

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD041219\
 Data File : VD061790.D
 Acq On : 12 Apr 2019 20:53
 Operator : VA/AP
 Sample : K2374-01
 Misc : 6.56g/5ml/MSVOA D/SOIL
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 P-1(0-10)

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	1.538	10	12	20	rVB2	70737	179424	6.34%	0.263%
2	1.637	20	22	31	rVB2	14203	33819	1.20%	0.049%
3	1.814	37	40	45	rVB4	16782	31760	1.12%	0.046%
4	3.202	176	181	188	rBV	30730	84231	2.98%	0.123%
5	3.360	193	197	202	rBV	24766	58813	2.08%	0.086%
6	3.803	237	242	253	rBV	49486	142807	5.05%	0.209%
7	6.126	473	478	489	rVB2	15835	55360	1.96%	0.081%
8	6.904	550	557	564	rBV	324013	920271	32.54%	1.347%
9	7.022	564	569	586	rVB	479433	1413310	49.97%	2.069%
10	7.435	604	611	623	rBV	207979	559757	19.79%	0.819%
11	8.203	683	689	707	rVB	679713	1655621	58.54%	2.423%
12	8.853	748	755	761	rBV2	26378	70982	2.51%	0.104%
13	10.398	905	912	920	rBV	771913	1830311	64.72%	2.679%
14	10.910	960	964	969	rVB3	19632	45968	1.63%	0.067%
15	11.471	1017	1021	1026	rBV	34836	68071	2.41%	0.100%
16	11.629	1031	1037	1038	rBV	19429	39841	1.41%	0.058%
17	11.658	1038	1040	1044	rVV2	22397	49096	1.74%	0.072%
18	11.737	1044	1048	1052	rVV	62027	116923	4.13%	0.171%
19	11.816	1052	1056	1061	rVV	91060	195674	6.92%	0.286%
20	11.885	1061	1063	1069	rVB3	20856	45998	1.63%	0.067%
21	11.993	1069	1074	1077	rBV	13298	29847	1.06%	0.044%
22	12.052	1077	1080	1083	rVV	42488	81345	2.88%	0.119%
23	12.111	1083	1086	1089	rVV	82996	164748	5.83%	0.241%
24	12.160	1089	1091	1096	rVB	56695	95692	3.38%	0.140%
25	12.298	1101	1105	1107	rBV	69100	133882	4.73%	0.196%
26	12.357	1107	1111	1116	rVB	1035798	1670769	59.08%	2.445%
27	12.446	1116	1120	1125	rVB4	23818	60530	2.14%	0.089%
28	12.534	1125	1129	1133	rBV2	49219	101316	3.58%	0.148%
29	12.623	1133	1138	1142	rBV3	78306	228063	8.06%	0.334%
30	12.692	1142	1145	1148	rVV	176500	292290	10.34%	0.428%
31	12.741	1148	1150	1157	rVB	144561	246105	8.70%	0.360%
32	12.840	1157	1160	1165	rBV2	50106	95711	3.38%	0.140%
33	12.987	1171	1175	1179	rBV2	242800	535941	18.95%	0.784%
34	13.037	1179	1180	1182	rVV2	70139	88412	3.13%	0.129%

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35	13.086	1182	1185	1188	rVV	125302	236378	8.36%	0.346%
36	13.145	1188	1191	1192	rVV3	45441	70339	2.49%	0.103%
37	13.214	1192	1198	1201	rVV2	536506	951513	33.64%	1.393%
38	13.302	1201	1207	1214	rVV3	469358	1350402	47.75%	1.977%
39	13.391	1214	1216	1220	rVV2	121966	200896	7.10%	0.294%
40	13.470	1220	1224	1227	rVV3	190041	479998	16.97%	0.703%
41	13.548	1227	1232	1235	rVV2	1002660	1601146	56.61%	2.344%
42	13.608	1235	1238	1241	rVV2	98663	210758	7.45%	0.308%
43	13.657	1241	1243	1245	rVV2	127462	216136	7.64%	0.316%
44	13.706	1245	1248	1252	rVV3	470485	820927	29.03%	1.202%
45	13.765	1252	1254	1256	rVV	69656	103966	3.68%	0.152%
46	13.814	1256	1259	1263	rVV4	180480	511519	18.09%	0.749%
47	13.883	1263	1266	1267	rVV2	274463	482946	17.08%	0.707%
48	13.923	1267	1270	1272	rVV2	533706	1039164	36.74%	1.521%
49	13.962	1272	1274	1277	rVV	322321	525855	18.59%	0.770%
50	14.021	1277	1280	1282	rVV3	121901	238372	8.43%	0.349%
51	14.080	1282	1286	1287	rVV2	215640	496329	17.55%	0.726%
52	14.110	1287	1289	1292	rVV2	297798	558521	19.75%	0.817%
53	14.159	1292	1294	1295	rVV	201294	310984	11.00%	0.455%
54	14.198	1295	1298	1299	rVV	504293	843038	29.81%	1.234%
55	14.228	1299	1301	1308	rVV	932479	1896372	67.05%	2.776%
56	14.316	1308	1310	1311	rVV	185542	248639	8.79%	0.364%
57	14.366	1311	1315	1319	rVV5	413669	1180180	41.73%	1.727%
58	14.434	1319	1322	1324	rVV	689315	1174683	41.54%	1.719%
59	14.513	1327	1330	1332	rVV	1332366	1908619	67.49%	2.794%
60	14.553	1332	1334	1337	rVV	484124	817056	28.89%	1.196%
61	14.592	1337	1338	1339	rVV	139292	150515	5.32%	0.220%
62	14.641	1339	1343	1345	rVV3	299135	716066	25.32%	1.048%
63	14.700	1346	1349	1350	rVV3	343865	557385	19.71%	0.816%
64	14.730	1350	1352	1354	rVV	411731	574427	20.31%	0.841%
65	14.769	1354	1356	1358	rVV	647720	1098387	38.84%	1.608%
66	14.799	1358	1359	1362	rVV3	748219	1126351	39.83%	1.649%
67	14.848	1362	1364	1366	rVV	306502	577621	20.42%	0.845%
68	14.907	1366	1370	1372	rVV2	563570	1361886	48.15%	1.993%
69	14.937	1372	1373	1375	rVV2	413474	598546	21.16%	0.876%
70	14.976	1375	1377	1378	rVV	583861	954742	33.76%	1.397%
71	14.996	1378	1379	1383	rVV3	851554	1973841	69.79%	2.889%

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72	15.045	1383	1384	1387	rVV	896032	1272727	45.00%	1.863%
73	15.084	1387	1388	1390	rVV	320723	445725	15.76%	0.652%
74	15.143	1390	1394	1398	rVV3	785817	2349312	83.07%	3.439%
75	15.212	1398	1401	1404	rVV4	580875	1364472	48.25%	1.997%
76	15.271	1404	1407	1409	rVV2	700321	1555251	54.99%	2.276%
77	15.311	1409	1411	1413	rVV2	797586	1495868	52.89%	2.189%
78	15.340	1413	1414	1415	rVV	761045	697025	24.65%	1.020%
79	15.370	1415	1417	1420	rVV	1013876	1604956	56.75%	2.349%
80	15.419	1420	1422	1425	rVV2	896683	1595115	56.40%	2.335%
81	15.468	1425	1427	1430	rVV3	479082	1224593	43.30%	1.792%
82	15.527	1430	1433	1436	rVV3	446601	1212195	42.86%	1.774%
83	15.576	1436	1438	1441	rVV3	376169	999910	35.36%	1.464%
84	15.626	1441	1443	1445	rVV	616333	951900	33.66%	1.393%
85	15.665	1445	1447	1452	rVV2	1389515	2828152	100.00%	4.139%
86	15.734	1452	1454	1457	rVV2	678018	1357765	48.01%	1.987%
87	15.783	1457	1459	1464	rVV4	374698	1188602	42.03%	1.740%
88	15.931	1464	1474	1475	rVV2	574650	2324833	82.20%	3.403%
89	15.990	1475	1480	1482	rVV3	565188	1417115	50.11%	2.074%
90	16.069	1482	1488	1489	rVV3	279891	861598	30.47%	1.261%
91	16.098	1489	1491	1493	rVV	190692	336772	11.91%	0.493%
92	16.137	1493	1495	1500	rVV3	263616	589597	20.85%	0.863%
93	16.216	1500	1503	1506	rVV2	199939	450317	15.92%	0.659%
94	16.256	1506	1507	1509	rVV	111426	131411	4.65%	0.192%
95	16.285	1509	1510	1515	rVV4	82139	157064	5.55%	0.230%
96	16.344	1515	1516	1519	rVV2	58317	88912	3.14%	0.130%
97	16.403	1519	1522	1525	rVB4	48447	110400	3.90%	0.162%
98	16.472	1525	1529	1531	rBV4	49315	87833	3.11%	0.129%
99	16.728	1551	1555	1558	rVB3	21558	35588	1.26%	0.052%

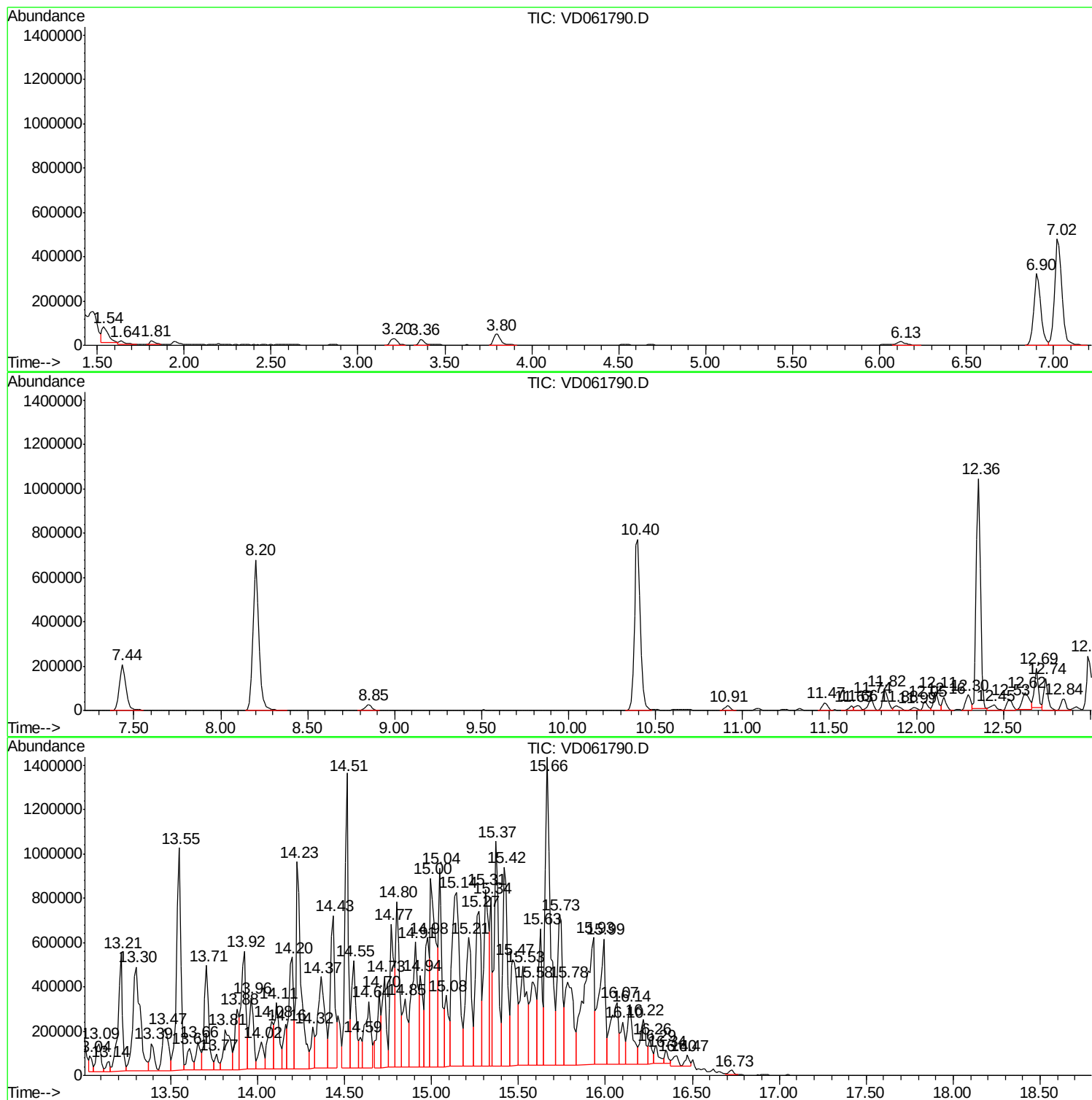
Sum of corrected areas: 68322199

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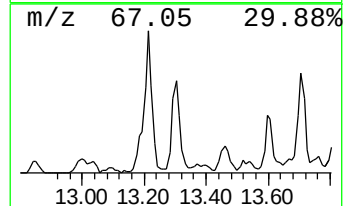
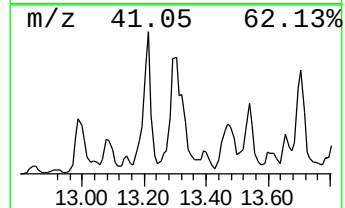
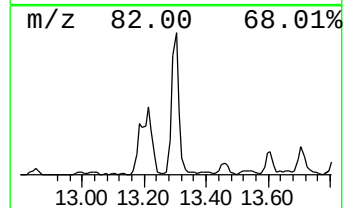
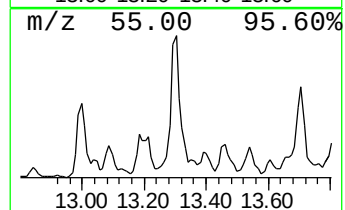
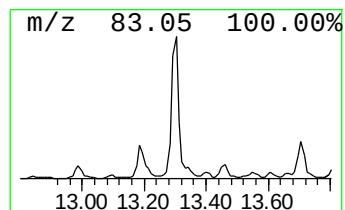
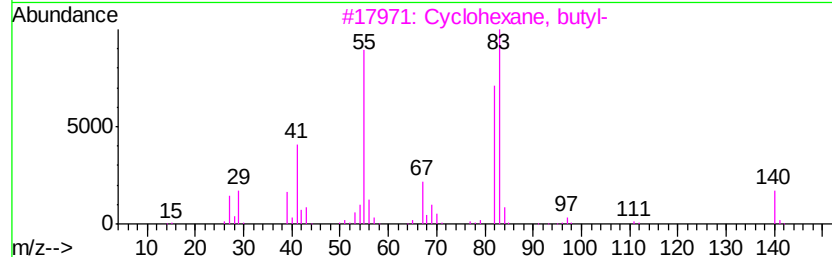
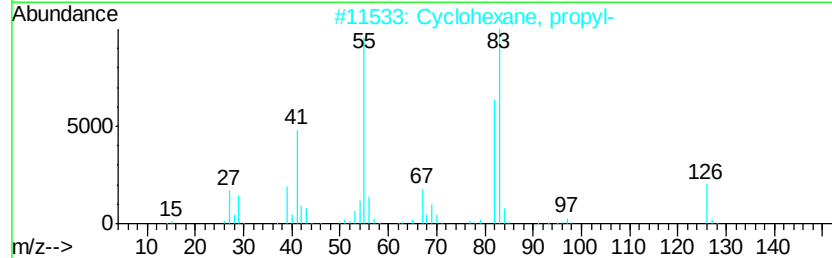
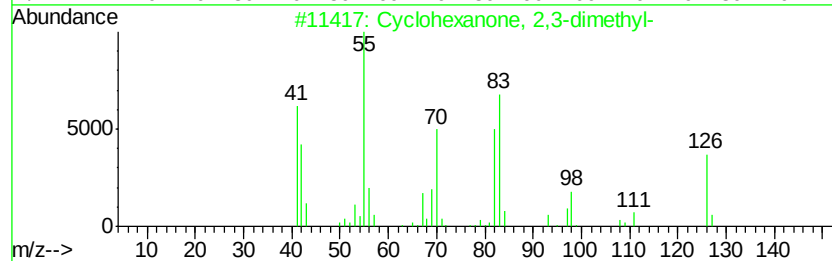
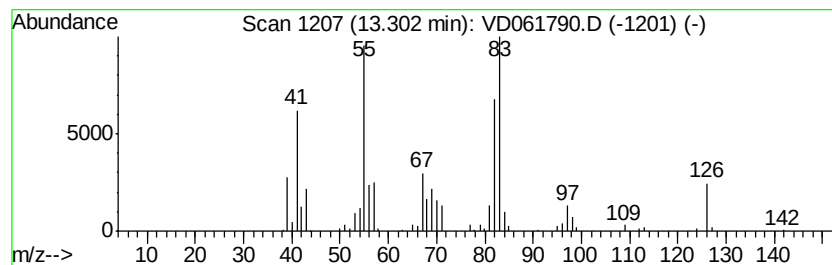
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D040619S.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclohexanone, 2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.30	40.41 ug/l	1350400	Chlorobenzene-d5	12.36		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexanone, 2,3-dimethyl-	126	C8H14O	013395-76-1	87
2		Cyclohexane, propyl-	126	C9H18	001678-92-8	74
3		Cyclohexane, butyl-	140	C10H20	001678-93-9	64
4		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	62
5		Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	58



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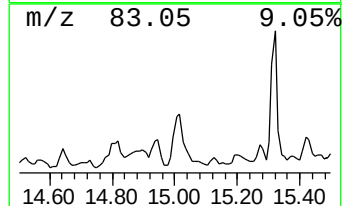
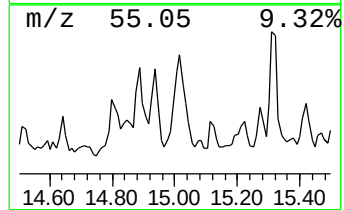
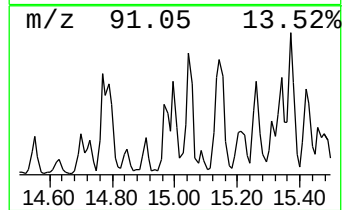
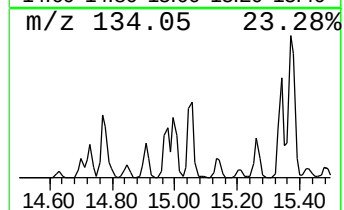
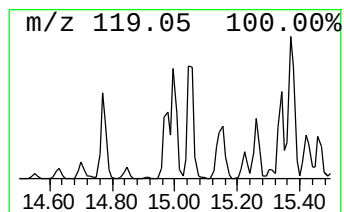
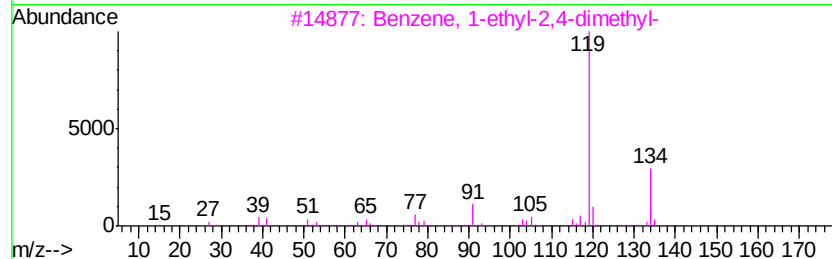
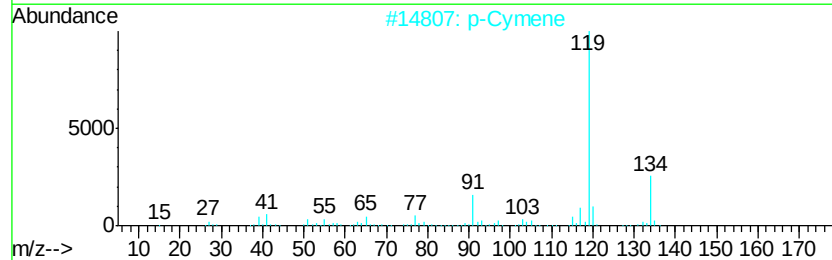
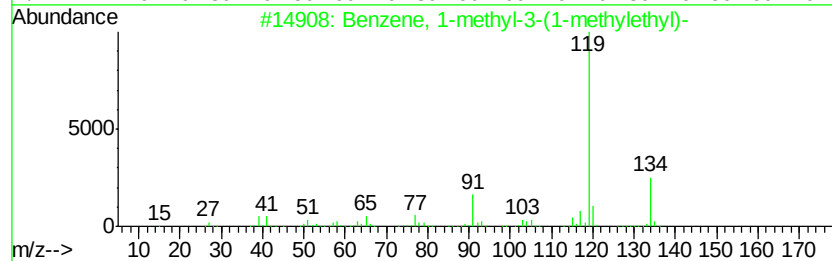
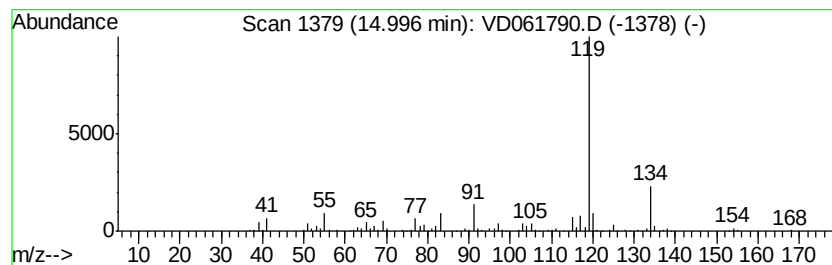
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1-methyl-3-(1-meth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.00	51.71 ug/l	1973840	1,4-Dichlorobenzene-d4	14.51
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1-methyl-3-(1-methyleth...	134 C10H14	000535-77-3	94
2	p-Cymene	134 C10H14	000099-87-6	94
3	Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9	93
4	o-Cymene	134 C10H14	000527-84-4	91
5	Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4	87



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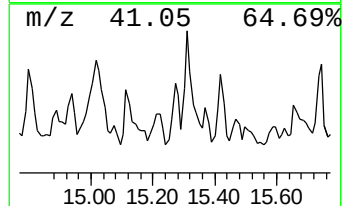
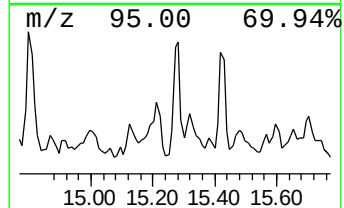
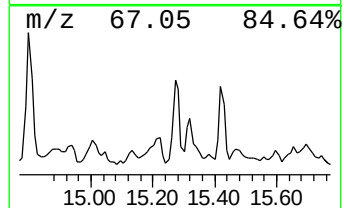
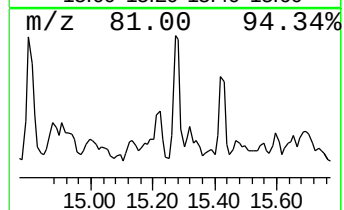
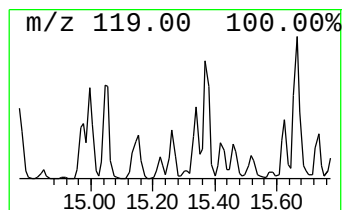
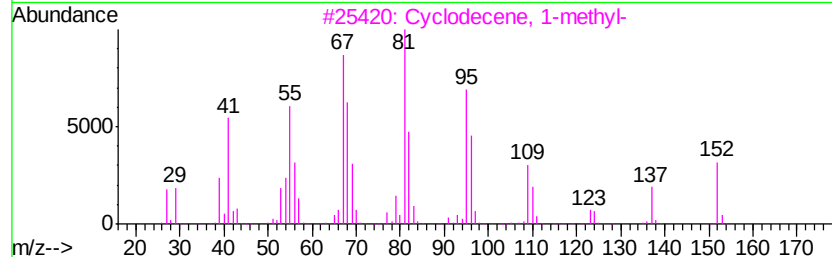
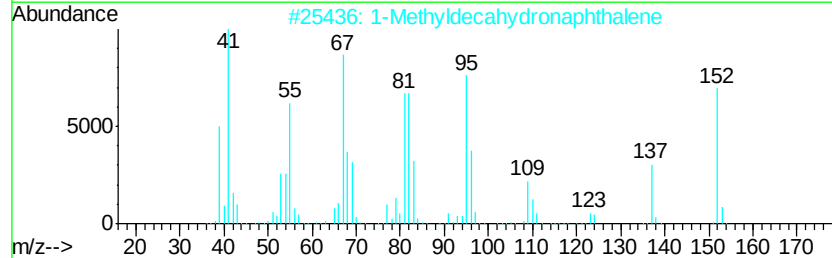
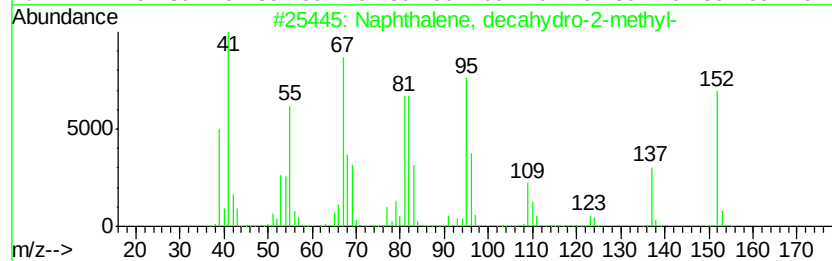
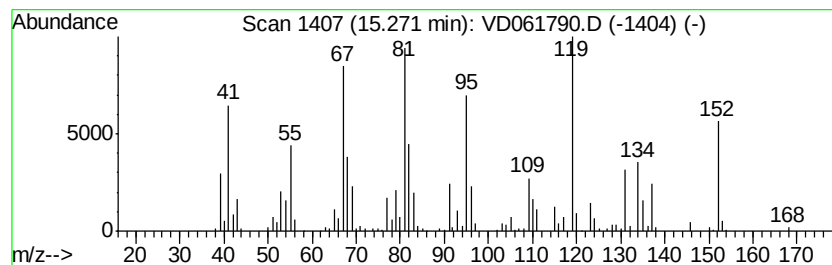
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Quant Title : SW846 8260

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

Peak Number 4 Naphthalene, decahydro-2-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.27	40.74 ug/l	1555250	1,4-Dichlorobenzene-d4	14.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	80
2		1-Methyldecahydronaphthalene	152	C11H20	002958-75-0	80
3		Cyclodecene, 1-methyl-	152	C11H20	066633-38-3	64
4		Benzofuran, octahydro-6-methyl-3...	152	C10H16O	077954-13-3	62
5		Decalin, syn-1-methyl-, cis-	152	C11H20	1000158-89-1	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041219\
Data File : VD061790.D
Acq On : 12 Apr 2019 20:53
Operator : VA/AP
Sample : K2374-01
Misc : 6.56g/5ml/MSVOA D/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
P-1(0-10)

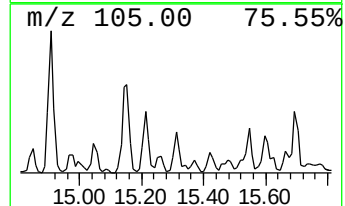
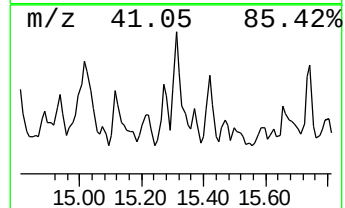
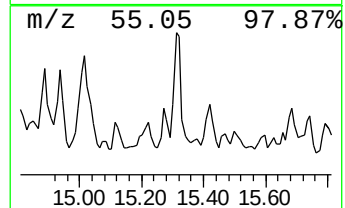
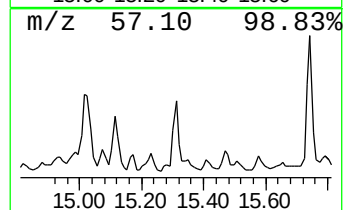
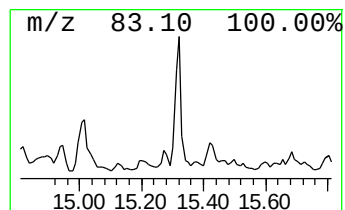
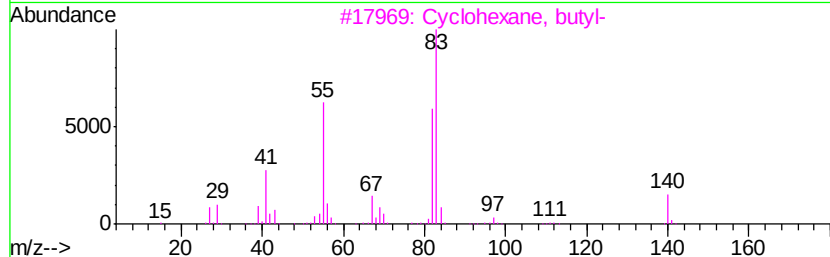
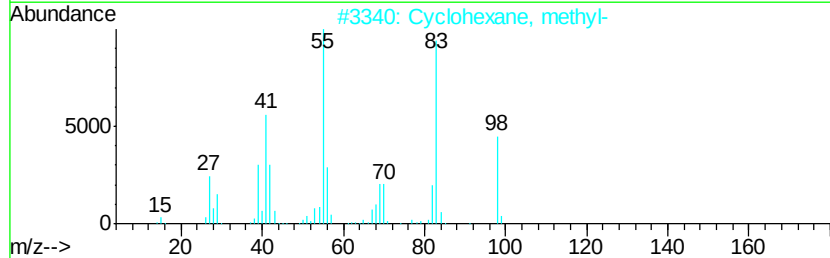
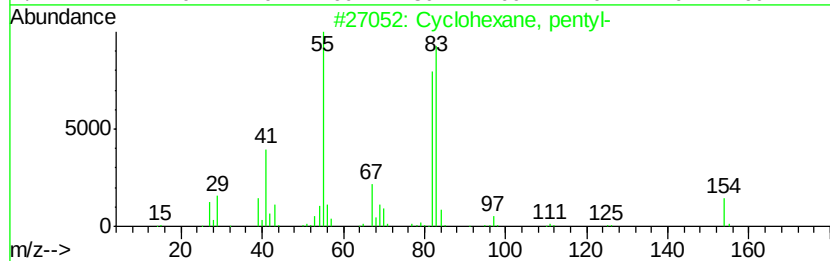
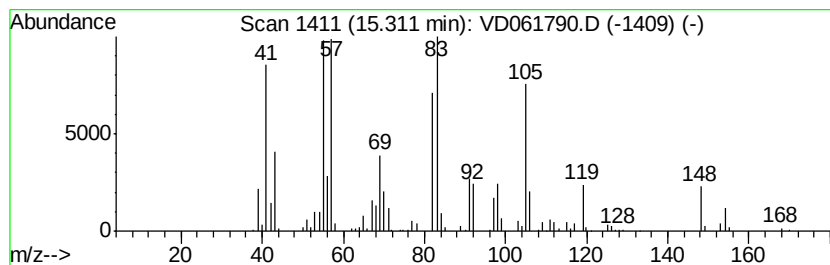
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Quant Title : SW846 8260

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

Peak Number 5 unknown15.31 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.31	39.19 ug/l	1495870	1,4-Dichlorobenzene-d4	14.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, pentyl-	154	C11H22	004292-92-6	45
2		Cyclohexane, methyl-	98	C7H14	000108-87-2	38
3		Cyclohexane, butyl-	140	C10H20	001678-93-9	35
4		Cyclohexanone, 3,3,5,5-tetramethyl-	154	C10H18O	014376-79-5	30
5		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	27



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041219\
Data File : VD061790.D
Acq On : 12 Apr 2019 20:53
Operator : VA/AP
Sample : K2374-01
Misc : 6.56g/5ml/MSVOA D/SOIL
ALS Vial : 22 Sample Multiplier: 1

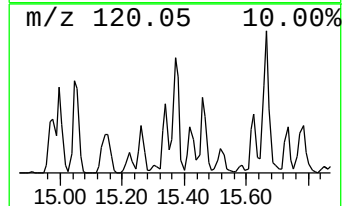
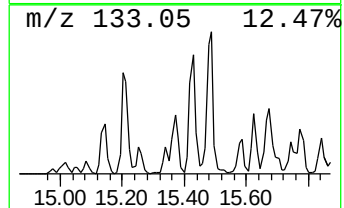
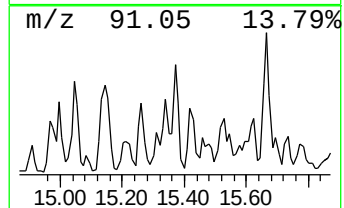
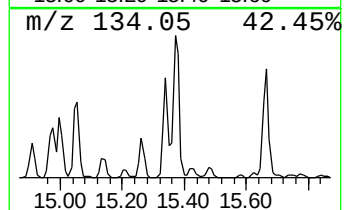
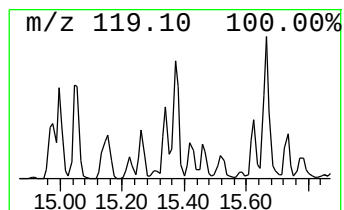
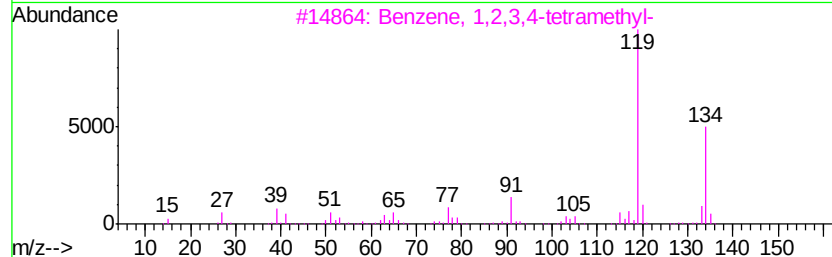
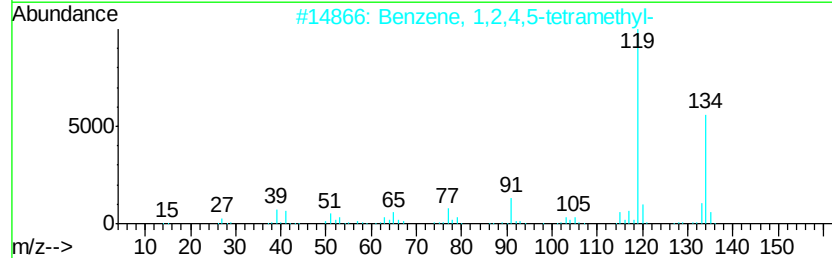
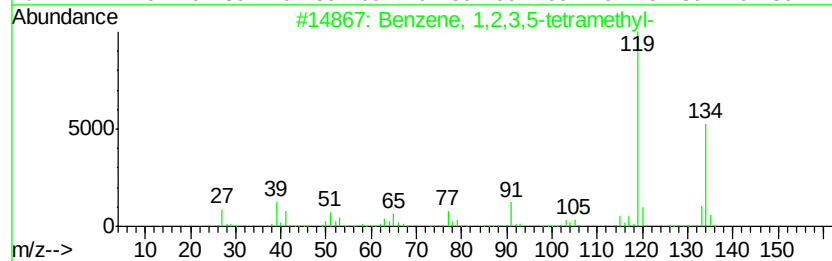
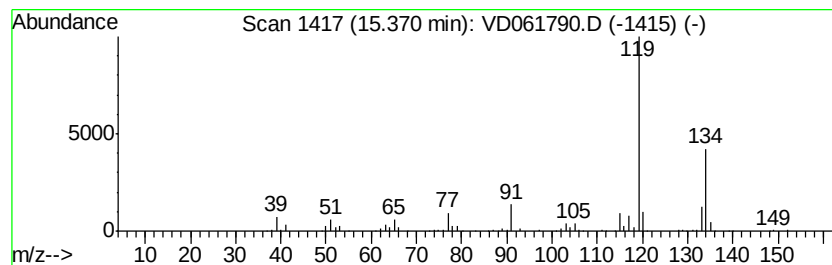
Instrument :
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ClientSampled :
P-1(0-10)

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, 1,2,3,5-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.37	42.05 ug/l	1604960	1,4-Dichlorobenzene-d4	14.51
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1,2,3,5-tetramethyl-	134 C10H14	000527-53-7	95
2	Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2	95
3	Benzene, 1,2,3,4-tetramethyl-	134 C10H14	000488-23-3	95
4	o-Cymene	134 C10H14	000527-84-4	94
5	Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041219\
Data File : VD061790.D
Acq On : 12 Apr 2019 20:53
Operator : VA/AP
Sample : K2374-01
Misc : 6.56g/5ml/MSVOA D/SOIL
ALS Vial : 22 Sample Multiplier: 1

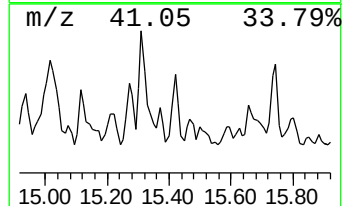
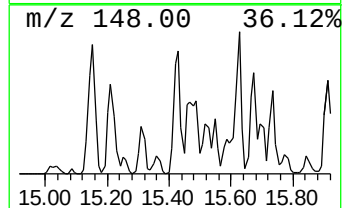
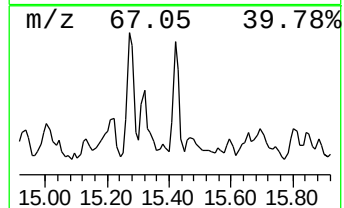
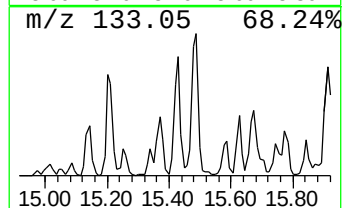
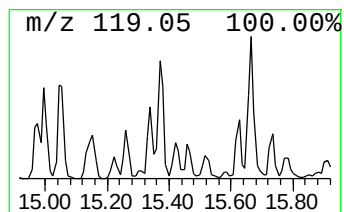
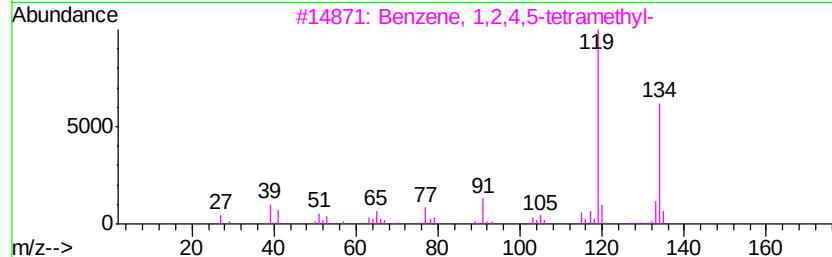
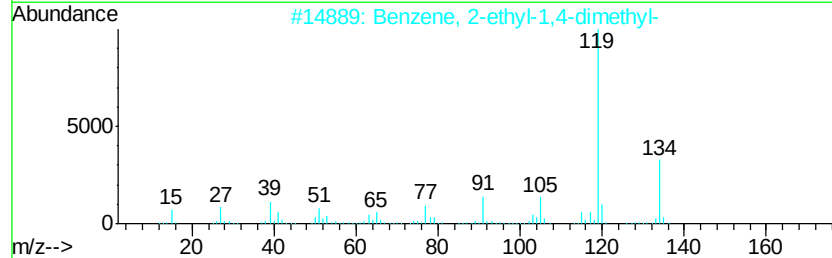
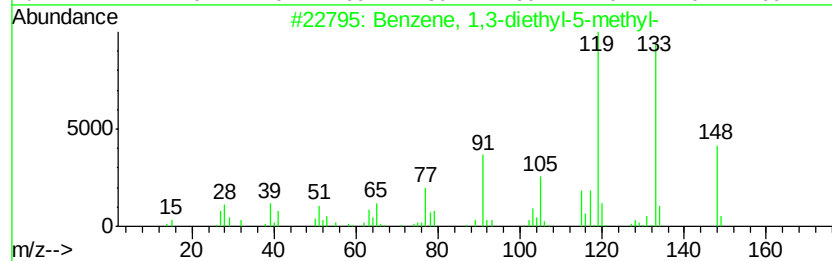
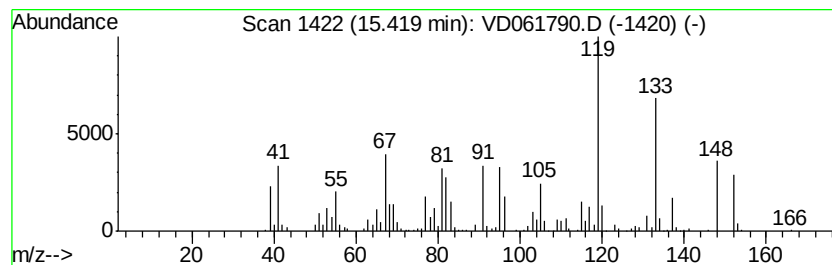
Instrument :
MSVOA_D
ClientSampleId :
P-1(0-10)

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 7 Benzene, 1,3-diethyl-5-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.42	41.79 ug/l	1595120	1,4-Dichlorobenzene-d4	14.51	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	90
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	46
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	42
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	42
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	42



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041219\
Data File : VD061790.D
Acq On : 12 Apr 2019 20:53
Operator : VA/AP
Sample : K2374-01
Misc : 6.56g/5ml/MSVOA D/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
P-1(0-10)

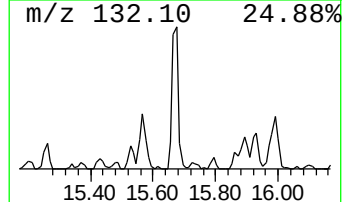
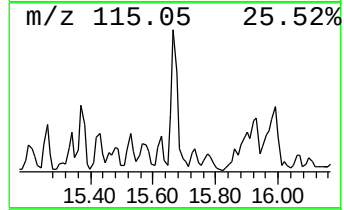
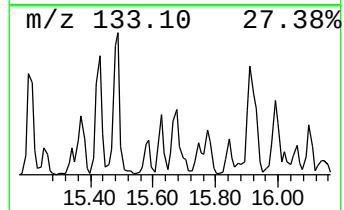
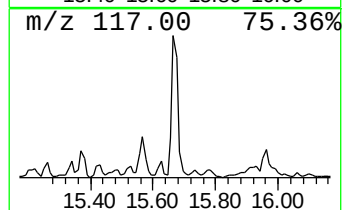
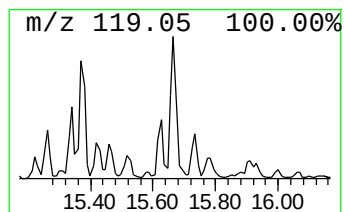
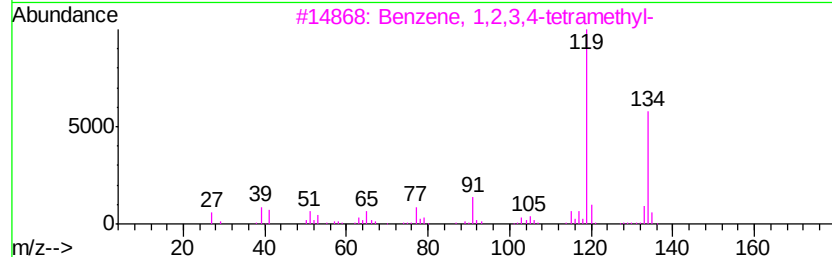
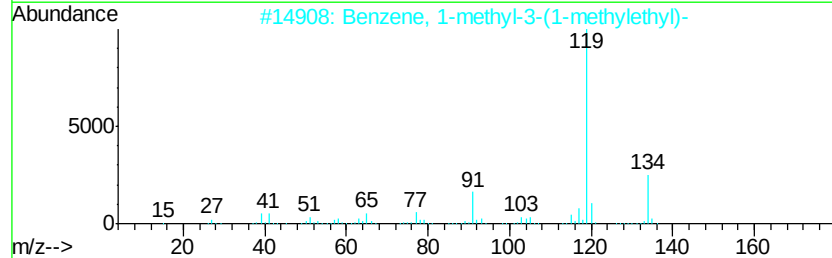
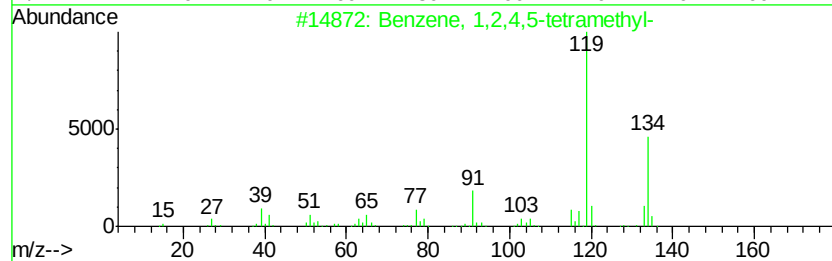
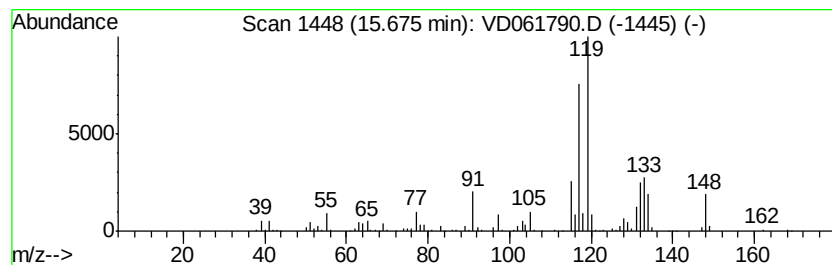
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.67	74.09 ug/l	2828150	1,4-Dichlorobenzene-d4	14.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
2		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	64
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	62
4		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	62
5		o-Cymene	134	C10H14	000527-84-4	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041219\
Data File : VD061790.D
Acq On : 12 Apr 2019 20:53
Operator : VA/AP
Sample : K2374-01
Misc : 6.56g/5ml/MSVOA D/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
P-1(0-10)

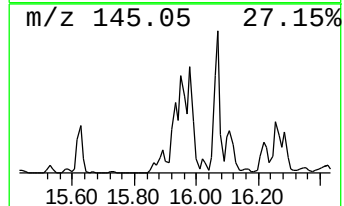
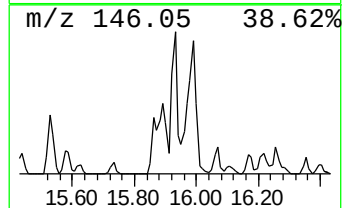
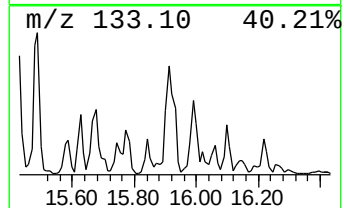
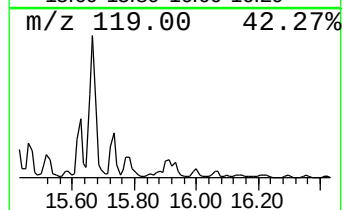
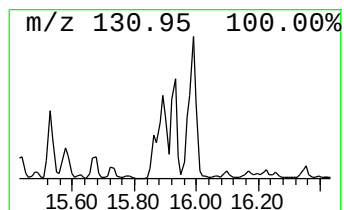
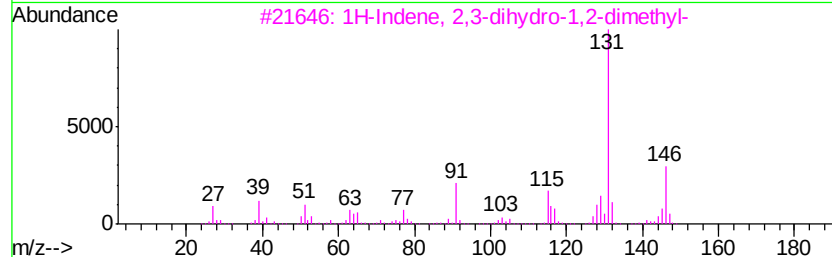
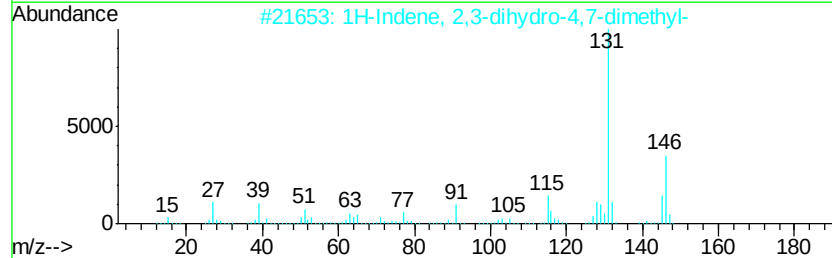
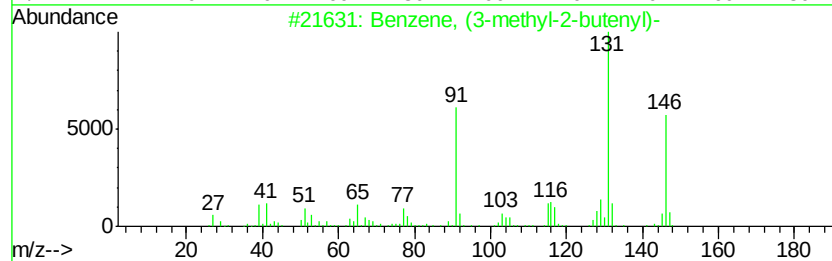
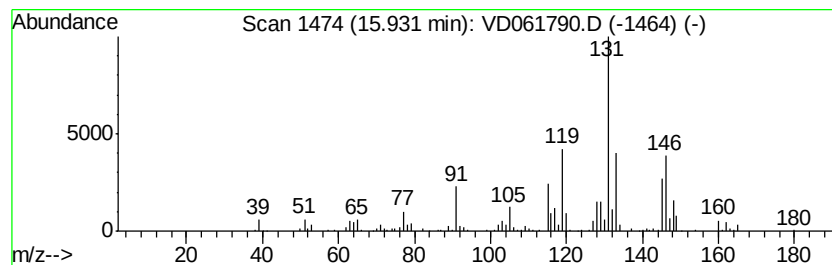
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, (3-methyl-2-butenyl)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.93	60.90 ug/l	2324830	1,4-Dichlorobenzene-d4	14.51

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	90
2		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	60
3		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	60
4		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	60
5		Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD041219\
Data File : VD061790.D
Acq On : 12 Apr 2019 20:53
Operator : VA/AP
Sample : K2374-01
Misc : 6.56g/5ml/MSVOA D/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
P-1(0-10)

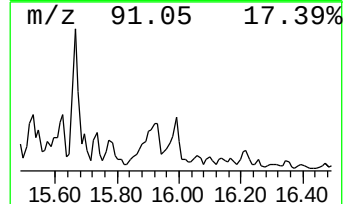
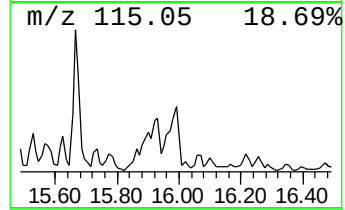
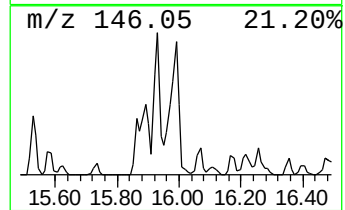
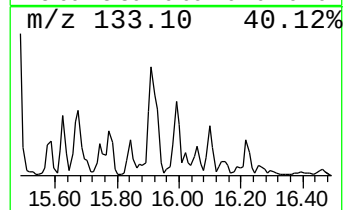
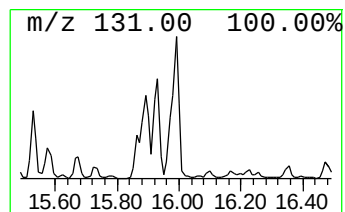
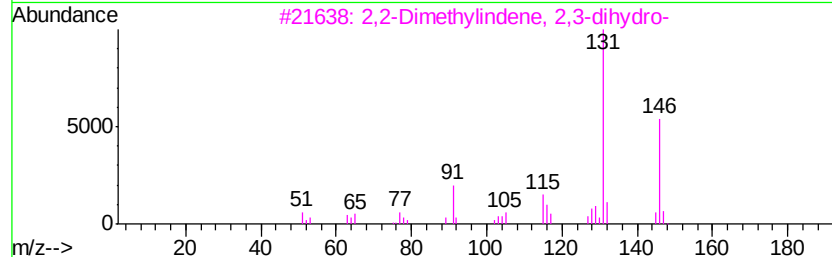
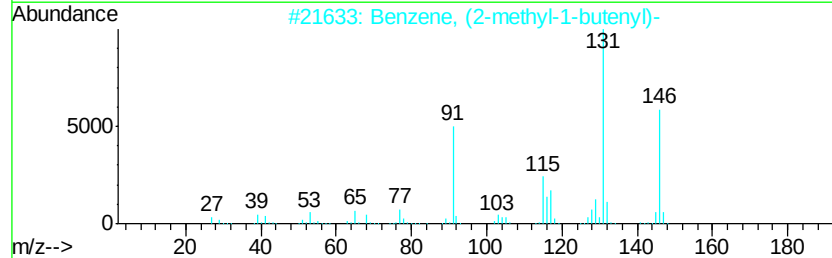
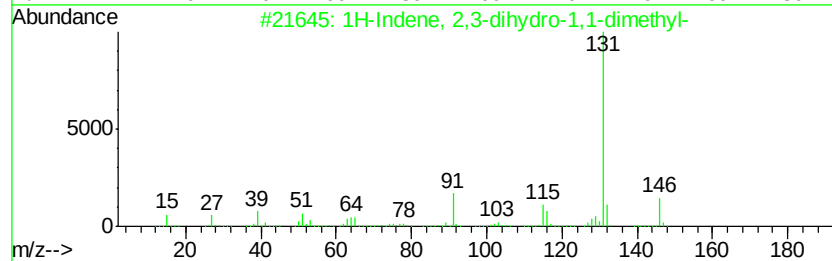
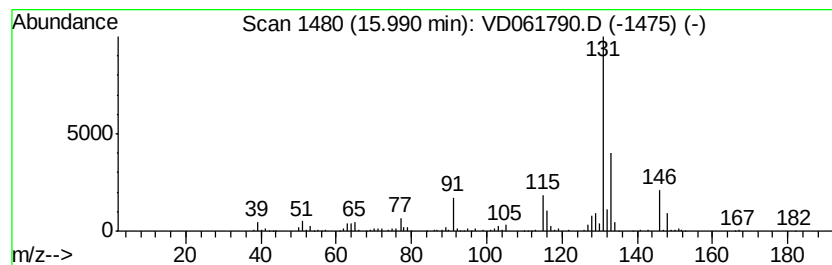
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.99	37.12 ug/l	1417120	1,4-Dichlorobenzene-d4	14.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	70
2		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	70
3		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	70
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	70
5		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD041219\
 Data File : VD061790.D
 Acq On : 12 Apr 2019 20:53
 Operator : VA/AP
 Sample : K2374-01
 Misc : 6.56g/5ml/MSVOA_D/SOIL
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 P-1(0-10)

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
Cyclohexanone, 2,...	13.30	40.4	ug/l	1350400	3	12.36	1670770	50.0
Benzene, 1-methyl...	15.00	51.7	ug/l	1973840	4	14.51	1908620	50.0
Naphthalene, deca...	15.27	40.7	ug/l	1555250	4	14.51	1908620	50.0
unknown15.31	15.31	39.2	ug/l	1495870	4	14.51	1908620	50.0
Benzene, 1,2,3,5-...	15.37	42.0	ug/l	1604960	4	14.51	1908620	50.0
Benzene, 1,3-diet...	15.42	41.8	ug/l	1595120	4	14.51	1908620	50.0
Benzene, 1,2,4,5-...	15.67	74.1	ug/l	2828150	4	14.51	1908620	50.0
Benzene, (3-methy...	15.93	60.9	ug/l	2324830	4	14.51	1908620	50.0
1H-Indene, 2,3-di...	15.99	37.1	ug/l	1417120	4	14.51	1908620	50.0