

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD092018\
 Data File : VD060054.D
 Acq On : 20 Sep 2018 18:08
 Operator : VA/AP
 Sample : J5017-02
 Misc : 5.10g/5ml/MSVOA D/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 RA-091918-N-FL-056

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\82D091318S.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	1.945	40	43	54	rVB	36456	116996	3.15%	0.215%
2	3.145	160	165	172	rBV	22954	62109	1.67%	0.114%
3	3.706	216	222	231	rVB	108637	289800	7.81%	0.533%
4	4.504	296	303	308	rBV2	16716	46405	1.25%	0.085%
5	5.635	410	418	432	rBV2	35786	136292	3.67%	0.251%
6	6.314	480	487	491	rBV	656002	1711667	46.15%	3.147%
7	6.393	491	495	512	rVB	1149011	2966870	80.00%	5.455%
8	6.767	525	533	549	rBV2	523783	1416128	38.18%	2.604%
9	7.426	595	600	617	rBV	1546821	3708691	100.00%	6.819%
10	9.444	799	805	823	rBV	1530744	3364276	90.71%	6.186%
11	10.182	875	880	887	rBV	191510	516624	13.93%	0.950%
12	10.772	936	940	944	rBV2	51385	102510	2.76%	0.188%
13	11.038	963	967	969	rBV2	19908	38772	1.05%	0.071%
14	11.077	969	971	976	rVB2	26551	55057	1.48%	0.101%
15	11.215	979	985	987	rBV	80733	155164	4.18%	0.285%
16	11.274	987	991	1004	rVB	2037156	3506688	94.55%	6.448%
17	11.520	1013	1016	1021	rBV5	23657	81557	2.20%	0.150%
18	11.589	1021	1023	1025	rVV	52670	91268	2.46%	0.168%
19	11.638	1025	1028	1032	rVB3	43738	103523	2.79%	0.190%
20	11.737	1035	1038	1042	rBV4	23606	42341	1.14%	0.078%
21	11.884	1048	1053	1056	rBV3	73747	196484	5.30%	0.361%
22	11.963	1059	1061	1065	rVV3	38937	83363	2.25%	0.153%
23	12.091	1069	1074	1077	rVV3	262371	489194	13.19%	0.899%
24	12.199	1077	1085	1088	rVV2	291583	736246	19.85%	1.354%
25	12.268	1090	1092	1096	rVB2	80229	129515	3.49%	0.238%
26	12.347	1096	1100	1103	rBV3	157861	383826	10.35%	0.706%
27	12.425	1103	1108	1112	rVB2	1475111	2443195	65.88%	4.492%
28	12.484	1112	1114	1116	rBV2	53676	71060	1.92%	0.131%
29	12.524	1116	1118	1120	rVV	228243	300132	8.09%	0.552%
30	12.573	1120	1123	1130	rVB2	373336	770001	20.76%	1.416%
31	12.681	1130	1134	1138	rBV3	135819	296071	7.98%	0.544%
32	12.780	1138	1144	1146	rBV4	282578	756762	20.41%	1.391%
33	12.819	1146	1148	1152	rVB	484801	827584	22.31%	1.522%
34	12.888	1152	1155	1156	rBV2	54057	90249	2.43%	0.166%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	12.917	1156	1158	1159	rVV	73956	87065	2.35%	0.160%
36	12.957	1159	1162	1166	rVV4	162904	421691	11.37%	0.775%
37	13.016	1166	1168	1169	rVV	138983	198176	5.34%	0.364%
38	13.055	1169	1172	1174	rVV	764425	1208266	32.58%	2.222%
39	13.104	1174	1177	1182	rVV4	212173	468496	12.63%	0.861%
40	13.173	1182	1184	1186	rVV	132072	220857	5.96%	0.406%
41	13.223	1186	1189	1193	rVV3	463075	918760	24.77%	1.689%
42	13.282	1193	1195	1201	rVV3	546292	1246321	33.61%	2.292%
43	13.370	1201	1204	1208	rVV	1999499	3161553	85.25%	5.813%
44	13.429	1208	1210	1212	rVV	271519	381331	10.28%	0.701%
45	13.498	1214	1217	1220	rVV2	354285	539487	14.55%	0.992%
46	13.557	1220	1223	1227	rVB2	222833	386165	10.41%	0.710%
47	13.626	1227	1230	1231	rBV	125501	191844	5.17%	0.353%
48	13.655	1231	1233	1236	rVV3	460952	787845	21.24%	1.449%
49	13.754	1236	1243	1245	rVV2	337806	1285901	34.67%	2.364%
50	13.793	1245	1247	1251	rVV3	295382	613283	16.54%	1.128%
51	13.862	1251	1254	1256	rVV3	381465	838165	22.60%	1.541%
52	13.902	1256	1258	1260	rVV2	287355	437922	11.81%	0.805%
53	13.941	1260	1262	1263	rVV	124835	212011	5.72%	0.390%
54	13.980	1263	1266	1271	rVV3	307835	844804	22.78%	1.553%
55	14.069	1271	1275	1277	rVV4	186629	397049	10.71%	0.730%
56	14.128	1277	1281	1282	rVV4	362194	553896	14.94%	1.018%
57	14.157	1282	1284	1286	rVV2	554722	907800	24.48%	1.669%
58	14.187	1286	1287	1289	rVV2	325831	429535	11.58%	0.790%
59	14.226	1289	1291	1293	rVB	289822	365032	9.84%	0.671%
60	14.275	1293	1296	1299	rBV2	427211	733635	19.78%	1.349%
61	14.315	1299	1300	1302	rVV2	90348	109443	2.95%	0.201%
62	14.354	1302	1304	1305	rVV	78702	101070	2.73%	0.186%
63	14.423	1309	1311	1314	rVV4	134570	304363	8.21%	0.560%
64	14.472	1314	1316	1318	rVV	176603	247488	6.67%	0.455%
65	14.512	1318	1320	1325	rVV3	482381	1101193	29.69%	2.025%
66	14.581	1325	1327	1330	rVV	511673	805107	21.71%	1.480%
67	14.649	1330	1334	1336	rVV4	208761	442483	11.93%	0.814%
68	14.708	1336	1340	1341	rVV3	93526	212701	5.74%	0.391%
69	14.768	1341	1346	1348	rVV4	259236	776060	20.93%	1.427%
70	14.836	1348	1353	1355	rVV4	335657	792058	21.36%	1.456%
71	14.886	1356	1358	1363	rVV3	302370	760917	20.52%	1.399%

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72	14.984	1365	1368	1371	rVV2	485115	847202	22.84%	1.558%
73	15.063	1371	1376	1379	rVV4	338034	843144	22.73%	1.550%
74	15.132	1382	1383	1385	rVV2	179660	273471	7.37%	0.503%
75	15.201	1388	1390	1391	rVV	152533	229712	6.19%	0.422%
76	15.230	1391	1393	1394	rVV2	180498	276541	7.46%	0.508%
77	15.260	1394	1396	1399	rVV3	245377	456990	12.32%	0.840%
78	15.319	1399	1402	1407	rVV5	242842	657066	17.72%	1.208%
79	15.417	1410	1412	1415	rVV2	119989	218122	5.88%	0.401%
80	15.466	1415	1417	1420	rVV2	105779	215550	5.81%	0.396%
81	15.574	1424	1428	1431	rVB5	193348	341737	9.21%	0.628%
82	15.624	1431	1433	1437	rVB5	53823	103346	2.79%	0.190%
83	15.683	1437	1439	1441	rBV3	60003	114813	3.10%	0.211%
84	15.899	1459	1461	1464	rVB3	110262	163200	4.40%	0.300%
85	16.598	1526	1532	1536	rVB5	29334	69418	1.87%	0.128%

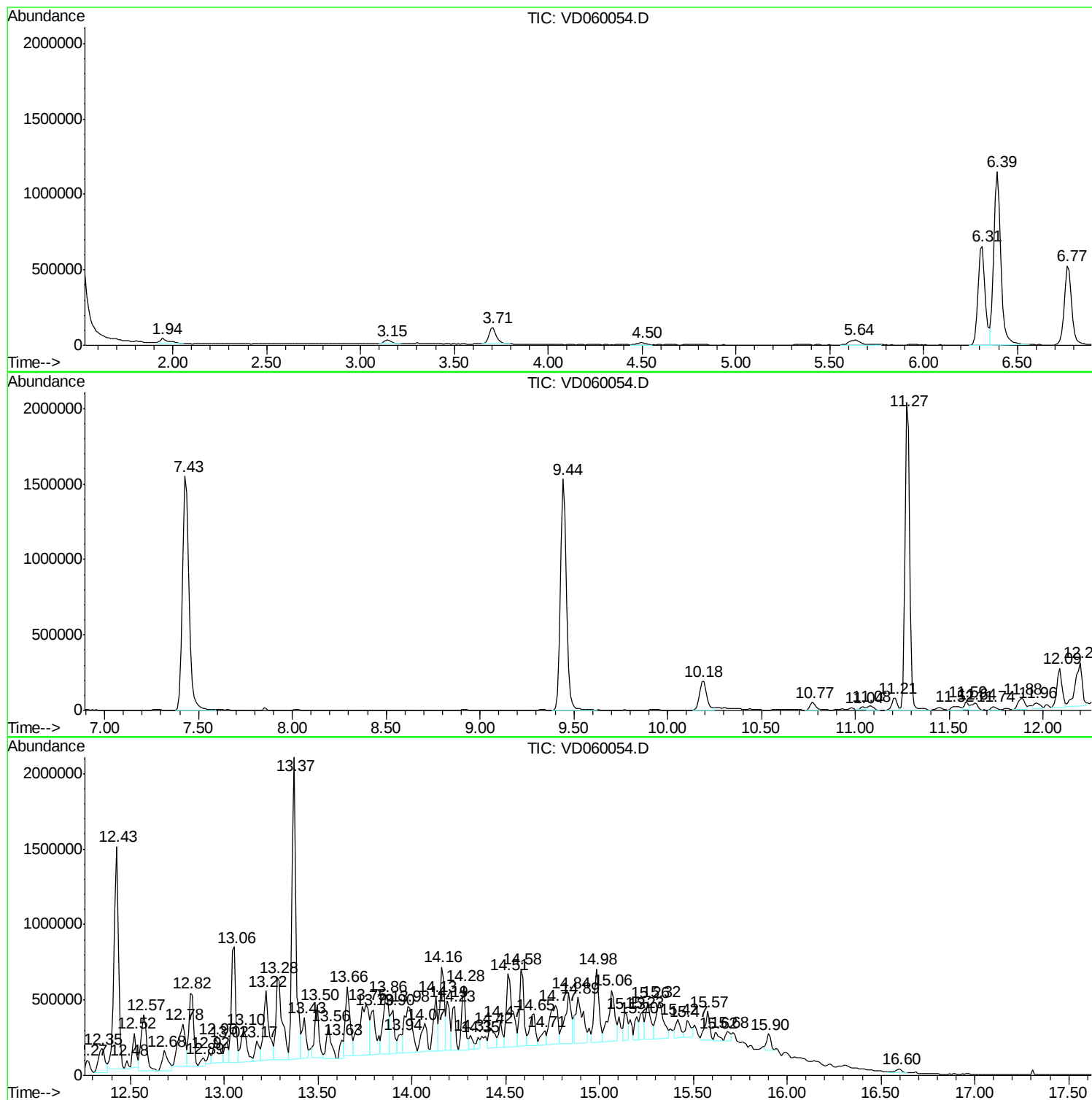
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D091318S.M
Quant Title : SW846 8260

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



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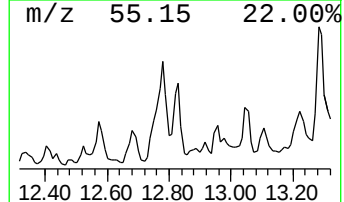
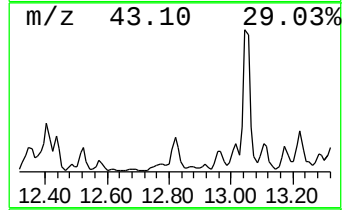
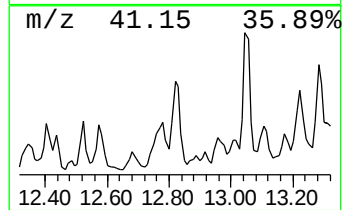
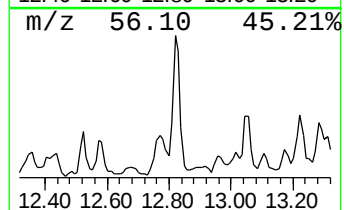
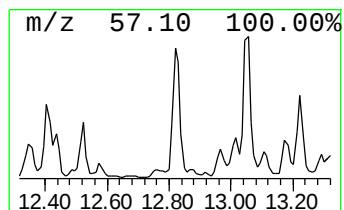
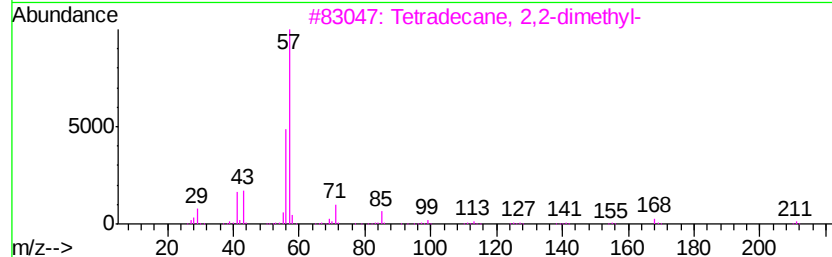
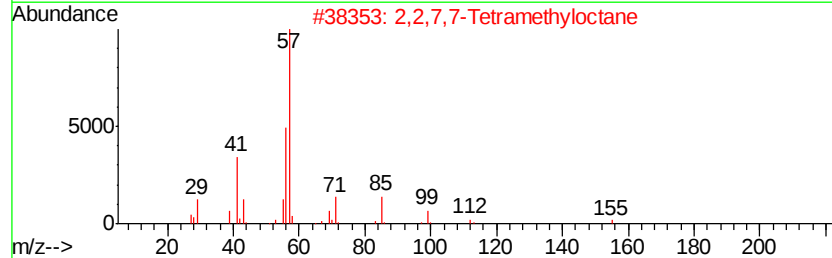
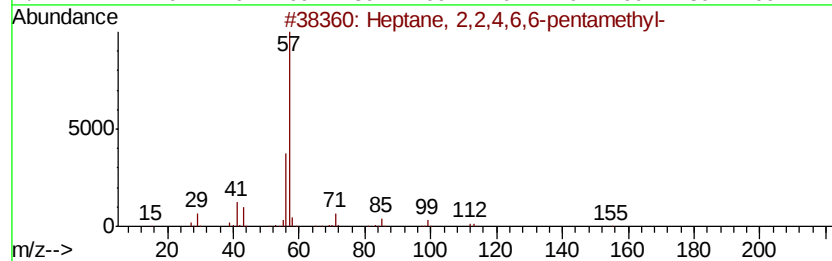
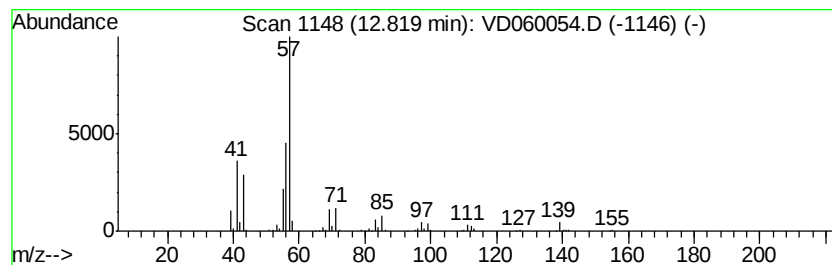
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D091318S.M
Quant Title : SW846 8260

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

Peak Number 1 Heptane, 2,2,4,6,6-pentamet... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.82	13.09 ug/l	827584	1,4-Dichlorobenzene-d4	13.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	64
2		2,2,7,7-Tetramethyloctane	170	C12H26	001071-31-4	64
3		Tetradecane, 2,2-dimethyl-	226	C16H34	059222-86-5	64
4		Decane, 2,2,8-trimethyl-	184	C13H28	062238-01-1	64
5		Decane, 2,2-dimethyl-	170	C12H26	017302-37-3	64



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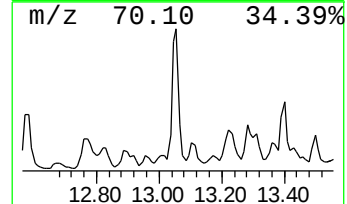
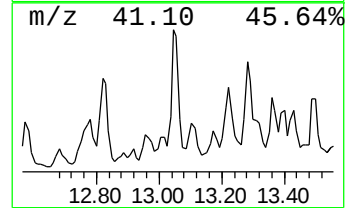
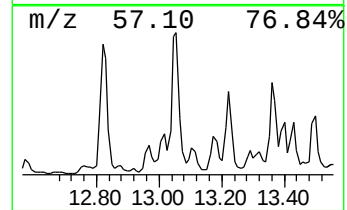
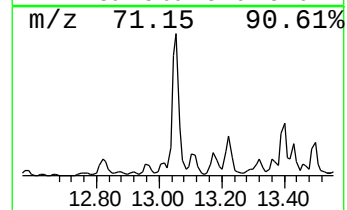
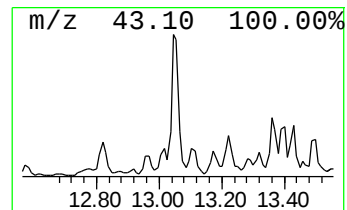
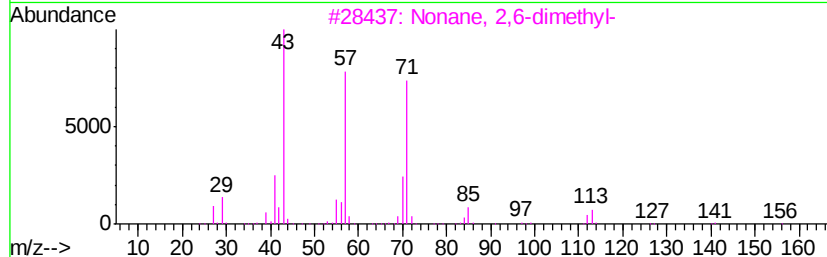
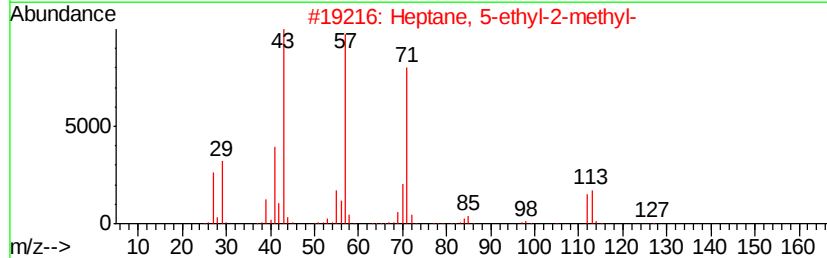
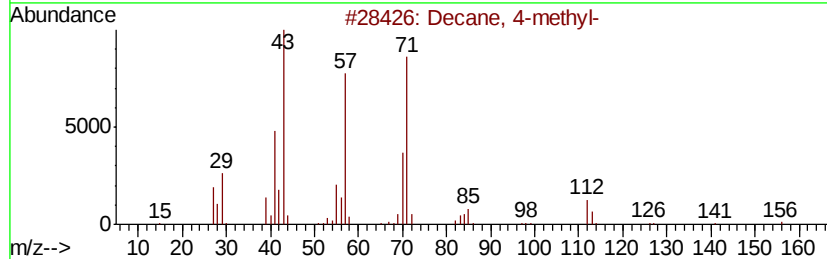
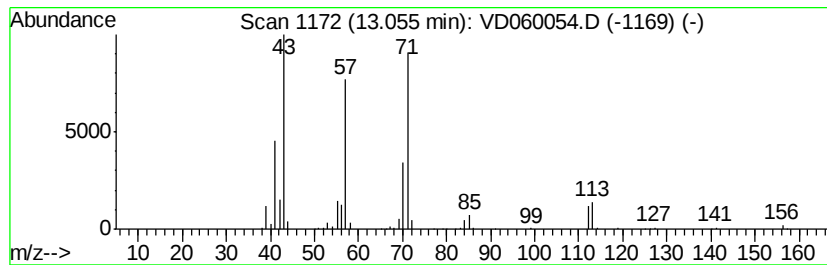
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Decane, 4-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.06	19.11 ug/l	1208270	1,4-Dichlorobenzene-d4	13.37	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 4-methyl-	156	C11H24	002847-72-5	94
2	Heptane, 5-ethyl-2-methyl-	142	C10H22	013475-78-0	78
3	Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	78
4	Heptane, 2,5,5-trimethyl-	142	C10H22	001189-99-7	72
5	Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	72



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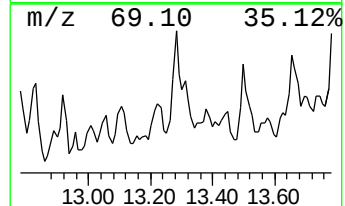
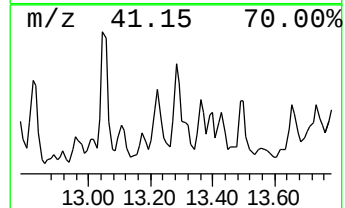
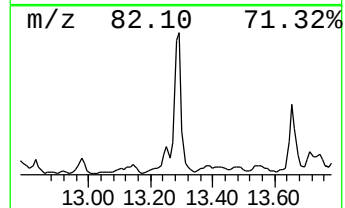
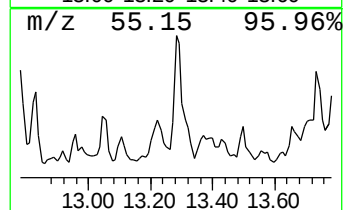
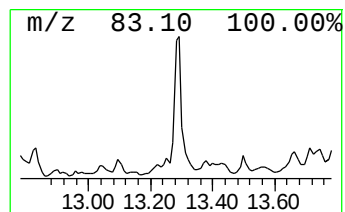
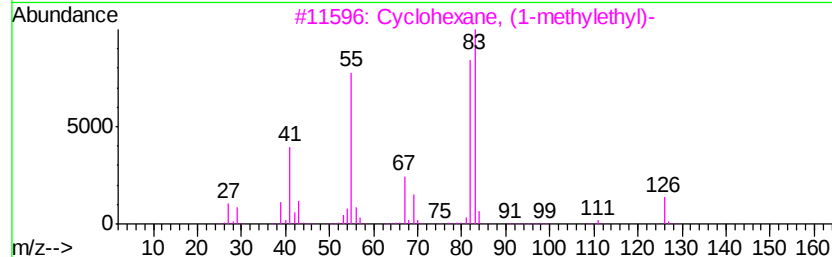
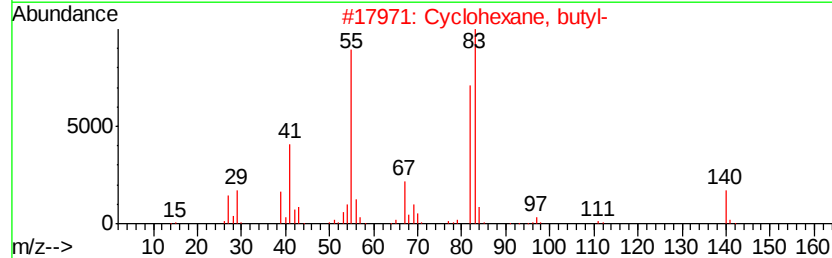
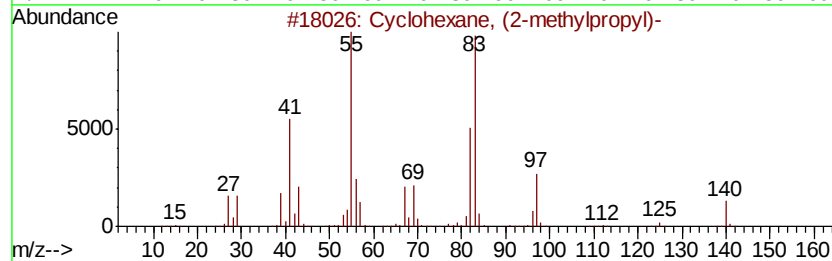
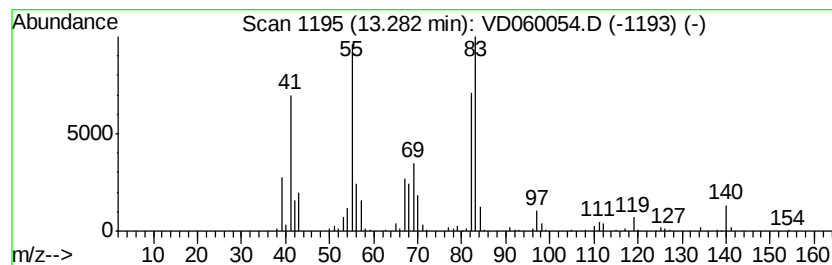
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Quant Title : SW846 8260

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

Peak Number 3 Cyclohexane, (2-methylpropyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.28	19.71 ug/l	1246320	1,4-Dichlorobenzene-d4	13.37		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	87
2		Cyclohexane, butyl-	140	C10H20	001678-93-9	72
3		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	64
4		Cyclohexane, (1-methylpropyl)-	140	C10H20	007058-01-7	53
5		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	38



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RA-091918-N-FL-056

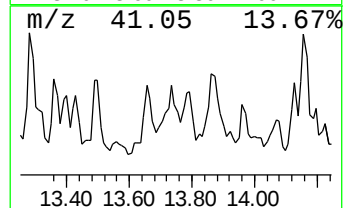
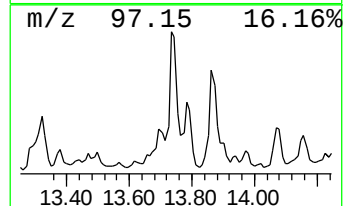
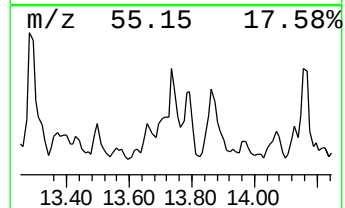
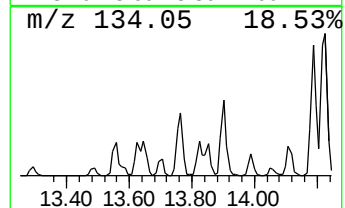
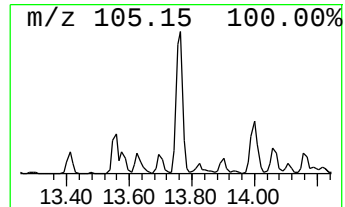
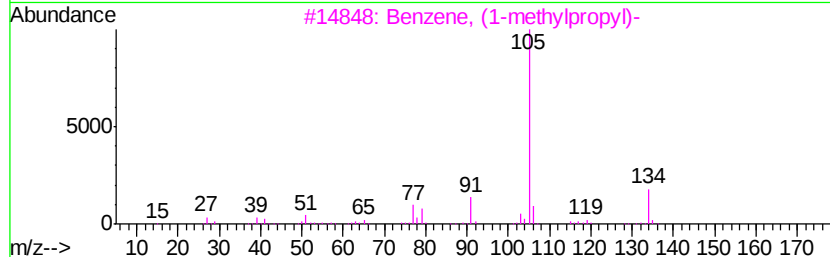
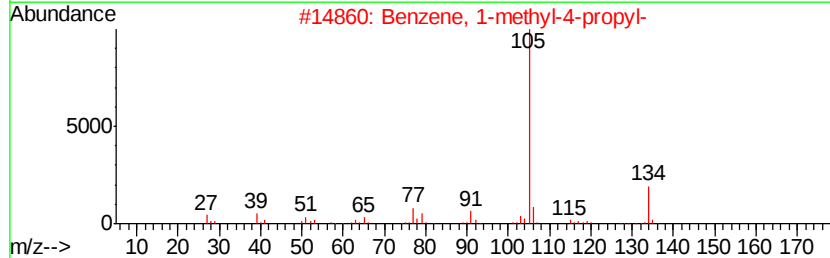
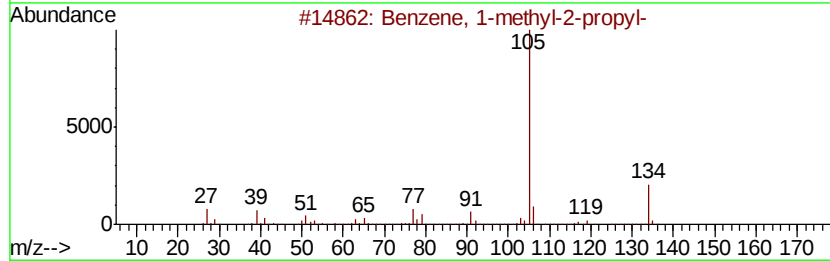
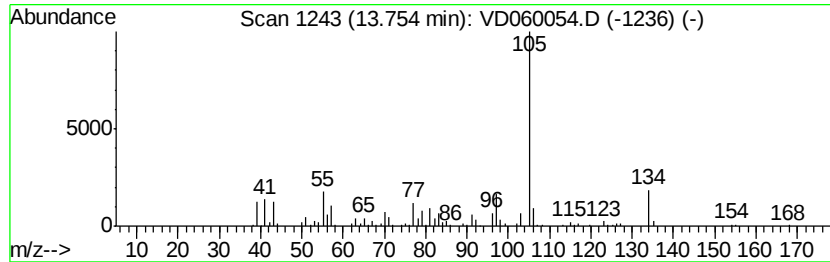
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-methyl-2-propyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.75	20.34 ug/l	1285900	1,4-Dichlorobenzene-d4	13.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	60
2		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	60
3		Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	53
4		Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	49
5		Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD092018\
Data File : VD060054.D
Acq On : 20 Sep 2018 18:08
Operator : VA/AP
Sample : J5017-02
Misc : 5.10g/5ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
RA-091918-N-FL-056

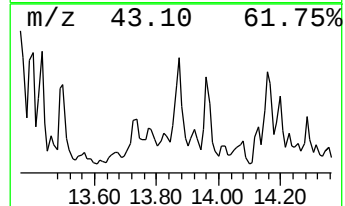
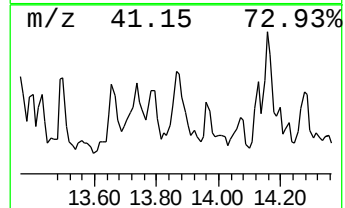
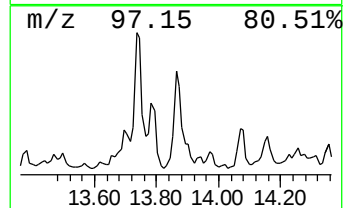
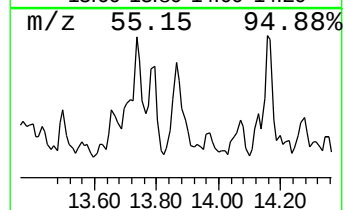
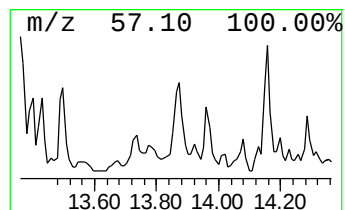
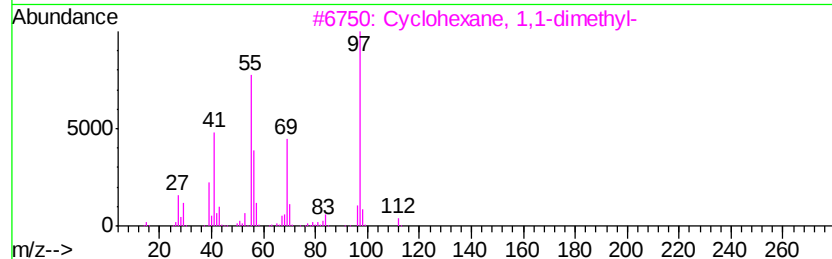
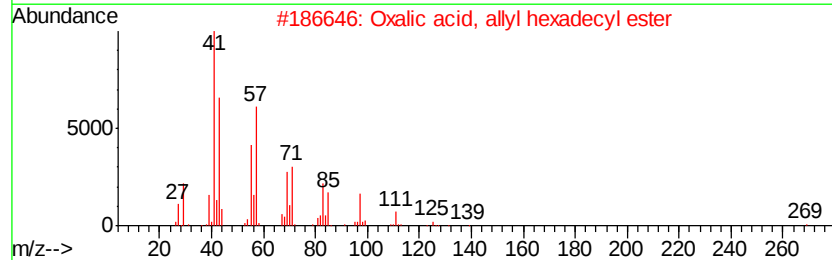
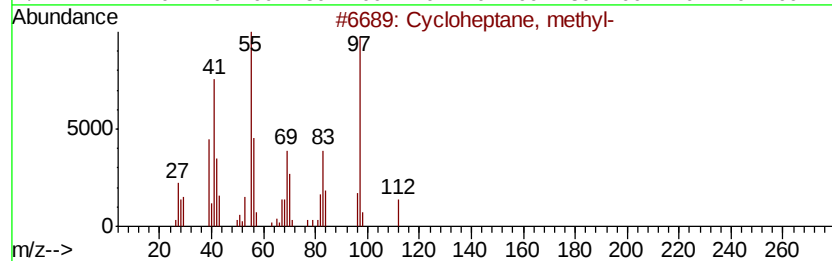
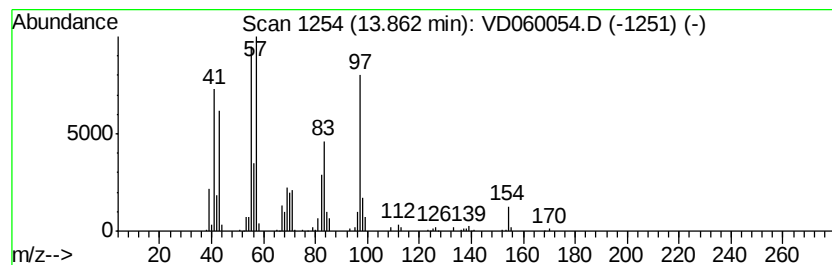
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Quant Title : SW846 8260

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

Peak Number 5 unknown13.86 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	13.26 ug/l	838165	1,4-Dichlorobenzene-d4	13.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cycloheptane, methyl-	112	C8H16	004126-78-7	38
2		Oxalic acid, allyl hexadecyl ester	354	C21H38O4	1000309-24-4	38
3		Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	38
4		Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	38
5		2-n-Butylacrolein	112	C7H12O	001070-66-2	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD092018\
Data File : VD060054.D
Acq On : 20 Sep 2018 18:08
Operator : VA/AP
Sample : J5017-02
Misc : 5.10g/5ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

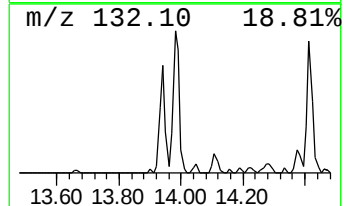
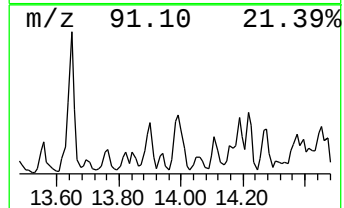
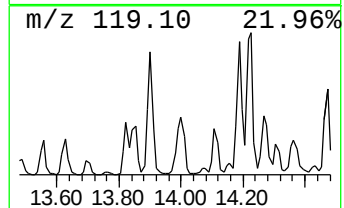
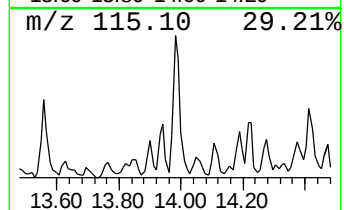
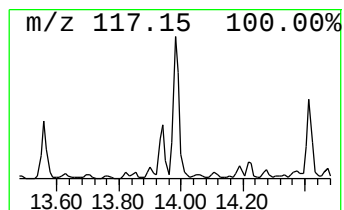
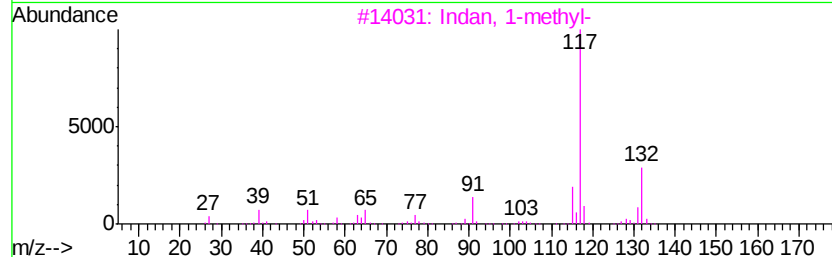
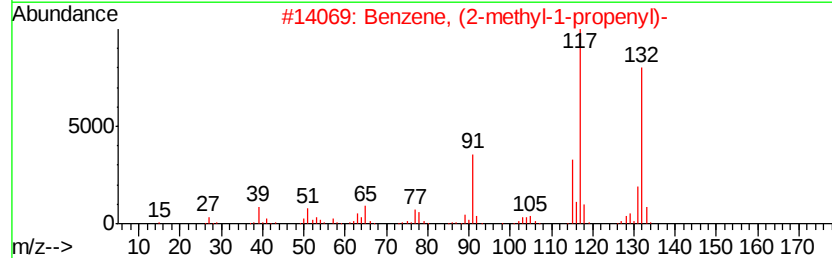
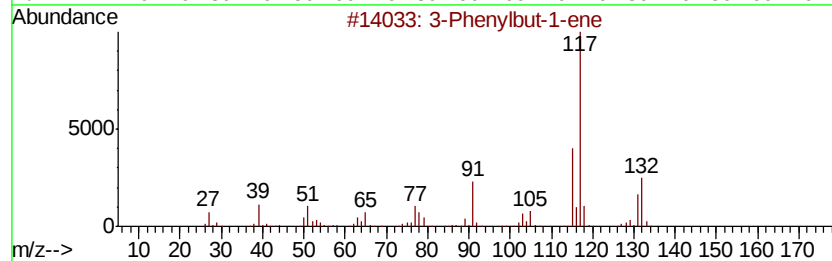
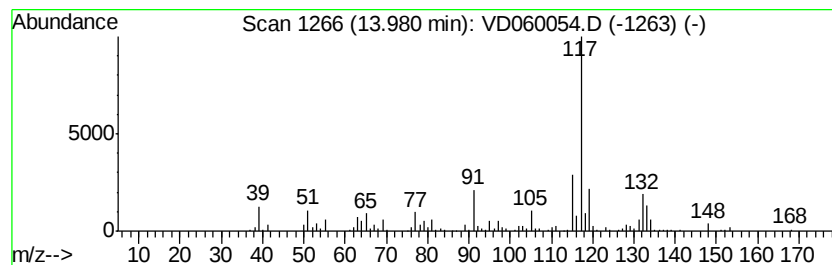
Instrument :
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ClientSampleId :
RA-091918-N-FL-056

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D091318S.M
Quant Title : SW846 8260

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

Peak Number 6 3-Phenylbut-1-ene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.98	13.36 ug/l	844804	1,4-Dichlorobenzene-d4	13.37
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	3-Phenylbut-1-ene	132 C10H12	000934-10-1	76
2	Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0	76
3	Indan, 1-methyl-	132 C10H12	000767-58-8	70
4	Benzene, (2-methylcyclopropyl)-	132 C10H12	1000327-39-0	62
5	1H-Indene, 2,3-dihydro-5-methyl-	132 C10H12	000874-35-1	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD092018\
Data File : VD060054.D
Acq On : 20 Sep 2018 18:08
Operator : VA/AP
Sample : J5017-02
Misc : 5.10g/5ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleID :
RA-091918-N-FL-056

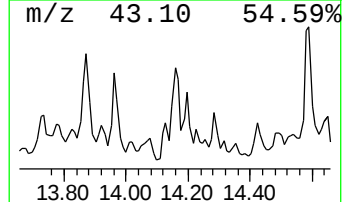
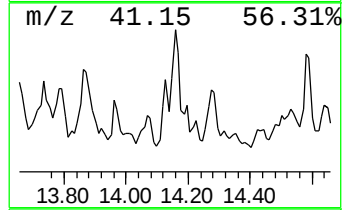
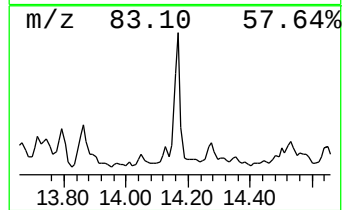
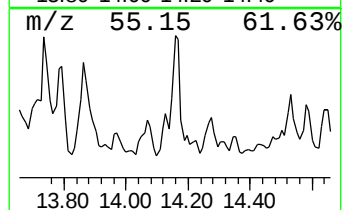
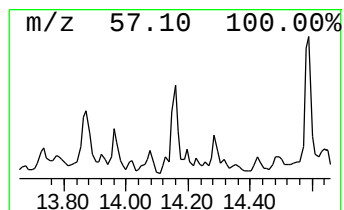
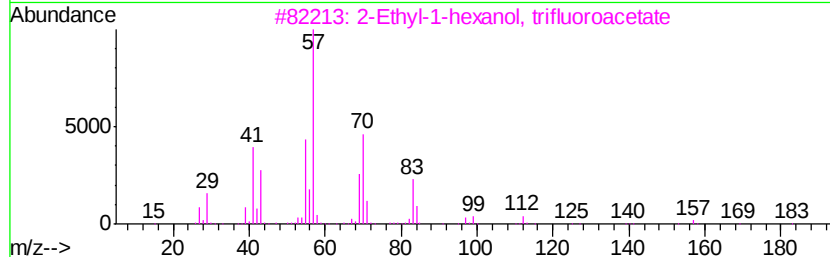
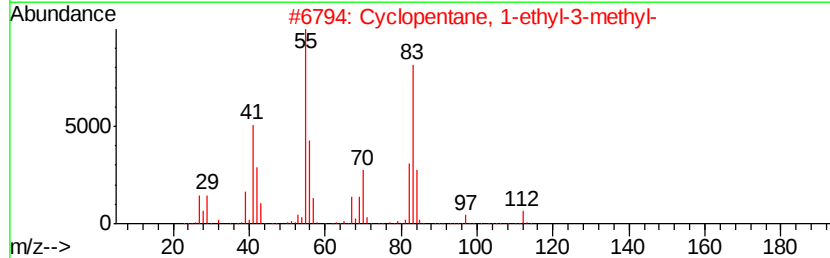
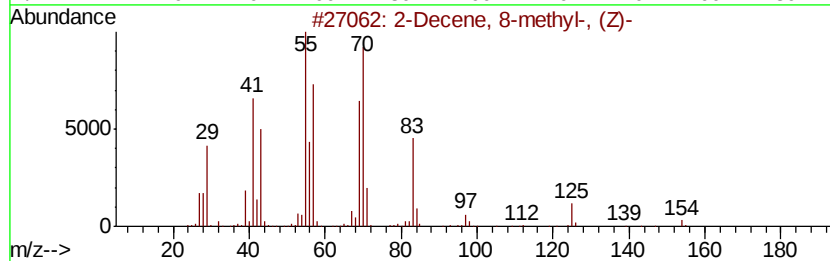
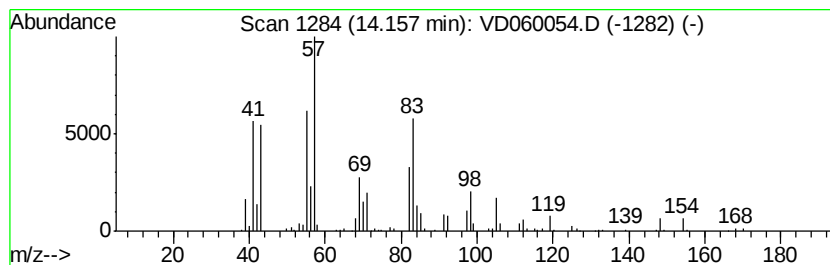
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Quant Title : SW846 8260

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

Peak Number 7 unknown14.16 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.16	14.36 ug/l	907800	1,4-Dichlorobenzene-d4	13.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Decene, 8-methyl-, (Z)-			154	C11H22	074630-25-4	49
2	Cyclopentane, 1-ethyl-3-methyl-			112	C8H16	003726-47-4	43
3	2-Ethyl-1-hexanol, trifluoroacetate			226	C10H17F3O2	1000365-19-6	38
4	Cyclohexane, methyl-			98	C7H14	000108-87-2	38
5	Decane, 2,5,6-trimethyl-			184	C13H28	062108-23-0	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD092018\
Data File : VD060054.D
Acq On : 20 Sep 2018 18:08
Operator : VA/AP
Sample : J5017-02
Misc : 5.10g/5ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

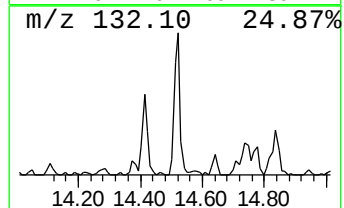
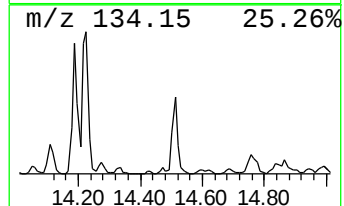
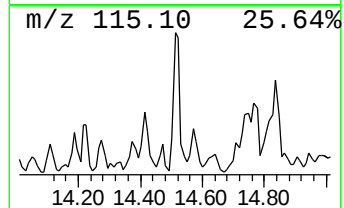
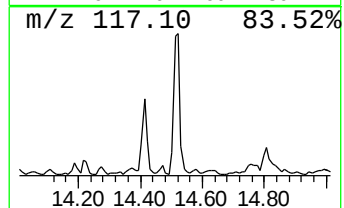
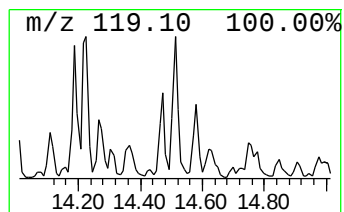
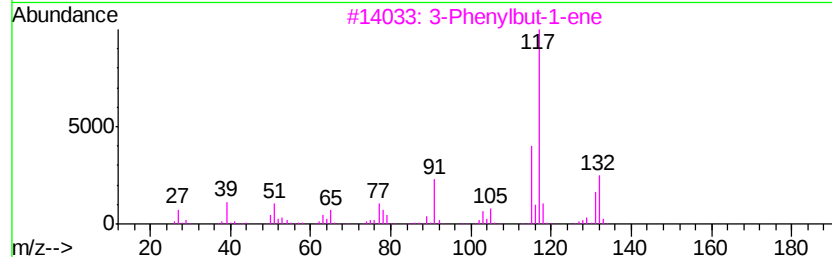
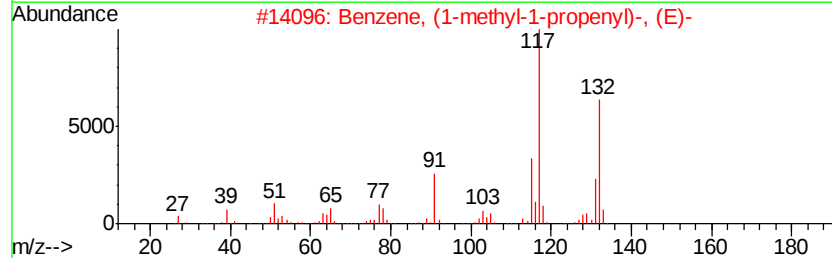
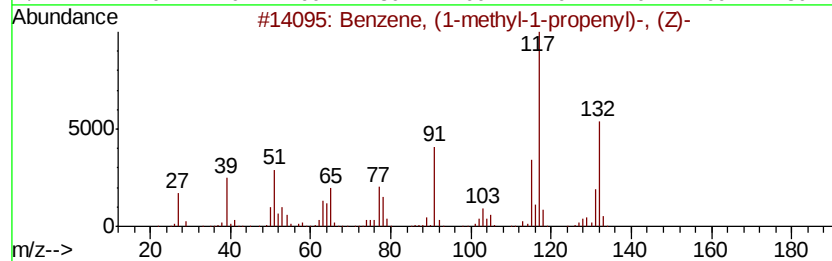
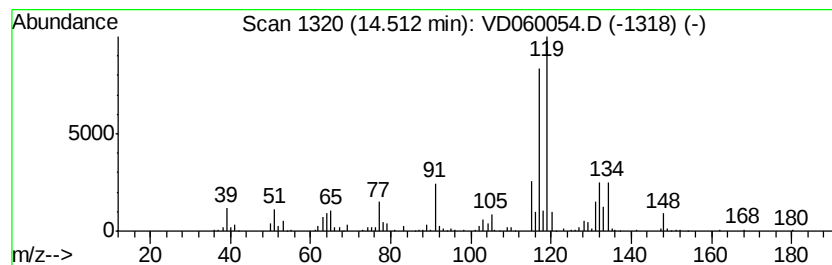
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ClientSampleId :
RA-091918-N-FL-056

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D091318S.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, (1-methyl-1-propen... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.51	17.42 ug/l	1101190	1,4-Dichlorobenzene-d4	13.37
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000767-99-7 87
2		Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000768-00-3 64
3		3-Phenylbut-1-ene	132 C10H12	000934-10-1 50
4		1H-Indene, 2,3-dihydro-5-methyl-	132 C10H12	000874-35-1 50
5		Benzene, 2-butenyl-	132 C10H12	001560-06-1 50



Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD092018\
Data File : VD060054.D
Acq On : 20 Sep 2018 18:08
Operator : VA/AP
Sample : J5017-02
Misc : 5.10g/5ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

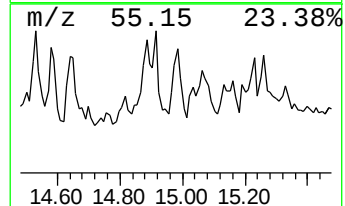
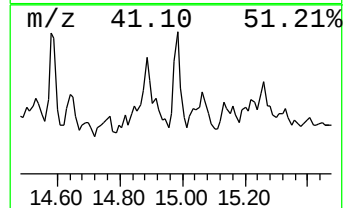
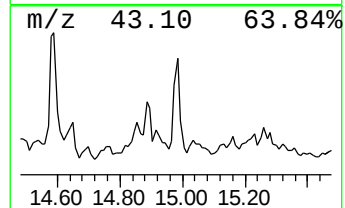
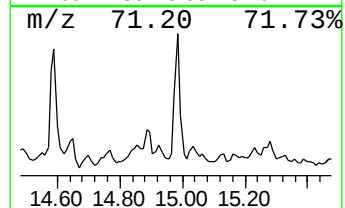
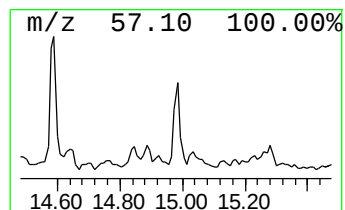
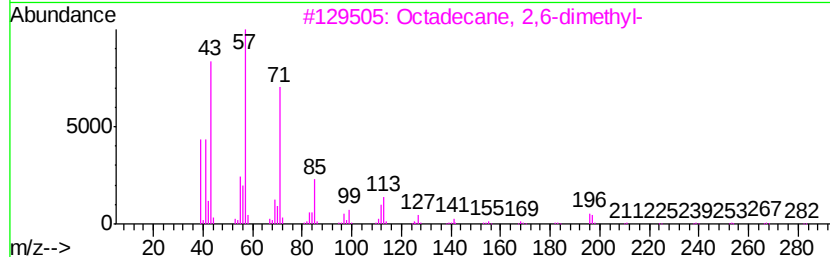
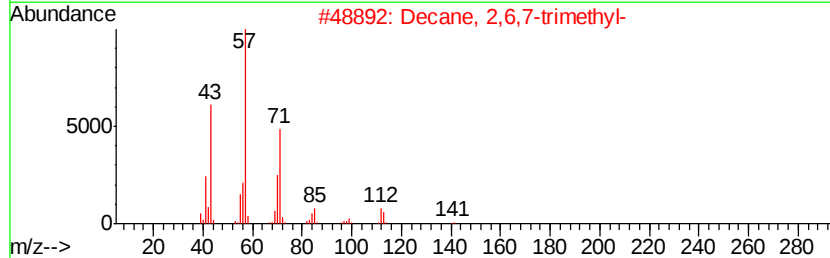
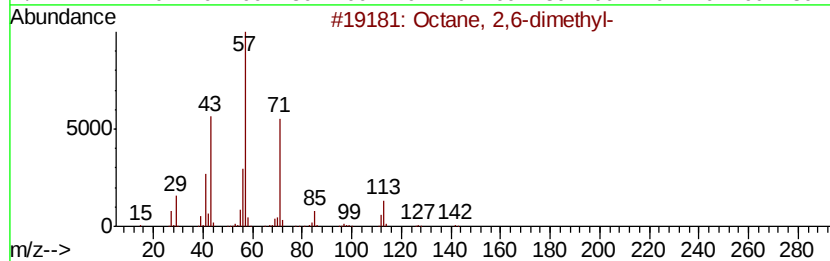
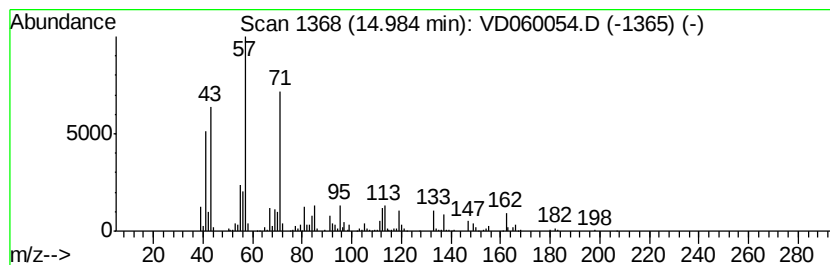
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D091318S.M
Quant Title : SW846 8260

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

Peak Number 9 Octane, 2,6-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.98	13.40 ug/l	847202	1,4-Dichlorobenzene-d4	13.37		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	53
2		Decane, 2,6,7-trimethyl-	184	C13H28	062108-25-2	53
3		Octadecane, 2,6-dimethyl-	282	C20H42	075163-97-2	50
4		Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1000309-19-2	50
5		Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD092018\
Data File : VD060054.D
Acq On : 20 Sep 2018 18:08
Operator : VA/AP
Sample : J5017-02
Misc : 5.10g/5ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
RA-091918-N-FL-056

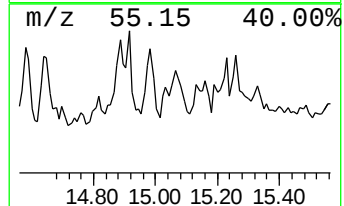
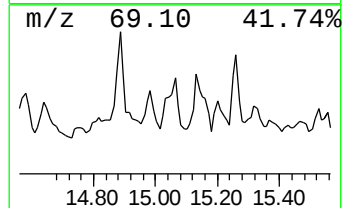
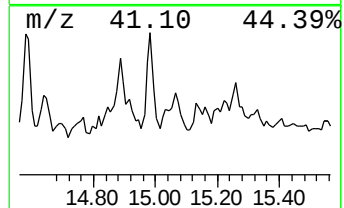
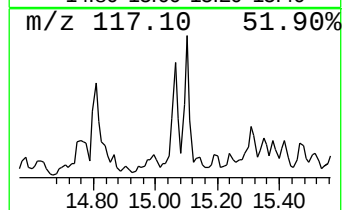
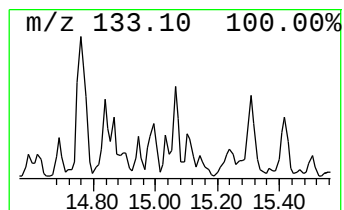
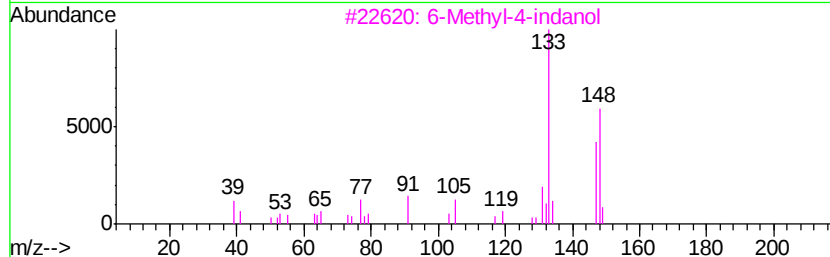
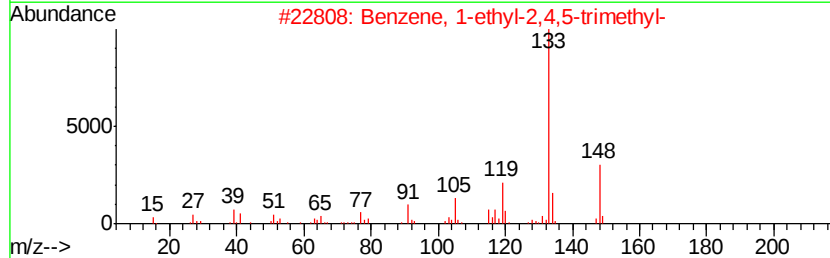
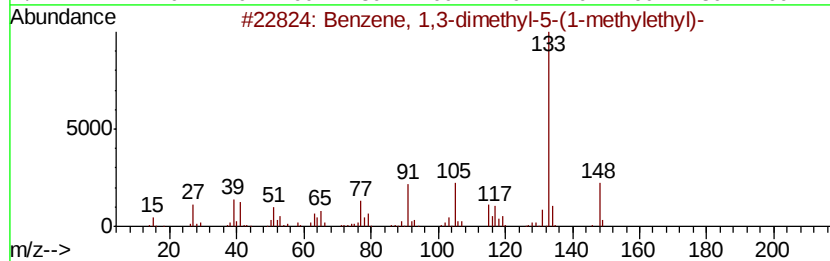
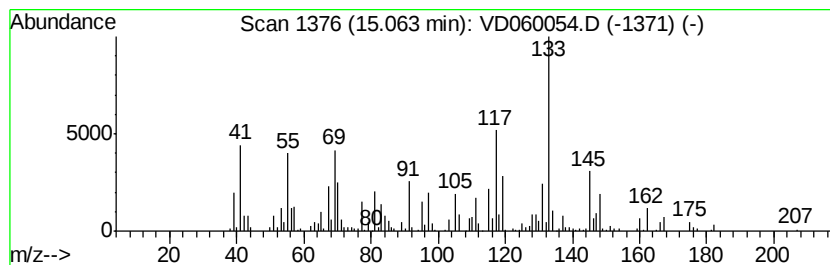
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D091318S.M
Quant Title : SW846 8260

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

Peak Number 10 unknown15.06 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.06	13.33 ug/l	843144	1,4-Dichlorobenzene-d4	13.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dimethyl-5-(1-methyl-...	148	C11H16	004706-90-5	46
2		Benzene, 1-ethyl-2,4,5-trimethyl-	148	C11H16	017851-27-3	43
3		6-Methyl-4-indanol	148	C10H12O	020294-32-0	43
4		1,5,6,7-Tetramethylbicyclo[3.2.0]...	148	C11H16	134329-46-7	35
5		4-tert-Butyltoluene	148	C11H16	000098-51-1	35



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD092018\
Data File : VD060054.D
Acq On : 20 Sep 2018 18:08
Operator : VA/AP
Sample : J5017-02
Misc : 5.10g/5ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
RA-091918-N-FL-056

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D091318S.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
Heptane, 2,2,4,6,...	12.82	13.1	ug/l	827584	4	13.37	3161550	50.0
Decane, 4-methyl-	13.06	19.1	ug/l	1208270	4	13.37	3161550	50.0
Cyclohexane, (2-m...	13.28	19.7	ug/l	1246320	4	13.37	3161550	50.0
Benzene, 1-methyl...	13.75	20.3	ug/l	1285900	4	13.37	3161550	50.0
unknown13.86	13.86	13.3	ug/l	838165	4	13.37	3161550	50.0
3-Phenylbut-1-ene	13.98	13.4	ug/l	844804	4	13.37	3161550	50.0
unknown14.16	14.16	14.4	ug/l	907800	4	13.37	3161550	50.0
Benzene, (1-methy...	14.51	17.4	ug/l	1101190	4	13.37	3161550	50.0
Octane, 2,6-dimet...	14.98	13.4	ug/l	847202	4	13.37	3161550	50.0
unknown15.06	15.06	13.3	ug/l	843144	4	13.37	3161550	50.0