

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD081919\
 Data File : VD063659.D
 Acq On : 19 Aug 2019 15:56
 Operator : JC/SY
 Sample : K4385-14
 Misc : 4.75a/5ml/MSVOA D/SOIL
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 T8-C-(0)

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\82D080919S.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	1.474	3	5	25	rVB	1067896	2718624	100.00%	12.084%
2	1.818	37	40	46	rVB	17724	46102	1.70%	0.205%
3	3.216	176	182	191	rVB	87328	241654	8.89%	1.074%
4	3.807	237	242	249	rBV	51928	145996	5.37%	0.649%
5	6.061	466	471	477	rBV2	10842	30935	1.14%	0.138%
6	6.455	507	511	517	rBV5	10283	29542	1.09%	0.131%
7	6.918	552	558	565	rBV2	548176	1599883	58.85%	7.111%
8	7.036	565	570	584	rVB	687680	2064849	75.95%	9.178%
9	7.449	606	612	622	rBV	421243	1116890	41.08%	4.964%
10	8.217	683	690	702	rBV	1038795	2605439	95.84%	11.581%
11	10.403	905	912	919	rBV	860711	1998261	73.50%	8.882%
12	10.501	919	922	928	rVB3	24093	60868	2.24%	0.271%
13	10.924	959	965	971	rVB4	18082	41677	1.53%	0.185%
14	11.486	1017	1022	1026	rBV2	21933	42358	1.56%	0.188%
15	11.633	1033	1037	1039	rBV2	13507	30255	1.11%	0.134%
16	11.673	1039	1041	1045	rVV2	13656	31899	1.17%	0.142%
17	11.751	1045	1049	1053	rVB3	23819	46281	1.70%	0.206%
18	11.830	1053	1057	1061	rBV2	55121	115124	4.23%	0.512%
19	11.909	1061	1065	1068	rVV3	15160	37363	1.37%	0.166%
20	11.988	1068	1073	1077	rVV2	11596	27673	1.02%	0.123%
21	12.066	1077	1081	1083	rVV2	23797	44045	1.62%	0.196%
22	12.116	1083	1086	1089	rVV2	42007	82747	3.04%	0.368%
23	12.165	1089	1091	1096	rVB3	21444	40291	1.48%	0.179%
24	12.303	1101	1105	1108	rBV	32832	68510	2.52%	0.305%
25	12.362	1108	1111	1117	rVB	1035869	1655472	60.89%	7.358%
26	12.450	1117	1120	1125	rVB3	13004	35153	1.29%	0.156%
27	12.539	1125	1129	1134	rBV3	27522	75923	2.79%	0.337%
28	12.647	1134	1140	1143	rBV5	41033	141352	5.20%	0.628%
29	12.696	1143	1145	1148	rVB	53535	75652	2.78%	0.336%
30	12.746	1148	1150	1157	rVB2	46596	94056	3.46%	0.418%
31	12.854	1157	1161	1165	rBV3	29353	54576	2.01%	0.243%
32	12.992	1171	1175	1178	rBV3	98135	206356	7.59%	0.917%
33	13.041	1178	1180	1183	rVV3	37354	60402	2.22%	0.268%
34	13.090	1183	1185	1189	rVB3	39791	76751	2.82%	0.341%

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35	13.218	1192	1198	1201	rBV2	149982	282314	10.38%	1.255%
36	13.307	1201	1207	1208	rBV4	104191	227815	8.38%	1.013%
37	13.326	1208	1209	1214	rVV2	142270	222753	8.19%	0.990%
38	13.395	1214	1216	1220	rVB4	38738	78423	2.88%	0.349%
39	13.474	1220	1224	1227	rBV4	100310	224976	8.28%	1.000%
40	13.553	1229	1232	1236	rVV	718327	1000855	36.81%	4.449%
41	13.612	1236	1238	1240	rVV2	34287	44104	1.62%	0.196%
42	13.661	1240	1243	1245	rVV3	35359	61704	2.27%	0.274%
43	13.710	1245	1248	1255	rVB2	156055	267223	9.83%	1.188%
44	13.819	1255	1259	1263	rBV5	63236	157225	5.78%	0.699%
45	13.888	1263	1266	1268	rVV3	73893	136357	5.02%	0.606%
46	13.966	1271	1274	1277	rVV2	92305	156511	5.76%	0.696%
47	14.025	1277	1280	1282	rVV4	36344	49830	1.83%	0.221%
48	14.114	1285	1289	1292	rBV4	45929	121913	4.48%	0.542%
49	14.163	1292	1294	1295	rVV2	37323	52402	1.93%	0.233%
50	14.193	1295	1297	1300	rVV	82568	140825	5.18%	0.626%
51	14.242	1300	1302	1306	rVB3	45741	103034	3.79%	0.458%
52	14.360	1311	1314	1319	rVB7	45366	120724	4.44%	0.537%
53	14.439	1319	1322	1327	rBV4	34867	103299	3.80%	0.459%
54	14.518	1327	1330	1333	rBV	596127	730478	26.87%	3.247%
55	14.646	1341	1343	1347	rVB2	35518	41091	1.51%	0.183%
56	14.695	1347	1348	1354	rVB5	21456	49018	1.80%	0.218%
57	14.803	1357	1359	1364	rBV6	47303	86928	3.20%	0.386%
58	14.882	1364	1367	1369	rBV3	37904	49427	1.82%	0.220%
59	14.941	1369	1373	1377	rVB4	36796	87042	3.20%	0.387%
60	15.010	1377	1380	1383	rBV4	25327	58208	2.14%	0.259%
61	15.138	1389	1393	1397	rBV6	55573	125213	4.61%	0.557%
62	15.217	1397	1401	1404	rVB6	21194	55955	2.06%	0.249%
63	15.286	1404	1408	1410	rBV2	104856	162084	5.96%	0.720%
64	15.394	1416	1419	1421	rBV	57087	92743	3.41%	0.412%
65	15.423	1421	1422	1424	rVV	39277	54081	1.99%	0.240%
66	15.463	1424	1426	1429	rVB4	36471	69905	2.57%	0.311%
67	15.660	1442	1446	1448	rBV4	38763	102757	3.78%	0.457%
68	15.699	1448	1450	1453	rVB3	47010	73184	2.69%	0.325%
69	15.748	1453	1455	1458	rBV3	25346	37562	1.38%	0.167%
70	15.807	1458	1461	1466	rBV4	108653	181448	6.67%	0.807%
71	15.876	1466	1468	1471	rVB	59095	76772	2.82%	0.341%

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Sampling : 1 Min Area: 3 % of largest Peak
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72	15.925	1471	1473	1474	rBV	45587	50591	1.86%	0.225%
73	16.044	1482	1485	1488	rVV4	30746	50508	1.86%	0.225%
74	16.103	1488	1491	1493	rVV4	17521	30651	1.13%	0.136%
75	16.142	1493	1495	1498	rVV2	101760	168299	6.19%	0.748%
76	16.201	1499	1501	1509	rVB8	67352	165603	6.09%	0.736%
77	16.378	1516	1519	1525	rVB7	62580	201524	7.41%	0.896%
78	16.487	1525	1530	1534	rBV8	54622	163064	6.00%	0.725%
79	16.575	1537	1539	1542	rBV3	26641	52396	1.93%	0.233%
80	17.038	1584	1586	1589	rVB3	42845	61950	2.28%	0.275%
81	17.126	1593	1595	1601	rVB6	25064	58114	2.14%	0.258%
82	17.254	1605	1608	1611	rVB5	38429	65844	2.42%	0.293%

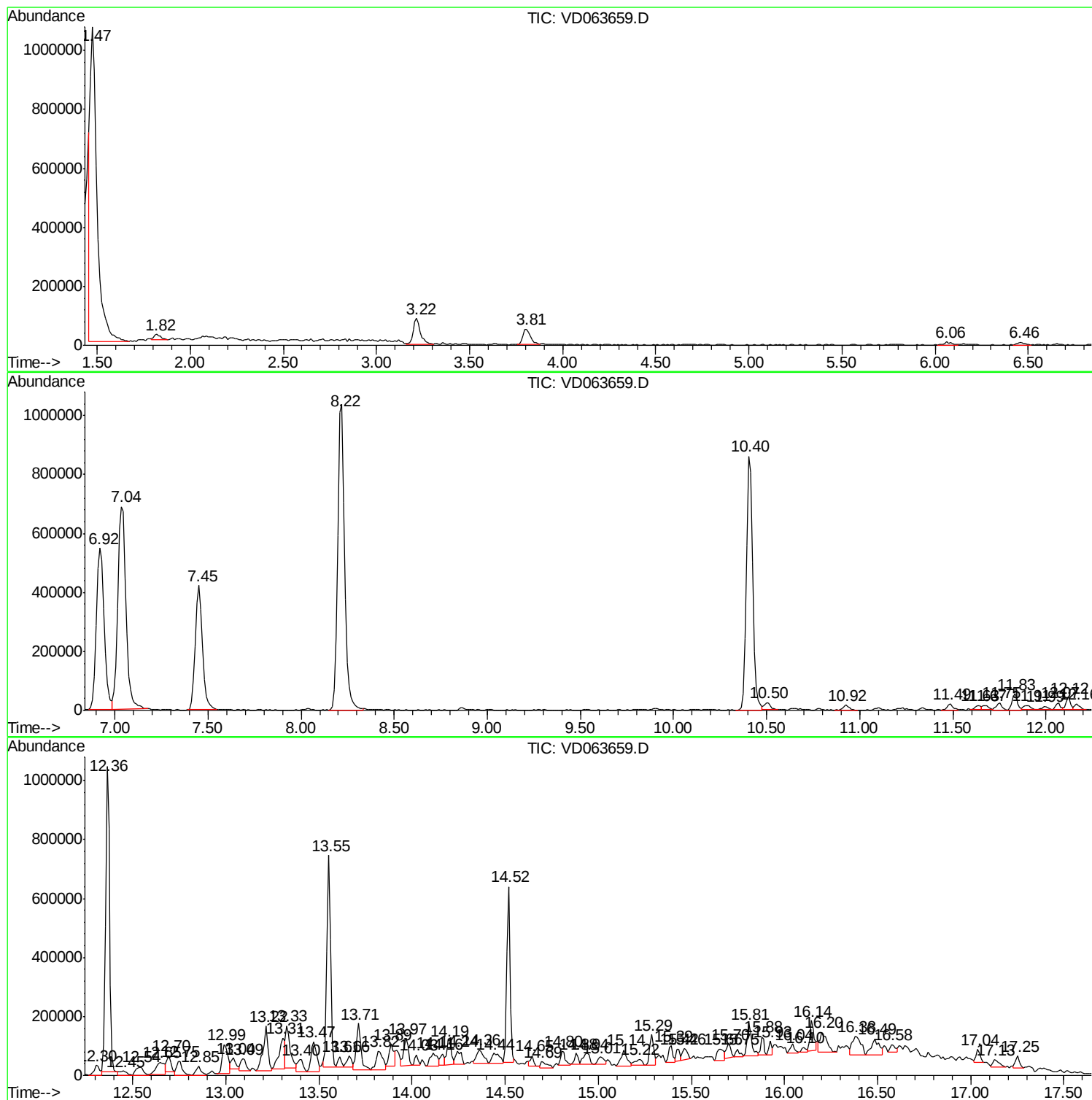
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080919S.M
Quant Title : SW846 8260

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



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Instrument :
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T8-C-(0)

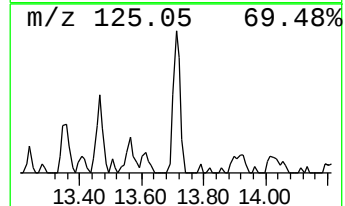
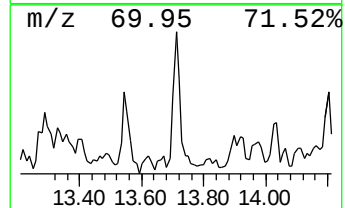
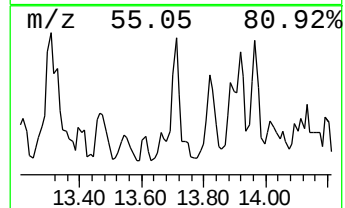
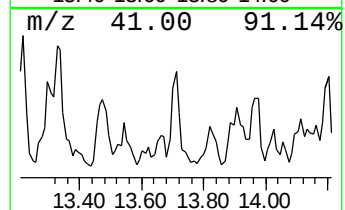
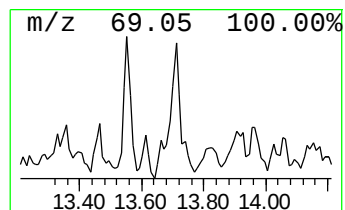
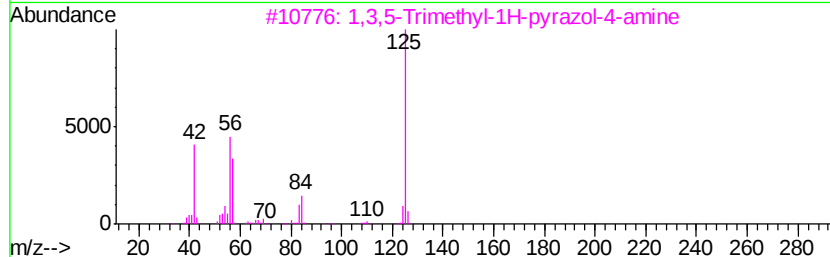
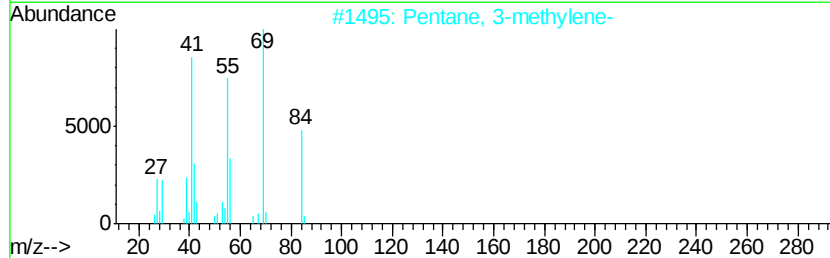
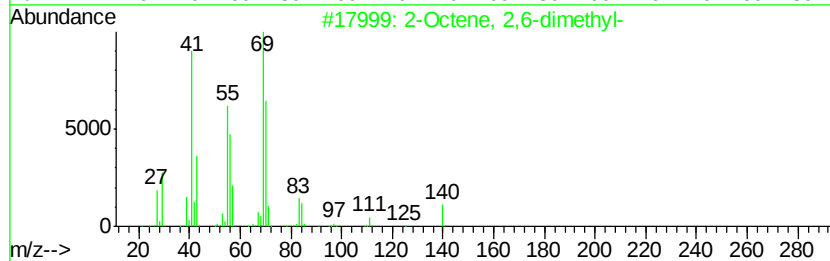
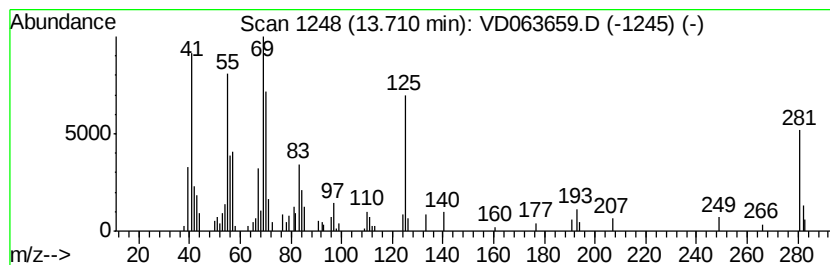
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080919S.M
Quant Title : SW846 8260

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

Peak Number 3 unknown13.71 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.71	18.29 ug/l	267223	1,4-Dichlorobenzene-d4	14.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	46
2		Pentane, 3-methylene-	84	C6H12	000760-21-4	43
3		1,3,5-Trimethyl-1H-pyrazol-4-amine	125	C6H11N3	028466-21-9	38
4		2-Pentene, 3-methyl-, (E)-	84	C6H12	000616-12-6	38
5		Thieno[2,3-d]-1,3-thiaselenol-2-...	238	C5H2S3Se	081803-09-0	30



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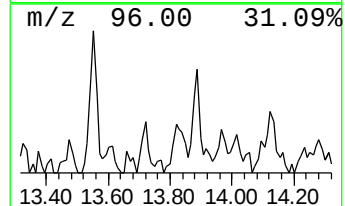
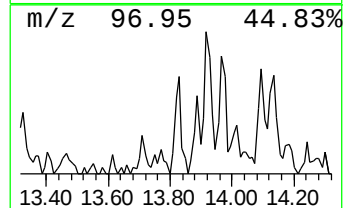
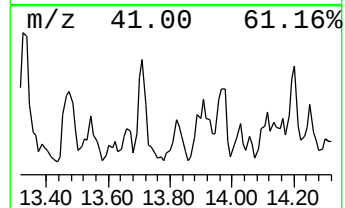
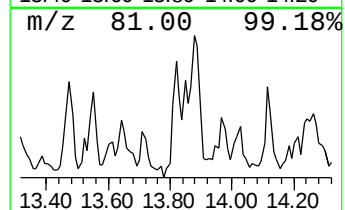
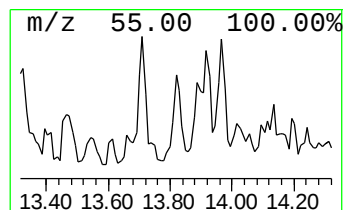
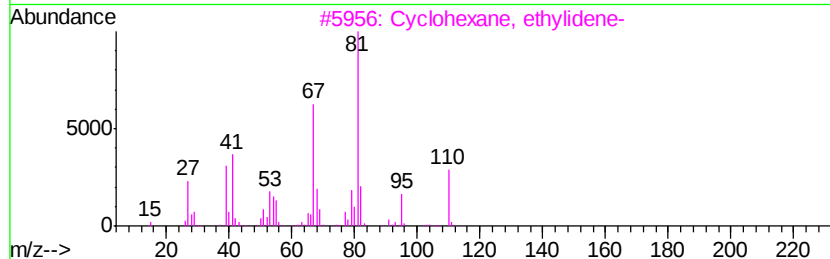
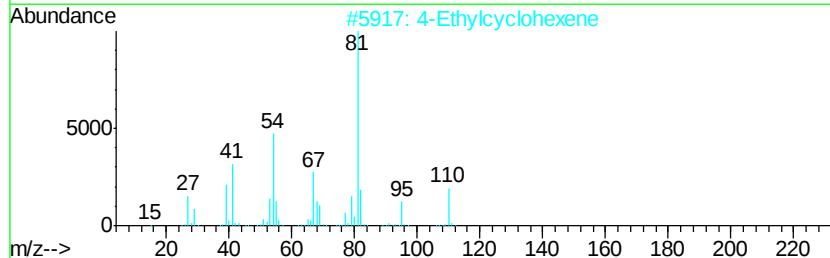
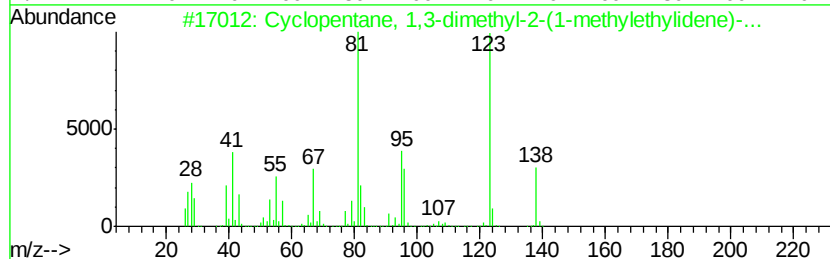
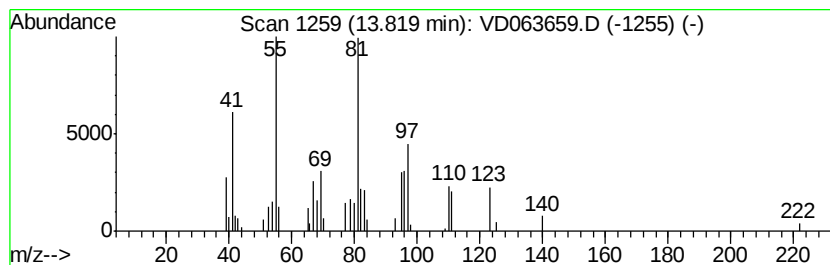
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown13.82 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.82	10.76 ug/l	157225	1,4-Dichlorobenzene-d4	14.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, 1,3-dimethyl-2-(1-...	138	C10H18	061142-30-1	47
2		4-Ethylcyclohexene	110	C8H14	003742-42-5	43
3		Cyclohexane, ethylidene-	110	C8H14	001003-64-1	41
4		Cyclohexene, 1-ethyl-	110	C8H14	001453-24-3	38
5		2-Propenyl-3-vinylloxirane	110	C7H10O	070597-49-8	38



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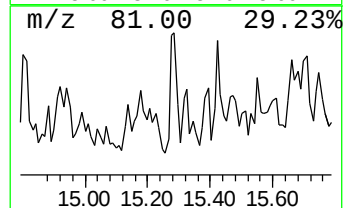
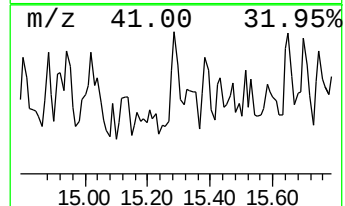
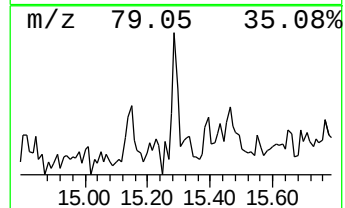
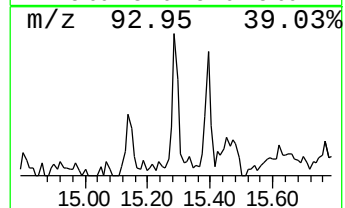
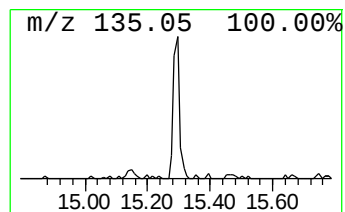
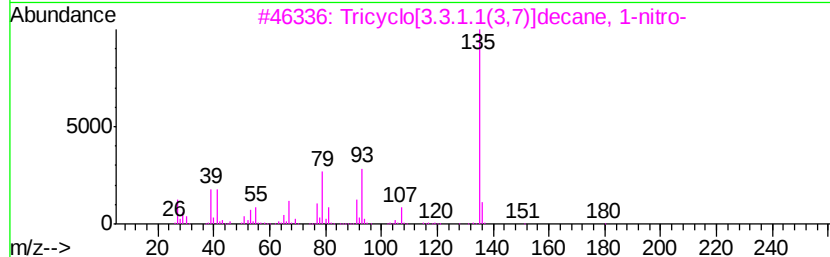
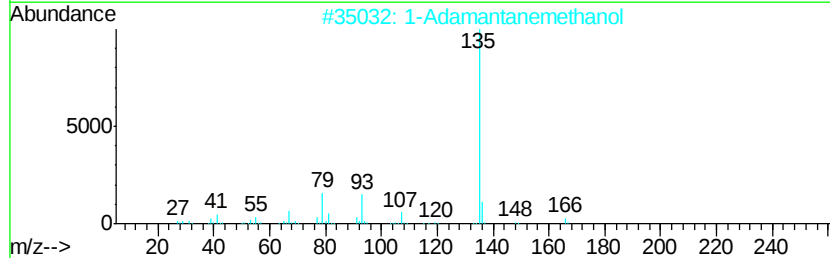
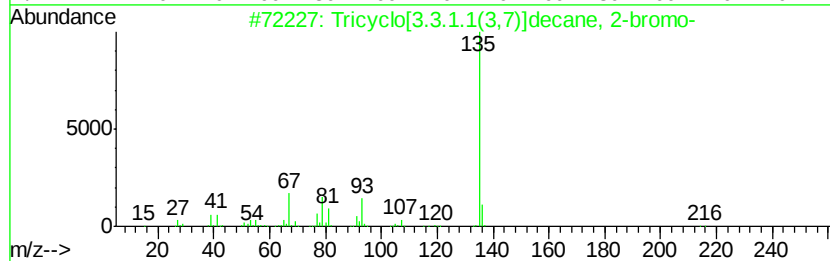
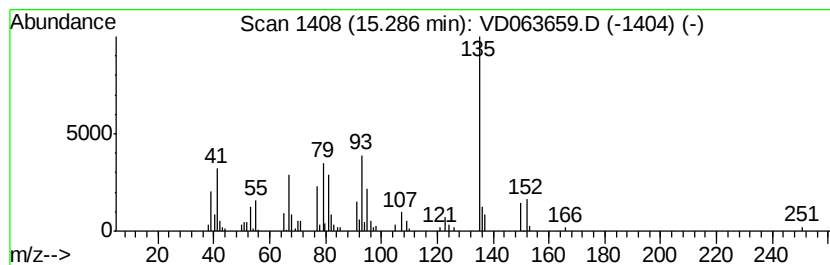
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Quant Title : SW846 8260

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

Peak Number 5 Tricyclo[3.3.1.1(3,7)]decan... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.29	11.09 ug/l	162084	1,4-Dichlorobenzene-d4	14.52

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tricyclo[3.3.1.1(3,7)]decane, 2-...	214	C10H15Br	007314-85-4	64
2		1-Adamantanemethanol	166	C11H18O	000770-71-8	62
3		Tricyclo[3.3.1.1(3,7)]decane, 1-...	181	C10H15NO2	007575-82-8	59
4		N-[2-(Trifluoromethyl)phenyl]-1-...	323	C18H20F3NO	314055-04-4	53
5		Phosphonic dichloride, 1-adamantyl-	252	C10H15Cl2OP	1000294-15-7	53



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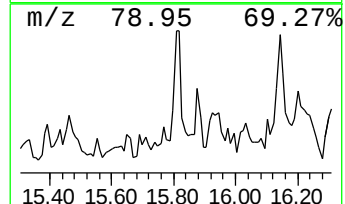
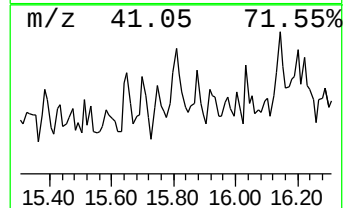
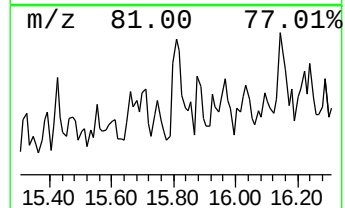
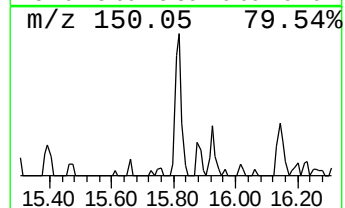
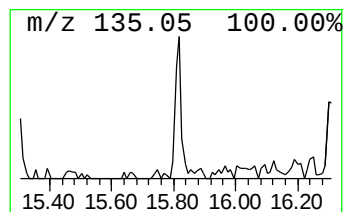
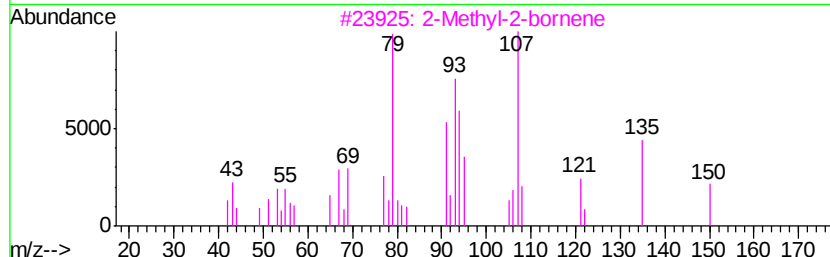
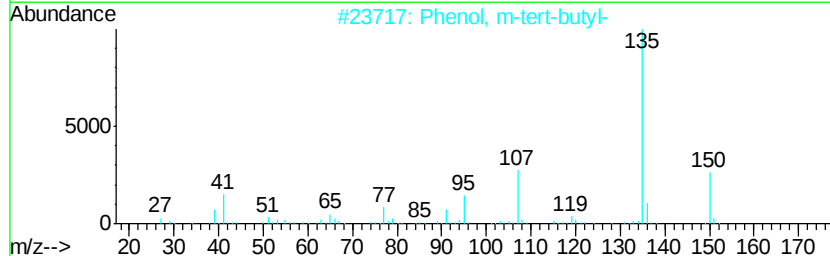
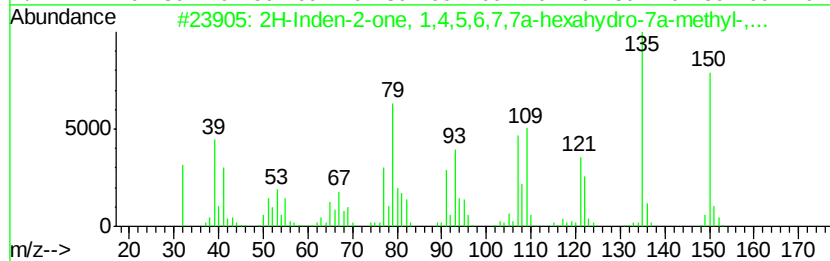
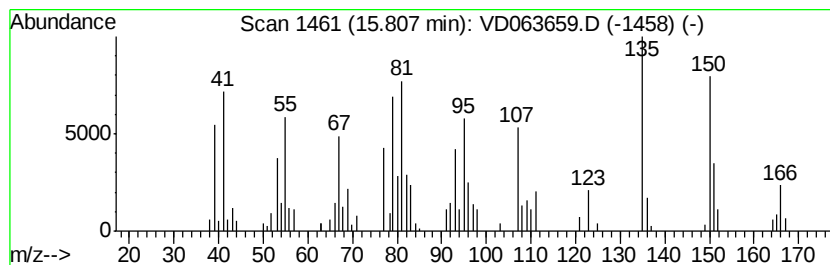
Instrument :
MSVOA_D
ClientSampleId :
T8-C-(0)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080919S.M
Quant Title : SW846 8260

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

Peak Number 6 2H-Inden-2-one, 1,4,5,6,7,7... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.81	12.42 ug/l	181448	1,4-Dichlorobenzene-d4	14.52	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Inden-2-one, 1,4,5,6,7,7a-hex...	150	C10H14O	054725-16-5	64
2	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	46
3	2-Methyl-2-bornene	150	C11H18	072540-93-3	45
4	Bicyclo[3.1.1]hept-3-en-2-one, 4...	150	C10H14O	001196-01-6	45
5	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD081919\
Data File : VD063659.D
Acq On : 19 Aug 2019 15:56
Operator : JC/SY
Sample : K4385-14
Misc : 4.75g/5ml/MSVOA D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
T8-C-(0)

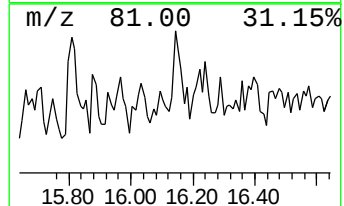
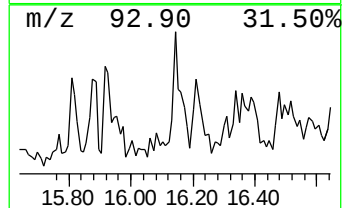
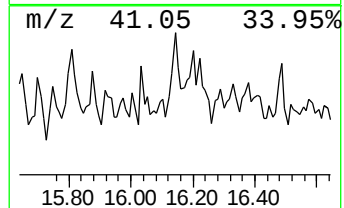
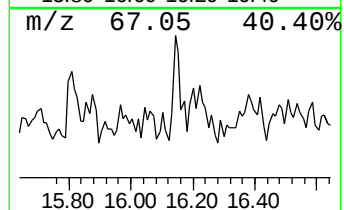
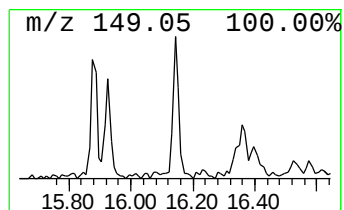
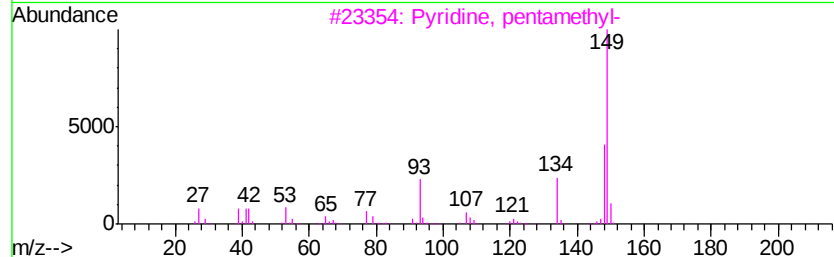
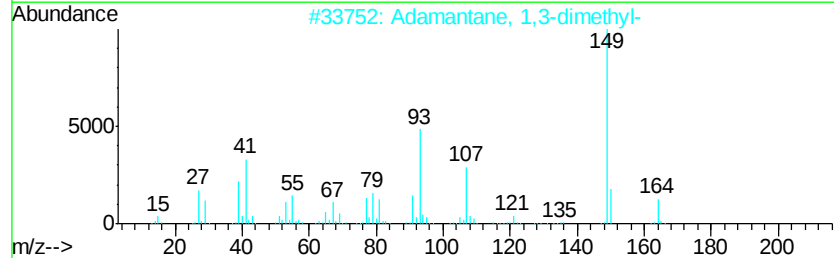
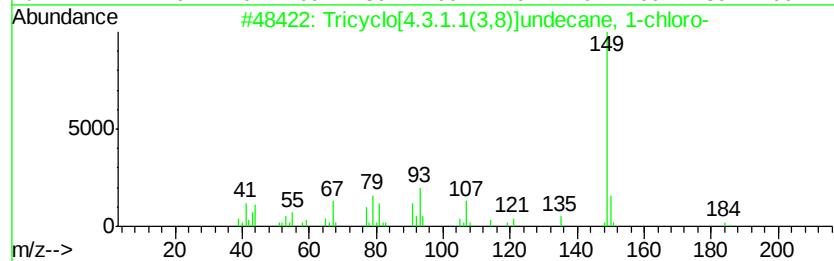
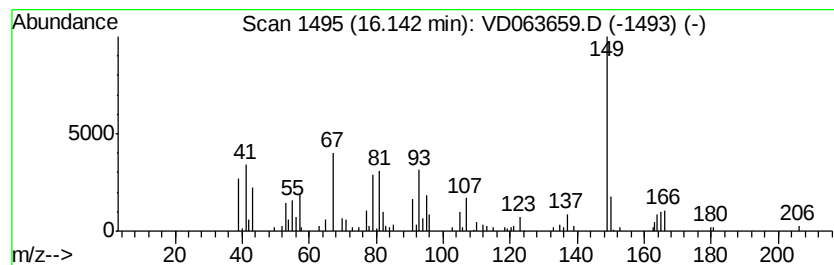
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080919S.M
Quant Title : SW846 8260

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

Peak Number 7 unknown16.14 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.14	11.52 ug/l	168299	1,4-Dichlorobenzene-d4	14.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricyclo[4.3.1.1(3,8)]undecane, ...	184	C11H17Cl	027011-46-7	40
2		Adamantane, 1,3-dimethyl-	164	C12H20	000702-79-4	38
3		Pyridine, pentamethyl-	149	C10H15N	003748-83-2	37
4		Phthalic acid, isobutyl 2-methyl...	276	C16H20O4	1000315-82-3	35
5		Phthalic acid, isobutyl 7-methyl...	344	C21H28O4	1000315-43-4	32



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD081919\
Data File : VD063659.D
Acq On : 19 Aug 2019 15:56
Operator : JC/SY
Sample : K4385-14
Misc : 4.75a/5ml/MSVOA D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleID :
T8-C-(0)

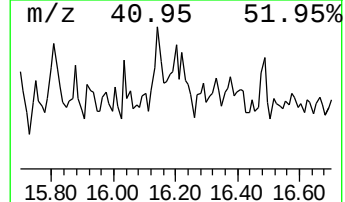
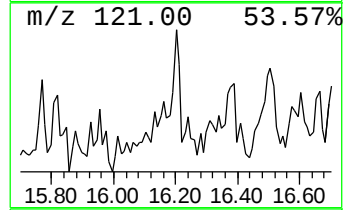
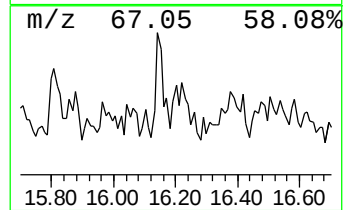
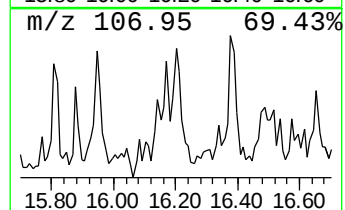
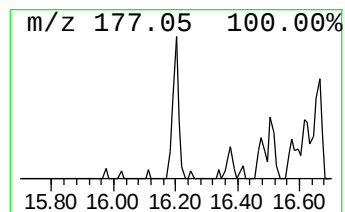
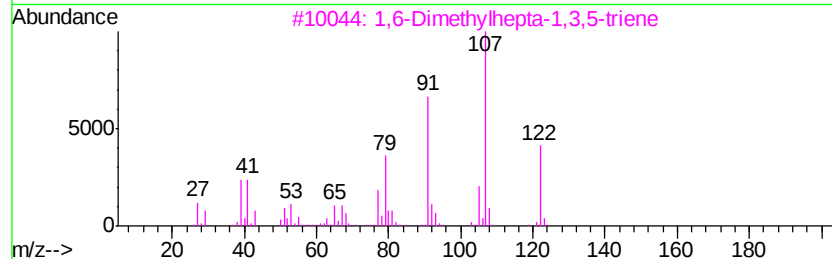
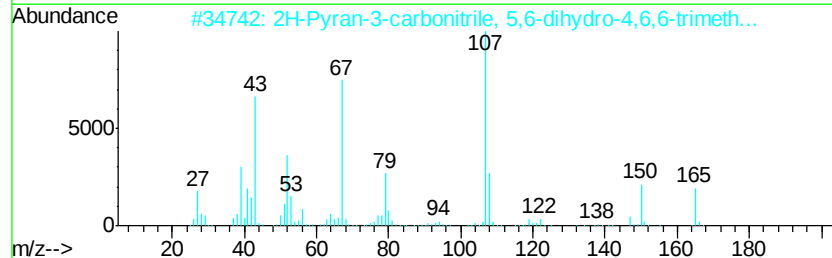
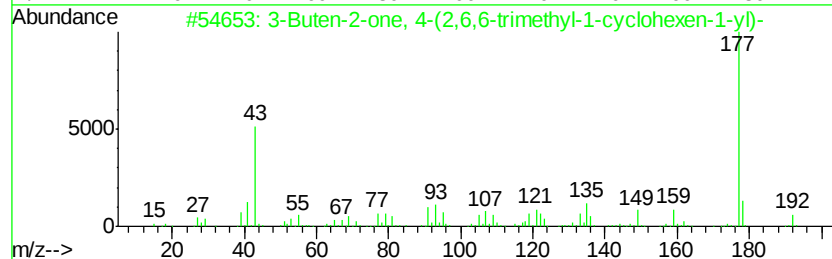
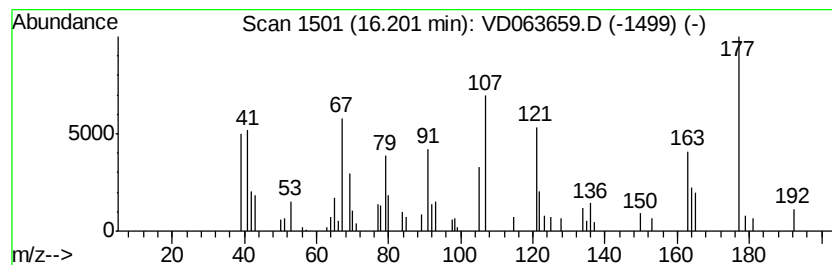
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Quant Title : SW846 8260

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

Peak Number 8 unknown16.20 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.20	11.34 ug/l	165603	1,4-Dichlorobenzene-d4	14.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Buten-2-one, 4-(2,6,6-trimethy...	192	C13H20O	014901-07-6	27
2		2H-Pyran-3-carbonitrile, 5,6-dih...	165	C9H11NO2	069245-73-4	22
3		1,6-Dimethylhepta-1,3,5-triene	122	C9H14	1000196-61-0	22
4		3,3-Dimethyl-6-methylenecyclohexene	122	C9H14	020185-16-4	18
5		5-Acetylpyrimidine	122	C6H6N2O	010325-70-9	18



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD081919\
Data File : VD063659.D
Acq On : 19 Aug 2019 15:56
Operator : JC/SY
Sample : K4385-14
Misc : 4.75g/5ml/MSVOA D/SOIL
ALS Vial : 20 Sample Multiplier: 1

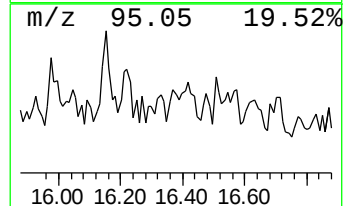
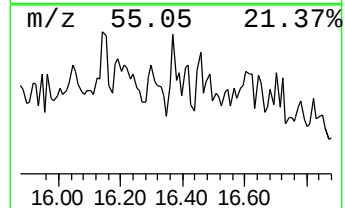
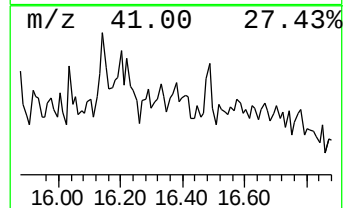
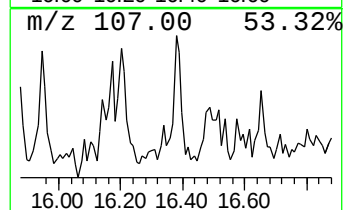
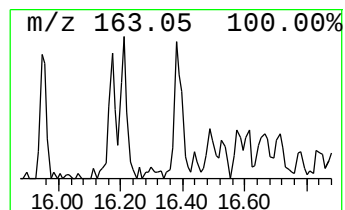
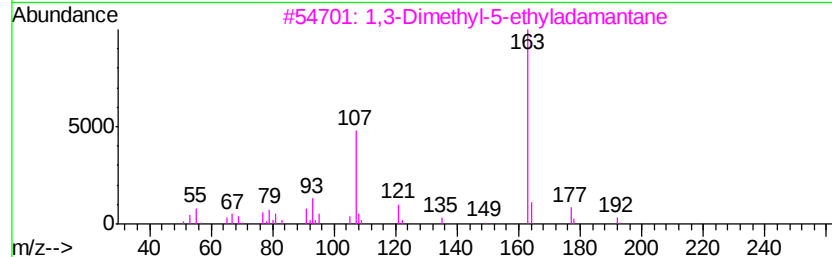
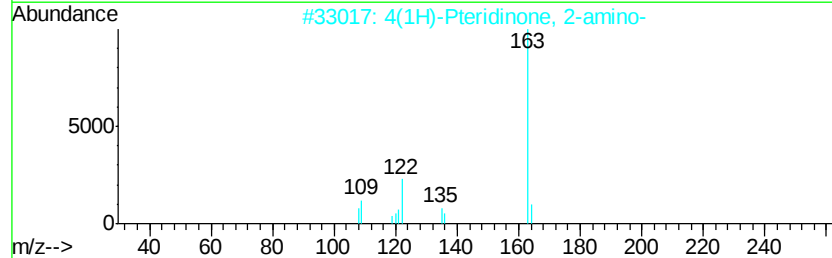
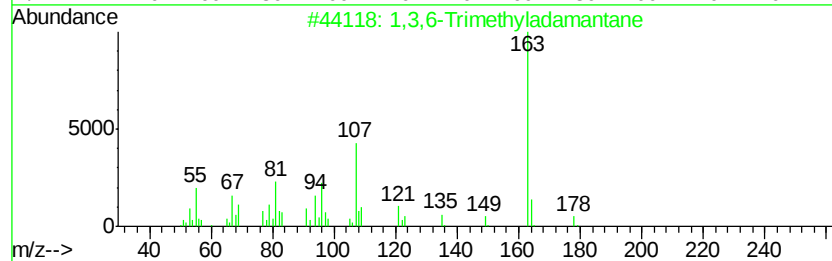
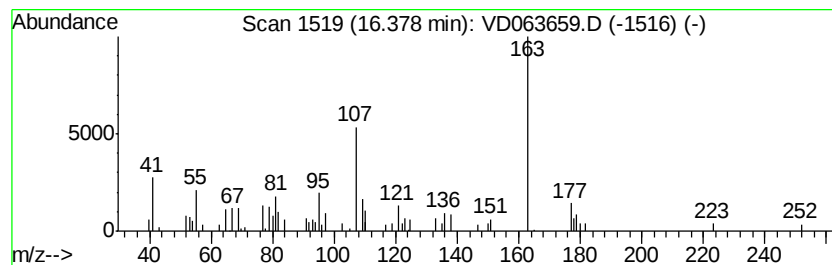
Instrument :
MSVOA_D
ClientSampleId :
T8-C-(0)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080919S.M
Quant Title : SW846 8260

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

Peak Number 9 4(1H)-Pteridinone, 2-amino- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.38	13.79 ug/l	201524	1,4-Dichlorobenzene-d4	14.52
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1,3,6-Trimethyladamantane	178 C13H22	024139-37-5	64
2	4(1H)-Pteridinone, 2-amino-	163 C6H5N5O	002236-60-4	64
3	1,3-Dimethyl-5-ethyladamantane	192 C14H24	001687-35-0	53
4	1,3,5-Trimethyladamantane	178 C13H22	000707-35-7	50
5	Tricyclo[3.3.1.1(3,7)-]decane, 1...	242 C12H19Br	000941-37-7	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD081919\
Data File : VD063659.D
Acq On : 19 Aug 2019 15:56
Operator : JC/SY
Sample : K4385-14
Misc : 4.75a/5ml/MSVOA D/SOIL
ALS Vial : 20 Sample Multiplier: 1

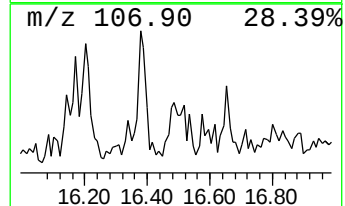
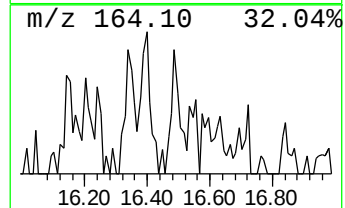
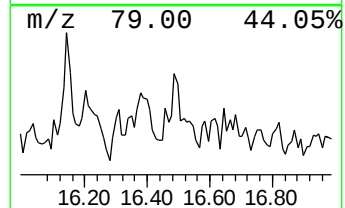
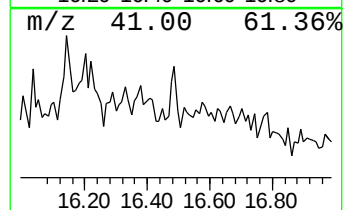
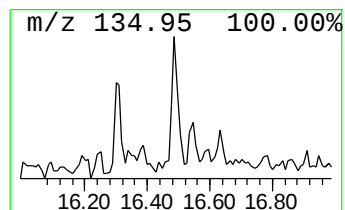
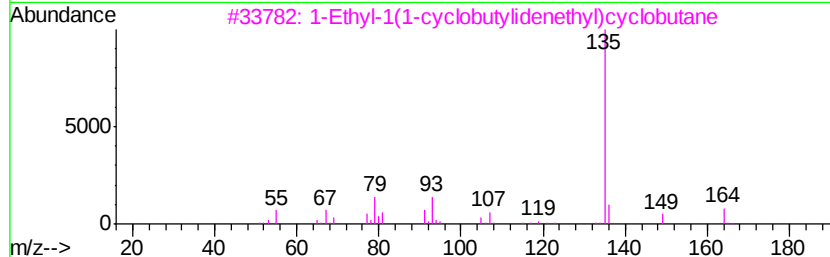
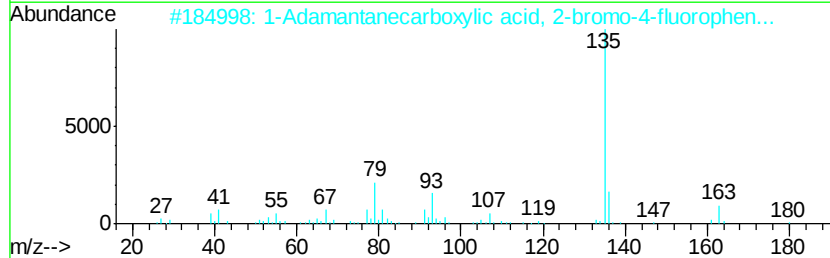
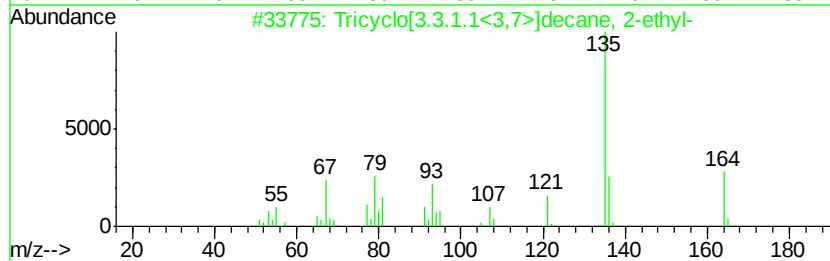
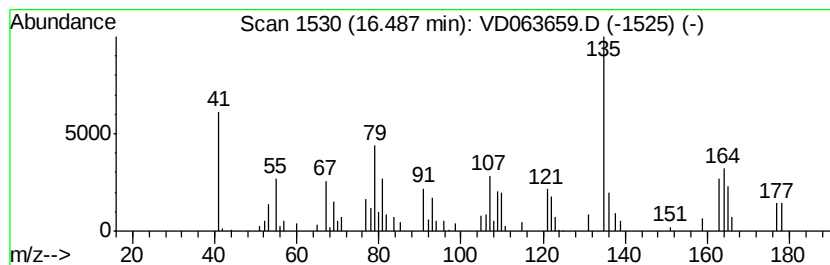
Instrument :
MSVOA_D
ClientSampleId :
T8-C-(0)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080919S.M
Quant Title : SW846 8260

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

Peak Number 10 unknown16.49 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.49	11.16 ug/l	163064	1,4-Dichlorobenzene-d4	14.52
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tricvclo[3.3.1.1<3,7>]decane, 2-...	164 C12H20	014451-87-7 49
2		1-Adamantanecarboxylic acid, 2-b...	352 C17H18BrF02	1000293-75-5 46
3		1-Ethyl-1(1-cvclobutylidenethyl)...	164 C12H20	1000215-13-7 43
4		4-n-Propylbenzoic acid	164 C10H12O2	002438-05-3 43
5		Pentafluoropropionic acid, 2-(1-...	326 C15H19F5O2	1000283-03-9 43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD081919\
Data File : VD063659.D
Acq On : 19 Aug 2019 15:56
Operator : JC/SY
Sample : K4385-14
Misc : 4.75g/5ml/MSVOA_D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
T8-C-(0)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080919S.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
Cyclohexane, 1-me...	13.47	15.4	ug/l	224976	4	14.52	730478	50.0
unknown13.71	13.71	18.3	ug/l	267223	4	14.52	730478	50.0
unknown13.82	13.82	10.8	ug/l	157225	4	14.52	730478	50.0
Tricyclo[3.3.1.1(...	15.29	11.1	ug/l	162084	4	14.52	730478	50.0
2H-Inden-2-one, 1...	15.81	12.4	ug/l	181448	4	14.52	730478	50.0
unknown16.14	16.14	11.5	ug/l	168299	4	14.52	730478	50.0
unknown16.20	16.20	11.3	ug/l	165603	4	14.52	730478	50.0
4(1H)-Pteridinone...	16.38	13.8	ug/l	201524	4	14.52	730478	50.0
unknown16.49	16.49	11.2	ug/l	163064	4	14.52	730478	50.0