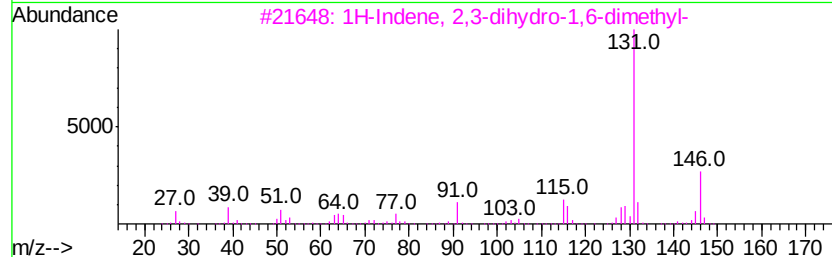
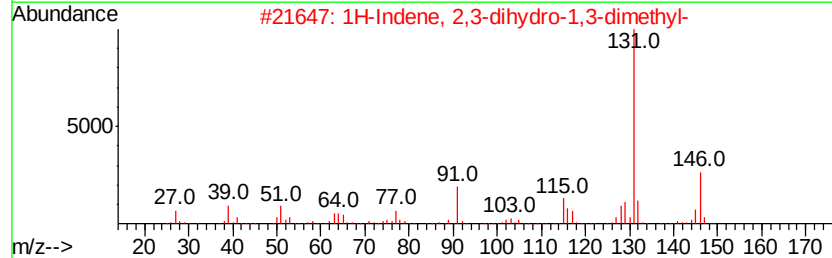
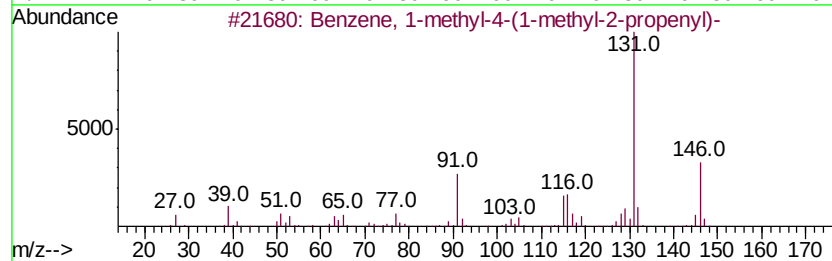
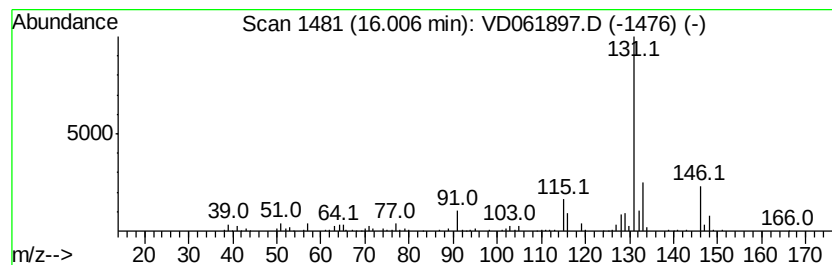
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-methyl-4-(1-meth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.01	75.64 ug/l	1072780	1,4-Dichlorobenzene-d4	14.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	87
2		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	81
3		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	81
4		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	81
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	74



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD041819\
Data File : VD061897.D
Acq On : 19 Apr 2019 5:29
Operator : VA/AP
Sample : K2197-07
Misc : 5.72g/5ml/MSVOA_D/SOIL
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
CL-01-041619-B

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D041519S.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, 2-propenyl-	14.72	71.6	ug/l	1016090	4	14.53	709141	50.0
Benzene, 4-ethyl-...	15.06	55.1	ug/l	781418	4	14.53	709141	50.0
Benzene, 1-etheny...	15.14	89.0	ug/l	1261750	4	14.53	709141	50.0
Benzene, 1,2,3,4-...	15.39	57.2	ug/l	810938	4	14.53	709141	50.0
1H-Indene, 2,3-di...	15.57	101.3	ug/l	1437120	4	14.53	709141	50.0
Dodecane	15.65	54.0	ug/l	765894	4	14.53	709141	50.0
Benzene, 1-etheny...	15.68	157.2	ug/l	2228970	4	14.53	709141	50.0
Naphthalene, 1,2,...	15.80	100.6	ug/l	1426710	4	14.53	709141	50.0
1H-Indene, 2,3-di...	15.91	70.6	ug/l	1001410	4	14.53	709141	50.0
Benzene, 1-methyl...	16.01	75.6	ug/l	1072780	4	14.53	709141	50.0