

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD122118\
 Data File : VD060698.D
 Acq On : 21 Dec 2018 21:56
 Operator : VA/AP
 Sample : J6469-01
 Misc : 5.08g/5ml/MSVOA_D/SOIL
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 DRUM-1

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\82D122018S.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	1.450	15	23	55	rBV	755688	3120567	11.18%	0.521%
2	6.360	515	522	526	rBV	861885	2134485	7.65%	0.357%
3	6.439	526	530	550	rVB	1151737	3094845	11.09%	0.517%
4	6.823	563	569	583	rBV	623791	1544830	5.53%	0.258%
5	7.482	630	636	651	rBV	1702110	4036067	14.46%	0.674%
6	9.490	835	840	848	rBV	1874455	4254643	15.24%	0.711%
7	9.824	870	874	883	rVB	157320	326757	1.17%	0.055%
8	10.474	932	940	945	rBV	136813	291780	1.05%	0.049%
9	10.611	950	954	957	rBV	170686	391354	1.40%	0.065%
10	10.739	963	967	970	rVB	169674	364658	1.31%	0.061%
11	10.808	970	974	977	rVV2	194280	426826	1.53%	0.071%
12	11.015	992	995	997	rBV	225545	347854	1.25%	0.058%
13	11.123	997	1006	1012	rVB	1921083	5315255	19.04%	0.888%
14	11.241	1012	1018	1022	rBV	2019891	3809767	13.65%	0.637%
15	11.310	1022	1025	1029	rVB	2164868	3384564	12.12%	0.566%
16	11.389	1029	1033	1037	rVB5	1000005	304923	1.09%	0.051%
17	11.458	1037	1040	1046	rBV2	777851	1768022	6.33%	0.295%
18	11.586	1046	1053	1056	rBV	4433812	9790886	35.08%	1.636%
19	11.655	1056	1060	1068	rVB2	6425131	13209388	47.32%	2.207%
20	11.763	1068	1071	1075	rBV2	215990	421662	1.51%	0.070%
21	11.832	1075	1078	1082	rBV2	284289	526479	1.89%	0.088%
22	11.920	1082	1087	1089	rBV3	1847273	4920815	17.63%	0.822%
23	11.960	1089	1091	1093	rBV	3956064	5349380	19.16%	0.894%
24	11.989	1093	1094	1098	rVB3	1537237	2551554	9.14%	0.426%
25	12.048	1098	1100	1103	rBV	769460	1131777	4.05%	0.189%
26	12.117	1103	1107	1111	rVB3	3682459	6201512	22.22%	1.036%
27	12.206	1111	1116	1121	rBV3	4034909	11645640	41.72%	1.946%
28	12.304	1124	1126	1129	rVB	1936721	2716843	9.73%	0.454%
29	12.373	1129	1133	1136	rBV3	3421953	8409790	30.13%	1.405%
30	12.442	1136	1140	1141	rBV2	5968245	10653469	38.17%	1.780%
31	12.471	1141	1143	1146	rVB	7517558	10551094	37.80%	1.763%
32	12.521	1146	1148	1149	rBV2	1285768	1647315	5.90%	0.275%
33	12.560	1149	1152	1155	rBV	6395877	10317515	36.96%	1.724%
34	12.609	1155	1157	1160	rVB3	2638764	3586081	12.85%	0.599%

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35	12.668	1160	1163	1165	rBV	4444766	6191249	22.18%	1.035%
36	12.737	1165	1170	1172	rBV	11638757	25120867	89.99%	4.198%
37	12.767	1172	1173	1176	rVB	5801552	7408142	26.54%	1.238%
38	12.826	1176	1179	1180	rBV2	7644722	14884571	53.32%	2.487%
39	12.914	1186	1188	1190	rVB2	853729	1182609	4.24%	0.198%
40	12.973	1190	1194	1196	rBV	9526463	19620668	70.29%	3.279%
41	13.042	1200	1201	1203	rVV	2972573	4273873	15.31%	0.714%
42	13.081	1203	1205	1207	rVV	7492582	11224889	40.21%	1.876%
43	13.131	1207	1210	1214	rVB	12600951	27914065	100.00%	4.665%
44	13.209	1214	1218	1220	rBV	3776546	7376619	26.43%	1.233%
45	13.249	1220	1222	1227	rVB4	5370873	10289713	36.86%	1.719%
46	13.318	1227	1229	1232	rBV	7937624	14766634	52.90%	2.468%
47	13.396	1235	1237	1239	rBV	6179559	8936428	32.01%	1.493%
48	13.446	1239	1242	1246	rVB2	7529752	22726755	81.42%	3.798%
49	13.534	1246	1251	1253	rVB3	7533764	15499484	55.53%	2.590%
50	13.593	1253	1257	1258	rBV	8620742	17716969	63.47%	2.961%
51	13.662	1261	1264	1270	rBV4	6842069	21955750	78.65%	3.669%
52	13.770	1272	1275	1282	rBV2	11376200	24156408	86.54%	4.037%
53	13.859	1282	1284	1285	rBV	3501993	5439095	19.49%	0.909%
54	13.938	1290	1292	1297	rVV2	5701354	14670064	52.55%	2.451%
55	14.016	1297	1300	1305	rVB3	5478269	14859001	53.23%	2.483%
56	14.105	1305	1309	1311	rVB4	5308875	9893039	35.44%	1.653%
57	14.164	1311	1315	1316	rBV2	5786495	13347920	47.82%	2.230%
58	14.193	1316	1318	1320	rVV2	4672777	8440384	30.24%	1.410%
59	14.262	1323	1325	1327	rVB2	5709668	7611311	27.27%	1.272%
60	14.312	1327	1330	1332	rBV3	5034947	10734338	38.45%	1.794%
61	14.449	1340	1344	1346	rBV2	3182048	7463843	26.74%	1.247%
62	14.508	1348	1350	1352	rVB2	3852855	5242370	18.78%	0.876%
63	14.548	1352	1354	1359	rBV3	5740511	10677133	38.25%	1.784%
64	14.676	1364	1367	1372	rVV	7669418	15545305	55.69%	2.598%
65	14.764	1372	1376	1379	rVV2	4811834	14545602	52.11%	2.431%
66	14.804	1379	1380	1382	rVV	4834868	4377546	15.68%	0.732%
67	14.863	1382	1386	1390	rVV2	5848885	14299950	51.23%	2.390%
68	14.912	1390	1391	1396	rVB4	2688041	5210316	18.67%	0.871%
69	14.991	1396	1399	1403	rBV2	2396035	5407115	19.37%	0.904%
70	15.040	1403	1404	1407	rVV	6375320	7576046	27.14%	1.266%
71	15.099	1407	1410	1415	rVB3	6036333	12260543	43.92%	2.049%

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Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
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72	15.227	1420	1423	1426	rVB	1404212	1965023	7.04%	0.328%
73	15.276	1426	1428	1431	rVB4	525724	907168	3.25%	0.152%
74	15.374	1431	1438	1443	rVB	3467540	6378416	22.85%	1.066%
75	15.492	1448	1450	1458	rVB	370310	690021	2.47%	0.115%
76	15.601	1458	1461	1464	rVB	532130	740689	2.65%	0.124%
77	15.788	1478	1480	1485	rVB	210544	353386	1.27%	0.059%

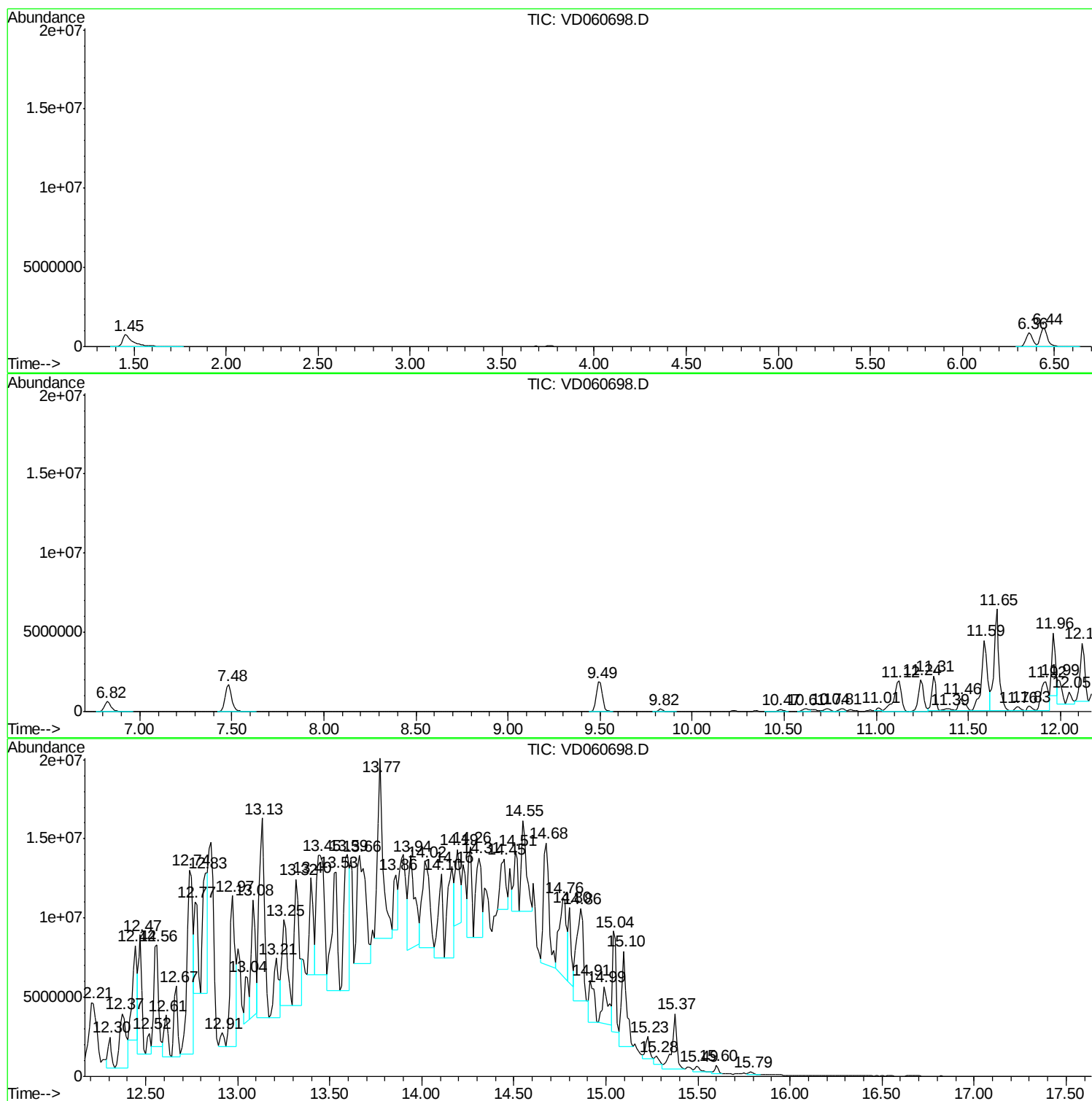
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Data Path : Z:\V0ASRV\HPCHEM1\MSV0A_D\DATA\VD122118\
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Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_D\METHOD\82D122018S.M
Quant Title : SW846 8260

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



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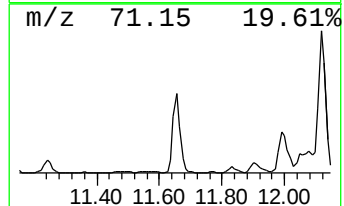
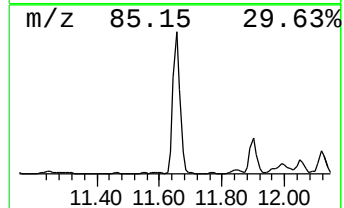
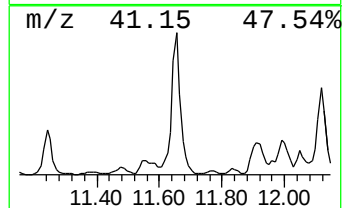
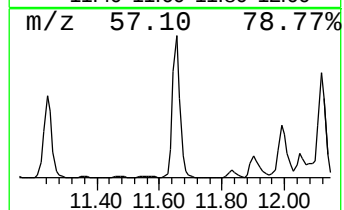
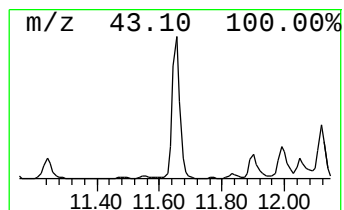
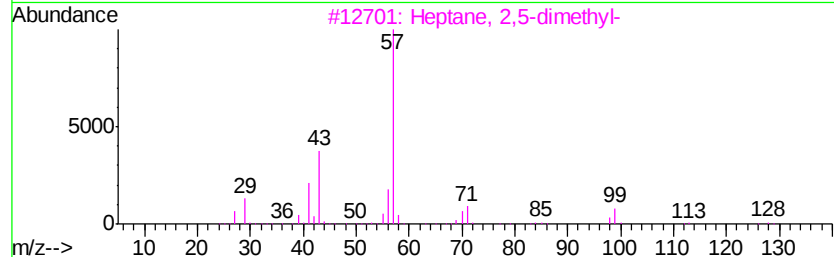
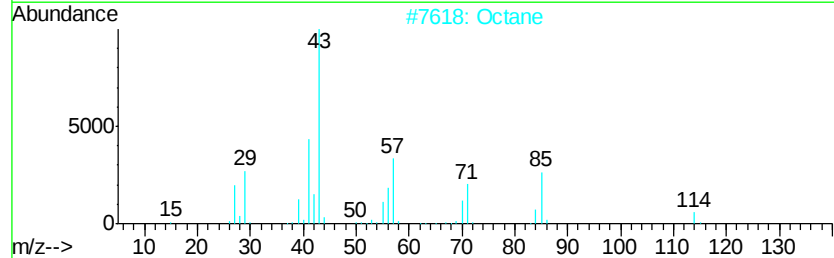
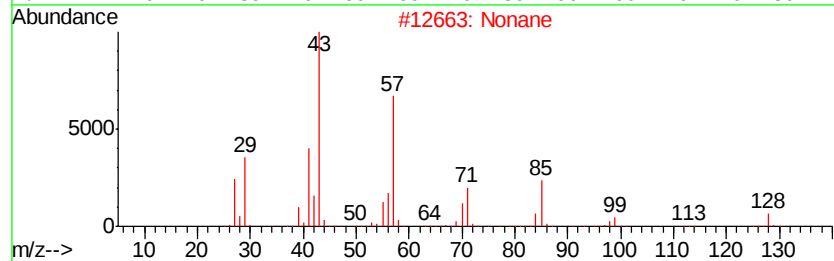
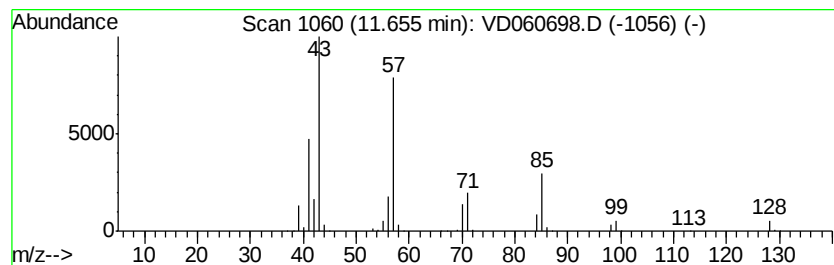
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D122018S.M
Quant Title : SW846 8260

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

Peak Number 1 Nonane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.65	195.14 ug/l	13209400	Chlorobenzene-d5	11.31		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane		128	C9H20	000111-84-2	97
2	Octane		114	C8H18	000111-65-9	64
3	Heptane, 2,5-dimethyl-		128	C9H20	002216-30-0	53
4	Heptane, 2,4-dimethyl-		128	C9H20	002213-23-2	50
5	Butanoic acid, 3-oxo-, 1,1-dimet...		158	C8H14O3	001694-31-1	50



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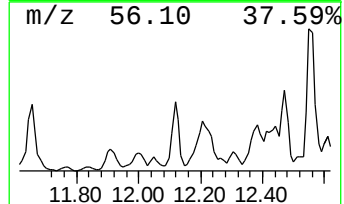
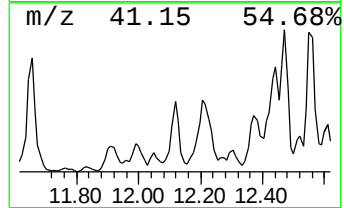
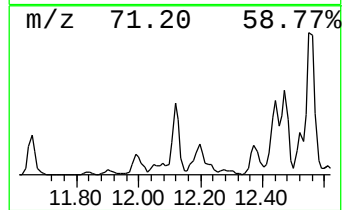
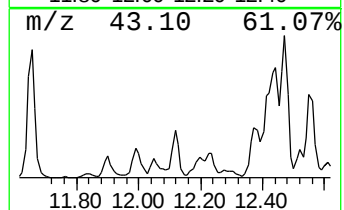
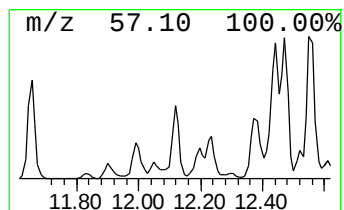
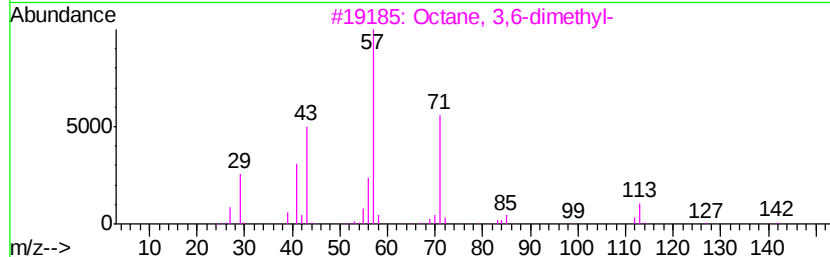
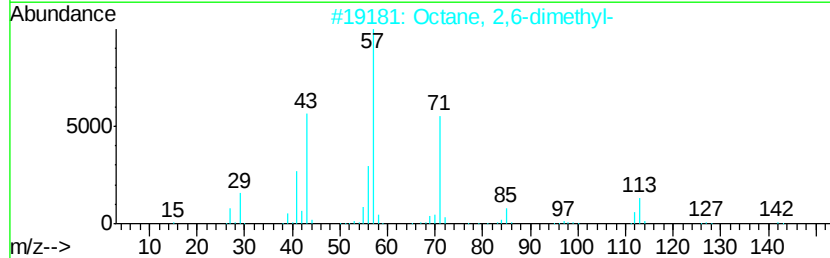
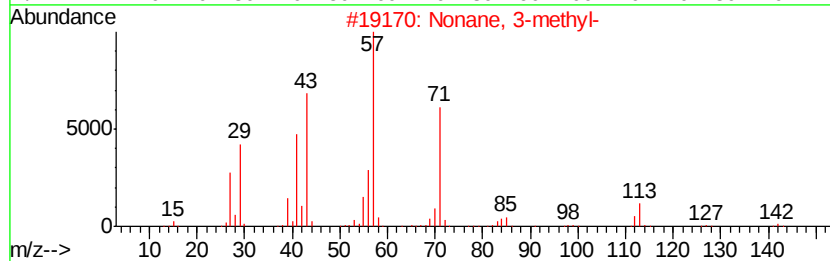
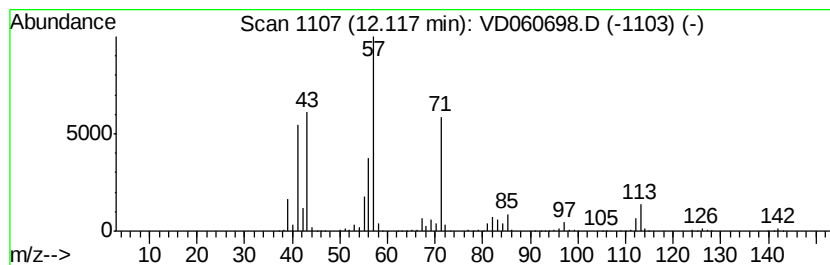
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Nonane, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.12	91.61 ug/l	6201510	Chlorobenzene-d5	11.31		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 3-methyl-	142	C10H22	005911-04-6	93
2		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	90
3		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	80
4		Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	72
5		Oxirane, pentyl-	114	C7H14O	005063-65-0	55



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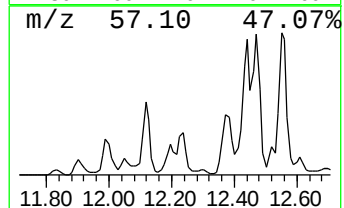
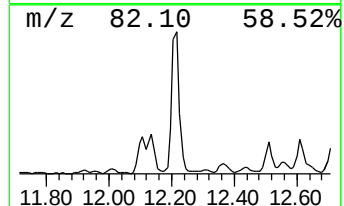
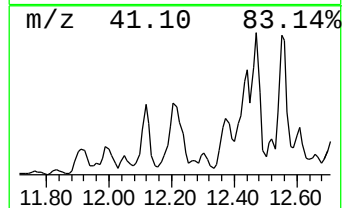
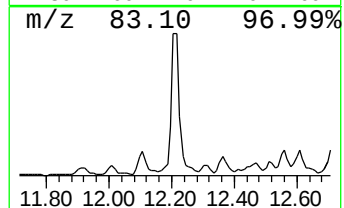
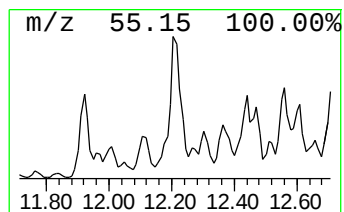
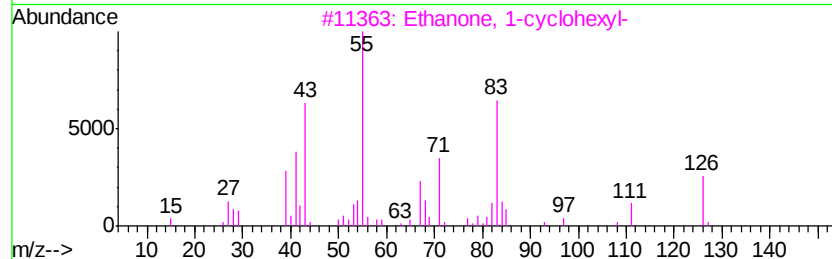
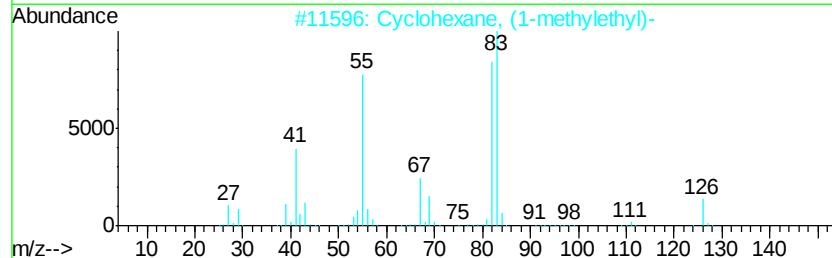
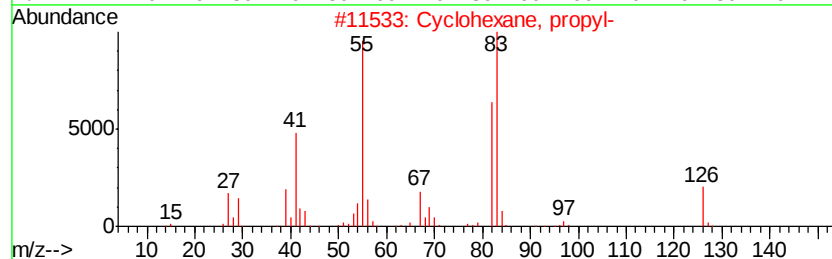
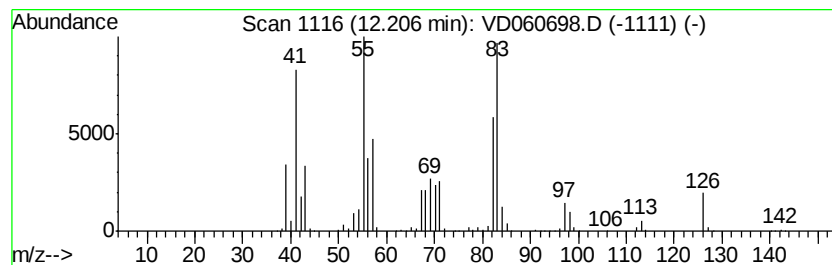
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Quant Title : SW846 8260

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

Peak Number 3 Cyclohexane, propyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.21	172.04 ug/l	11645600	Chlorobenzene-d5	11.31

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, propyl-	126	C9H18	001678-92-8	70
2		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	58
3		Ethanone, 1-cyclohexyl-	126	C8H14O	000823-76-7	43
4		2,3-Dimethyl-3-heptene	126	C9H18	1000113-49-3	38
5		4-Nonene	126	C9H18	002198-23-4	35



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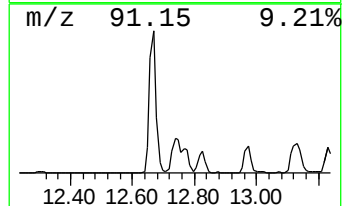
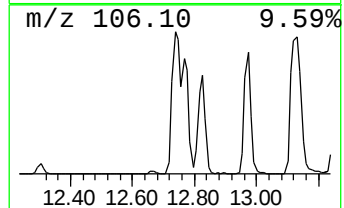
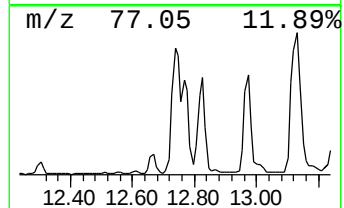
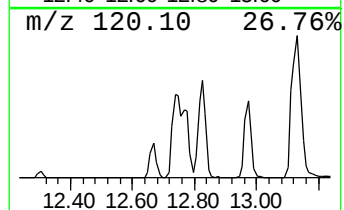
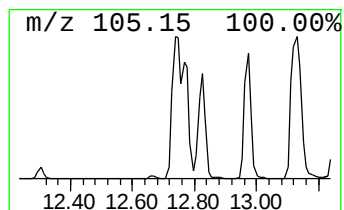
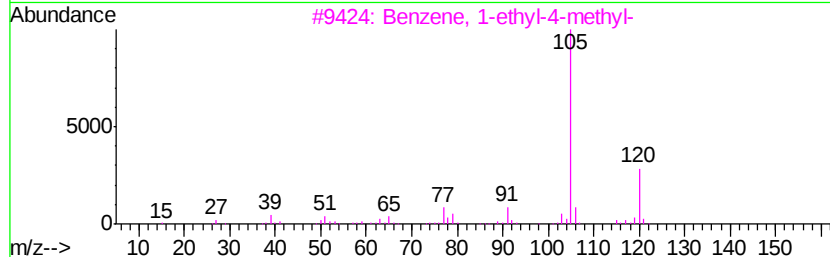
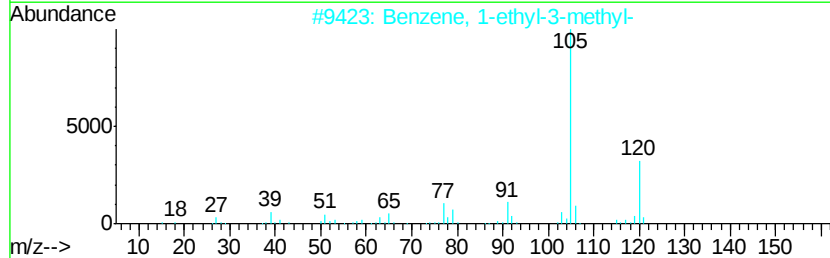
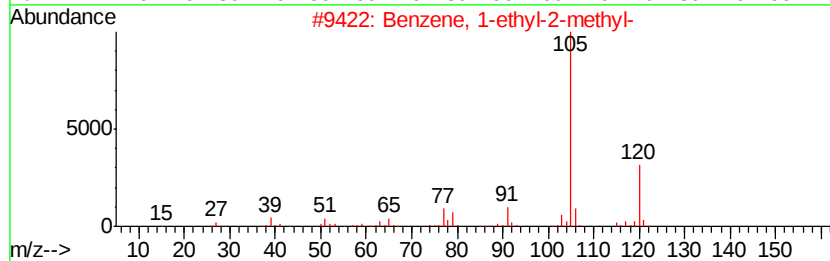
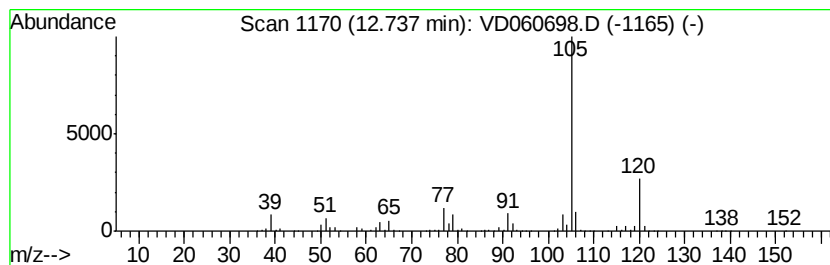
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	140.55 ug/l	25120900	1,4-Dichlorobenzene-d4	13.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD122118\
Data File : VD060698.D
Acq On : 21 Dec 2018 21:56
Operator : VA/AP
Sample : J6469-01
Misc : 5.08g/5ml/MSVOA_D/SOIL
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
DRUM-1

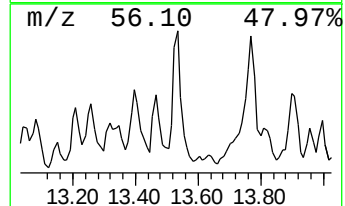
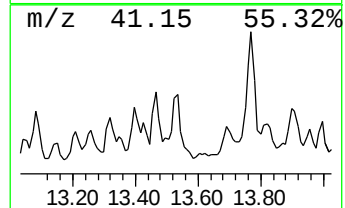
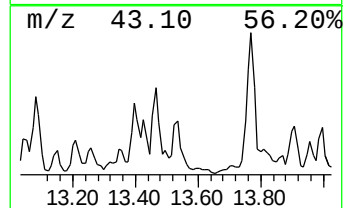
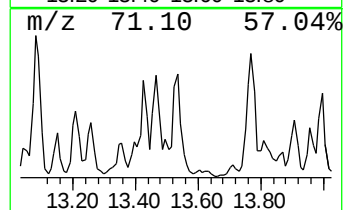
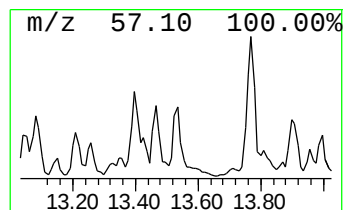
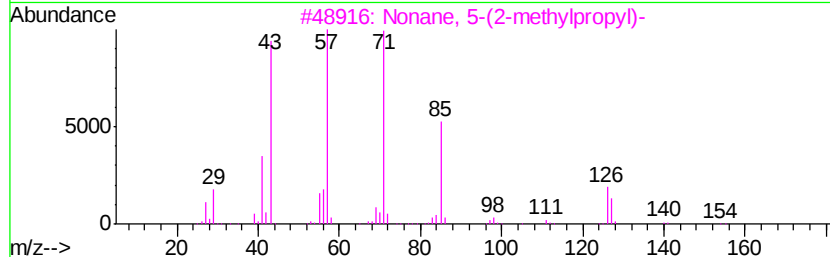
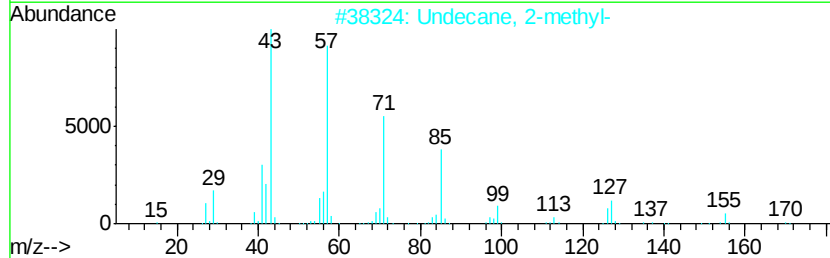
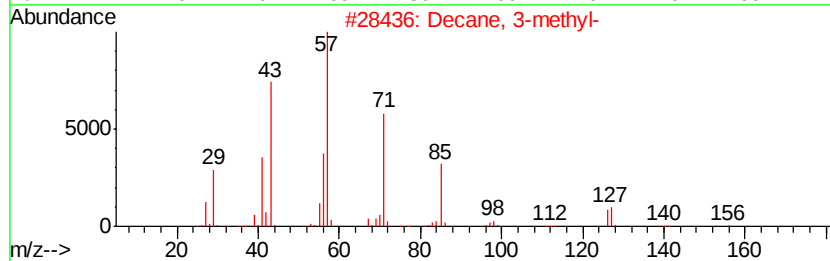
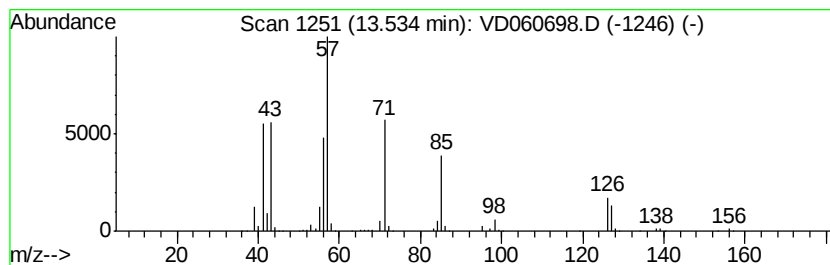
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Quant Title : SW846 8260

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

Peak Number 6 Decane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.53	86.72 ug/l	15499500	1,4-Dichlorobenzene-d4	13.40

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3-methyl-	156	C11H24	013151-34-3	87
2			Undecane, 2-methyl-	170	C12H26	007045-71-8	59
3			Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	53
4			Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	50
5			Dodecane, 3-methyl-	184	C13H28	017312-57-1	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD122118\
Data File : VD060698.D
Acq On : 21 Dec 2018 21:56
Operator : VA/AP
Sample : J6469-01
Misc : 5.08g/5ml/MSVOA_D/SOIL
ALS Vial : 25 Sample Multiplier: 1

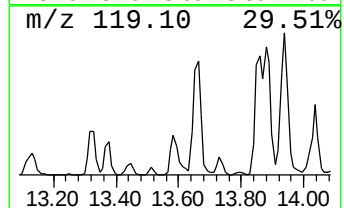
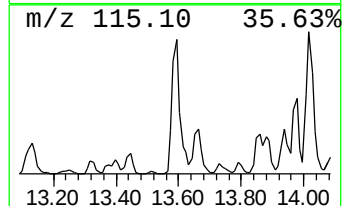
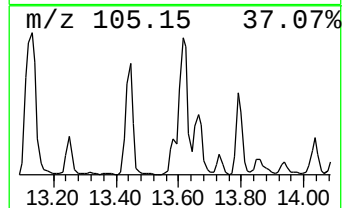
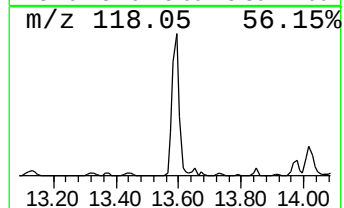
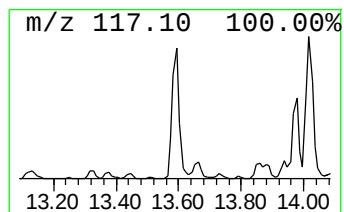
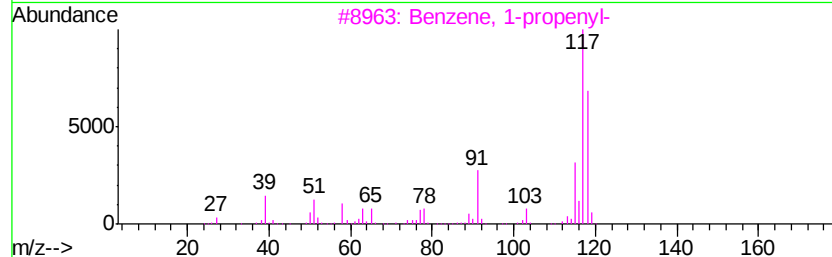
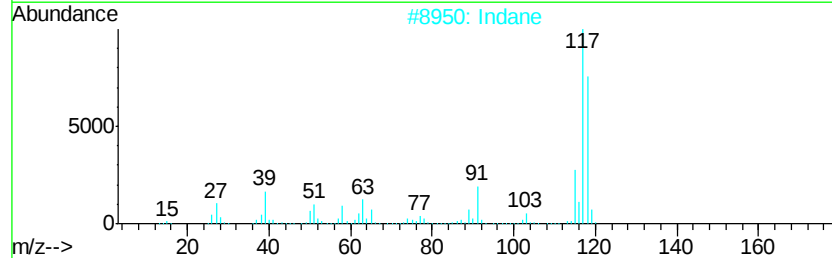
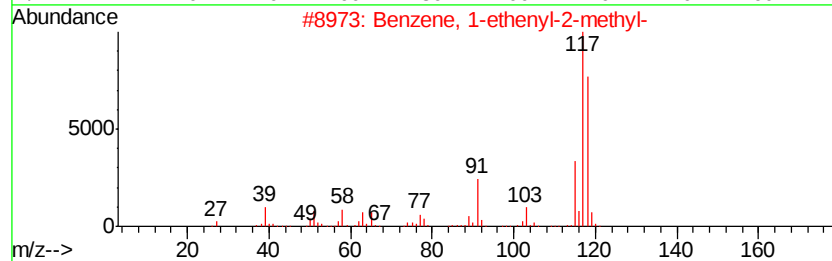
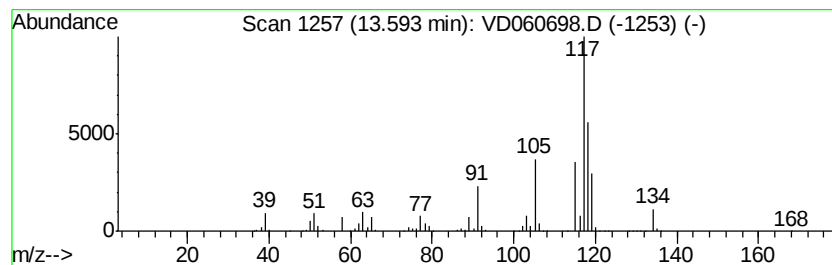
Instrument :
MSVOA_D
ClientSampleId :
DRUM-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D122018S.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethenyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.59	99.13 ug/l	17717000	1,4-Dichlorobenzene-d4	13.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 91
2		Indane	118 C9H10	000496-11-7 64
3		Benzene, 1-propenyl-	118 C9H10	000637-50-3 64
4		Benzene, cyclopropyl-	118 C9H10	000873-49-4 64
5		Benzene, 2-propenyl-	118 C9H10	000300-57-2 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD122118\
Data File : VD060698.D
Acq On : 21 Dec 2018 21:56
Operator : VA/AP
Sample : J6469-01
Misc : 5.08g/5ml/MSVOA_D/SOIL
ALS Vial : 25 Sample Multiplier: 1

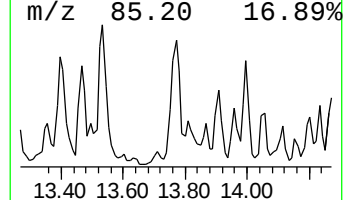
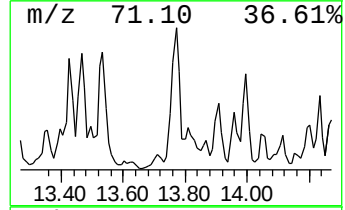
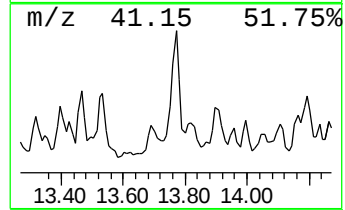
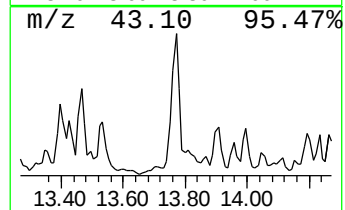
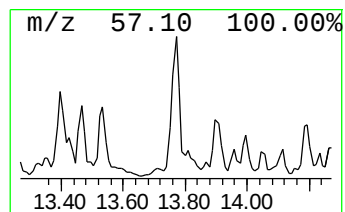
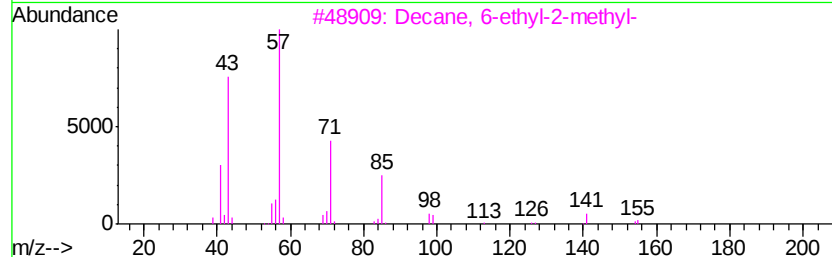
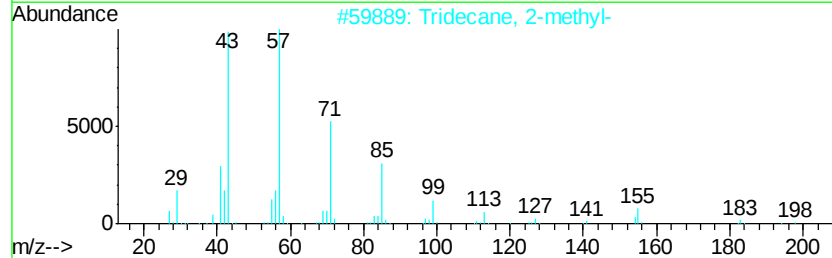
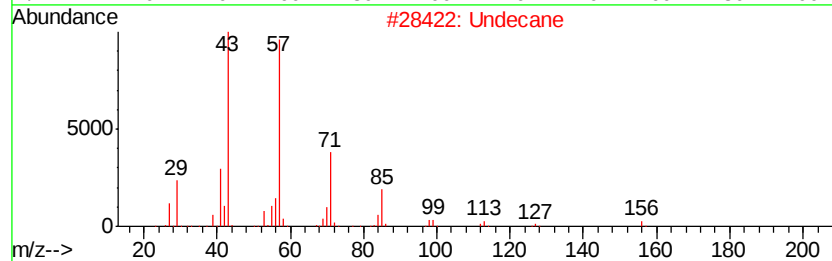
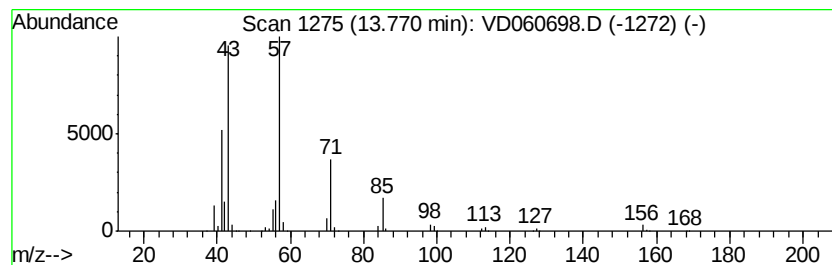
Instrument :
MSVOA_D
ClientSampleId :
DRUM-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D122018S.M
Quant Title : SW846 8260

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

Peak Number 8 Undecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	135.16 ug/l	24156400	1,4-Dichlorobenzene-d4	13.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Undecane		156 C11H24	001120-21-4 91
2	Tridecane, 2-methyl-		198 C14H30	001560-96-9 78
3	Decane, 6-ethyl-2-methyl-		184 C13H28	062108-21-8 72
4	Decane, 2-methyl-		156 C11H24	006975-98-0 64
5	Tridecane		184 C13H28	000629-50-5 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD122118\
Data File : VD060698.D
Acq On : 21 Dec 2018 21:56
Operator : VA/AP
Sample : J6469-01
Misc : 5.08g/5ml/MSVOA_D/SOIL
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
DRUM-1

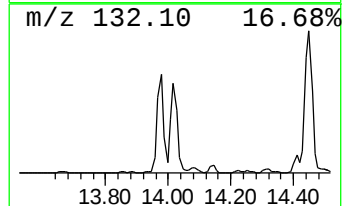
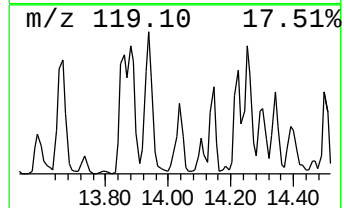
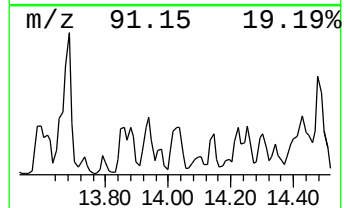
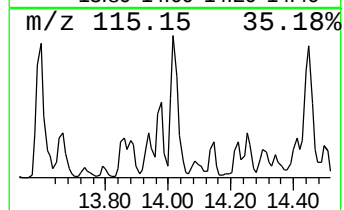
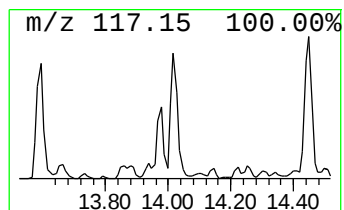
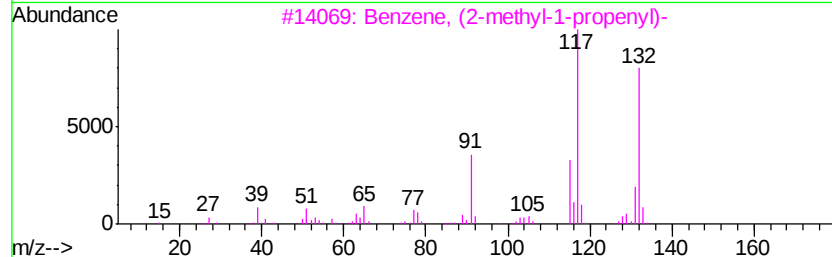
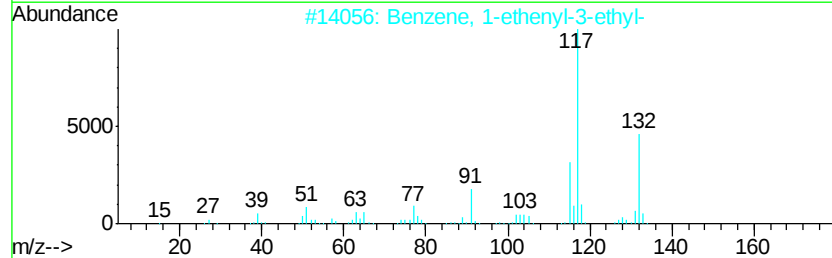
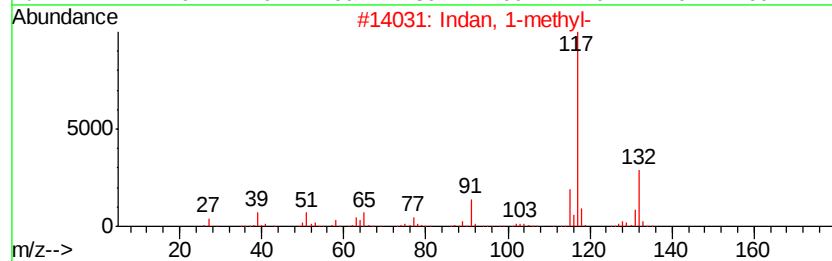
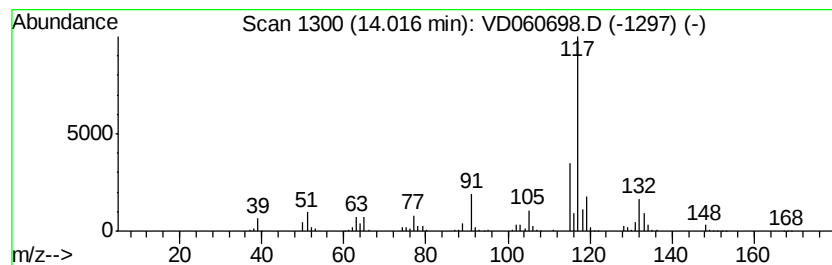
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Quant Title : SW846 8260

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

Peak Number 9 Indan, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.02	83.14 ug/l	14859000	1,4-Dichlorobenzene-d4	13.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indan, 1-methyl-	132	C10H12	000767-58-8	81
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	74
3		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	74
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	72
5		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD122118\
Data File : VD060698.D
Acq On : 21 Dec 2018 21:56
Operator : VA/AP
Sample : J6469-01
Misc : 5.08g/5ml/MSVOA_D/SOIL
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
DRUM-1

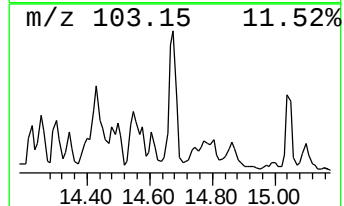
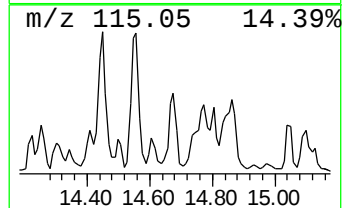
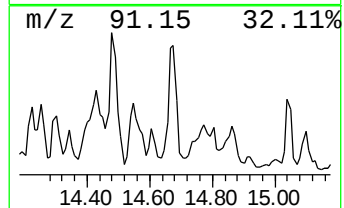
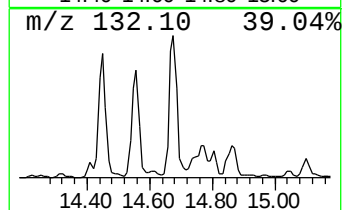
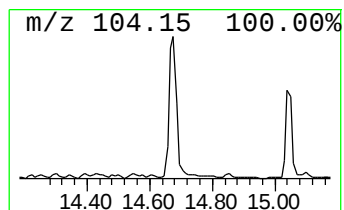
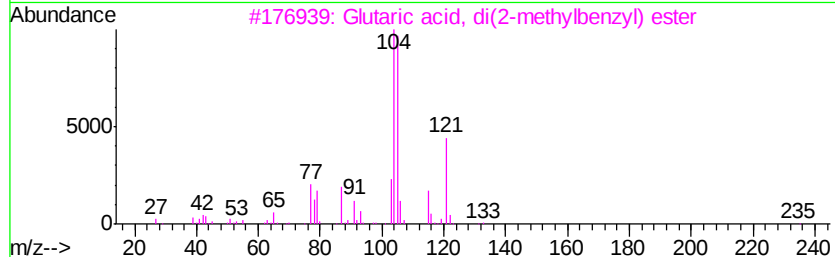
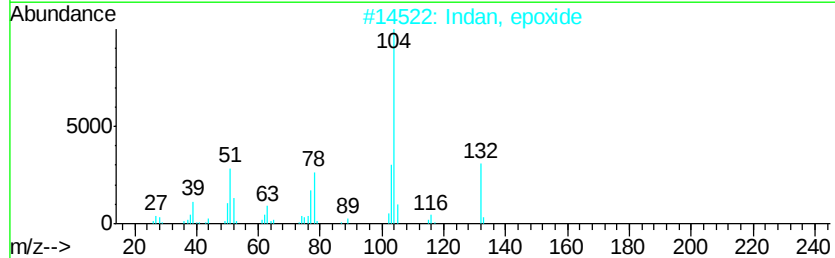
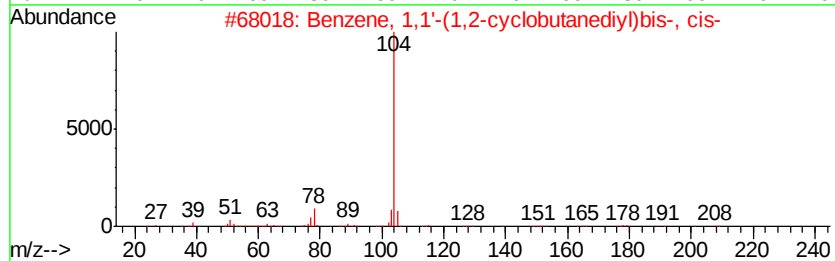
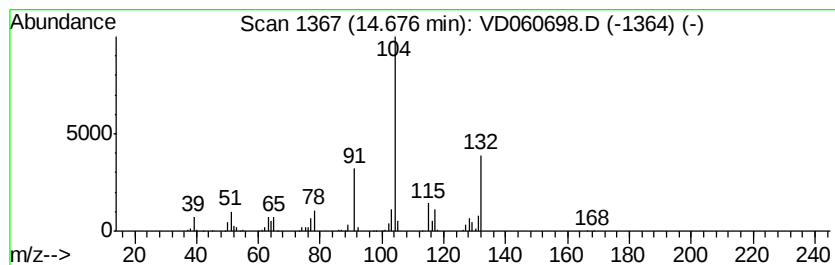
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D122018S.M
Quant Title : SW846 8260

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

Peak Number 10 unknown14.68 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.68	86.98 ug/l	15545300	1,4-Dichlorobenzene-d4	13.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	007694-30-6	38
2		Indan, epoxide	132	C9H8O	000768-22-9	38
3		Glutaric acid, di(2-methylbenzyl...	340	C21H24O4	1000371-34-1	37
4		6-Methyl-3-phenethylsulfanyl-[1,...	247	C12H13N3OS	1000296-66-5	37
5		3(2H)-Furanone, dihydro-2,2-dime...	190	C12H14O2	063678-00-2	37



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD122118\
Data File : VD060698.D
Acq On : 21 Dec 2018 21:56
Operator : VA/AP
Sample : J6469-01
Misc : 5.08g/5ml/MSVOA_D/SOIL
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
DRUM-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D122018S.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
Nonane	11.65	195.1	ug/l	13209400	3	11.31	3384560	50.0
Nonane, 3-methyl-	12.12	91.6	ug/l	6201510	3	11.31	3384560	50.0
Cyclohexane, propyl-	12.21	172.0	ug/l	11645600	3	11.31	3384560	50.0
Benzene, 1-ethyl-...	12.74	140.6	ug/l	25120900	4	13.40	8936430	50.0
Decane, 3-methyl-	13.53	86.7	ug/l	15499500	4	13.40	8936430	50.0
Benzene, 1-etheny...	13.59	99.1	ug/l	17717000	4	13.40	8936430	50.0
Undecane	13.77	135.2	ug/l	24156400	4	13.40	8936430	50.0
Indan, 1-methyl-	14.02	83.1	ug/l	14859000	4	13.40	8936430	50.0
unknown14.68	14.68	87.0	ug/l	15545300	4	13.40	8936430	50.0