

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD071019\
 Data File : VD063066.D
 Acq On : 10 Jul 2019 18:59
 Operator : VA/AP
 Sample : K3726-03
 Misc : 5.48g/5ml/MSVOA D/SOIL
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 PX-3 070919

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\82D061919S.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	3.216	177	182	193	rBV	317546	824618	1.81%	0.145%
2	11.486	1017	1022	1026	rBV	335262	649952	1.43%	0.114%
3	11.633	1031	1037	1039	rBV	261743	604167	1.33%	0.106%
4	11.742	1045	1048	1052	rVB3	208550	482217	1.06%	0.085%
5	11.830	1052	1057	1061	rBV	687372	1476268	3.24%	0.259%
6	12.057	1077	1080	1083	rBV	438517	783295	1.72%	0.137%
7	12.116	1083	1086	1089	rVV	681245	1348874	2.96%	0.236%
8	12.175	1089	1092	1097	rVB2	768202	1760351	3.87%	0.309%
9	12.303	1101	1105	1109	rBV	951644	1914217	4.21%	0.336%
10	12.441	1115	1119	1125	rBV4	157416	481747	1.06%	0.084%
11	12.539	1125	1129	1134	rBV2	466220	1040044	2.28%	0.182%
12	12.657	1134	1141	1143	rBV4	822974	2793438	6.14%	0.490%
13	12.697	1143	1145	1147	rVB	877206	1104886	2.43%	0.194%
14	12.746	1147	1150	1157	rVB2	1638954	3323746	7.30%	0.583%
15	12.854	1157	1161	1165	rBV2	453515	927862	2.04%	0.163%
16	13.002	1172	1176	1179	rBV2	2118241	4954992	10.89%	0.869%
17	13.051	1179	1181	1183	rVB	843700	1056757	2.32%	0.185%
18	13.090	1183	1185	1189	rVB2	1430211	2189807	4.81%	0.384%
19	13.149	1189	1191	1193	rBV2	399635	495047	1.09%	0.087%
20	13.218	1193	1198	1201	rBV	4180893	7025280	15.43%	1.232%
21	13.307	1201	1207	1208	rBV2	3109813	6233600	13.69%	1.093%
22	13.336	1208	1210	1215	rVB	3058102	4439209	9.75%	0.778%
23	13.405	1215	1217	1220	rVB2	1173549	1788855	3.93%	0.314%
24	13.484	1220	1225	1228	rBV3	2846240	7314530	16.07%	1.282%
25	13.553	1228	1232	1233	rBV2	3058867	4821728	10.59%	0.845%
26	13.583	1233	1235	1237	rVB	2609076	3075001	6.76%	0.539%
27	13.622	1237	1239	1241	rVB3	307047	469070	1.03%	0.082%
28	13.661	1241	1243	1246	rBV	3395751	5153809	11.32%	0.903%
29	13.711	1246	1248	1253	rVB2	4128654	6214559	13.65%	1.089%
30	13.848	1256	1262	1264	rBV2	3302488	9167685	20.14%	1.607%
31	13.927	1264	1270	1272	rBV4	7266807	18237294	40.07%	3.197%
32	13.966	1272	1274	1277	rVB	17115505	22978233	50.48%	4.028%
33	14.026	1277	1280	1282	rBV3	1117382	1708016	3.75%	0.299%
34	14.085	1282	1286	1287	rBV2	3384949	6583907	14.46%	1.154%

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Integrator: RTE
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35	14.114	1287	1289	1292	rVB2	3177797	5060592	11.12%	0.887%
36	14.163	1292	1294	1296	rBV	2046202	3518762	7.73%	0.617%
37	14.203	1296	1298	1300	rVV	11390877	14911677	32.76%	2.614%
38	14.242	1300	1302	1308	rVB2	7512155	16889388	37.10%	2.961%
39	14.331	1308	1311	1313	rBV	3639262	6163438	13.54%	1.080%
40	14.380	1313	1316	1319	rVB5	4943051	8066018	17.72%	1.414%
41	14.439	1319	1322	1324	rBV2	10530988	14146593	31.08%	2.480%
42	14.478	1324	1326	1328	rVB3	2452129	3026363	6.65%	0.531%
43	14.528	1328	1331	1332	rBV	6022463	8436460	18.53%	1.479%
44	14.557	1332	1334	1336	rVV	11726972	15641202	34.36%	2.742%
45	14.587	1336	1337	1339	rVB	6955771	6403457	14.07%	1.123%
46	14.656	1339	1344	1347	rVB2	9373742	13103912	28.79%	2.297%
47	14.705	1347	1349	1350	rBV2	3200906	3808101	8.37%	0.668%
48	14.734	1350	1352	1354	rVB	6968430	8315321	18.27%	1.458%
49	14.784	1354	1357	1358	rBV	10479573	13980534	30.71%	2.451%
50	14.813	1358	1360	1363	rVV3	6493452	10297721	22.62%	1.805%
51	14.892	1365	1368	1372	rVV2	25288616	45518827	100.00%	7.979%
52	14.941	1372	1373	1375	rVV2	5435265	7631152	16.76%	1.338%
53	14.980	1375	1377	1378	rVV	5315170	6253205	13.74%	1.096%
54	15.030	1378	1382	1383	rVV2	8914535	20989873	46.11%	3.680%
55	15.059	1383	1385	1387	rVV	10508698	14595240	32.06%	2.559%
56	15.138	1390	1393	1398	rBV4	10535172	32765044	71.98%	5.744%
57	15.227	1398	1402	1405	rVB3	6658333	13269079	29.15%	2.326%
58	15.286	1405	1408	1410	rBV3	8989666	15999892	35.15%	2.805%
59	15.325	1410	1412	1414	rVV	10325417	17164858	37.71%	3.009%
60	15.384	1416	1418	1420	rVB2	11203268	13884954	30.50%	2.434%
61	15.433	1420	1423	1426	rBV4	10885752	16331530	35.88%	2.863%
62	15.473	1426	1427	1431	rVB3	3857645	4399692	9.67%	0.771%
63	15.532	1431	1433	1434	rBV	2396516	2325403	5.11%	0.408%
64	15.650	1441	1445	1446	rBV2	5883754	13389700	29.42%	2.347%
65	15.670	1446	1447	1452	rVV2	13658885	22133556	48.63%	3.880%
66	15.738	1452	1454	1457	rVB2	4132676	5703777	12.53%	1.000%
67	15.798	1457	1460	1463	rVB3	5892359	8971859	19.71%	1.573%
68	15.896	1465	1470	1472	rBV	2459222	5371889	11.80%	0.942%
69	15.935	1472	1474	1476	rVB	2630812	2804372	6.16%	0.492%
70	15.994	1476	1480	1482	rVB2	3132070	5399745	11.86%	0.947%
71	16.044	1483	1485	1490	rVB4	1287420	2583430	5.68%	0.453%

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Integration Parameters: RTEINT.P

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Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Title : SW846 8260

72	16.152	1493	1496	1497	rBV3	572533	990098	2.18%	0.174%
73	16.172	1497	1498	1500	rVB	1781730	1588807	3.49%	0.279%
74	16.231	1500	1504	1509	rVB3	2099124	4421527	9.71%	0.775%
75	16.496	1530	1531	1537	rVB	498606	769414	1.69%	0.135%

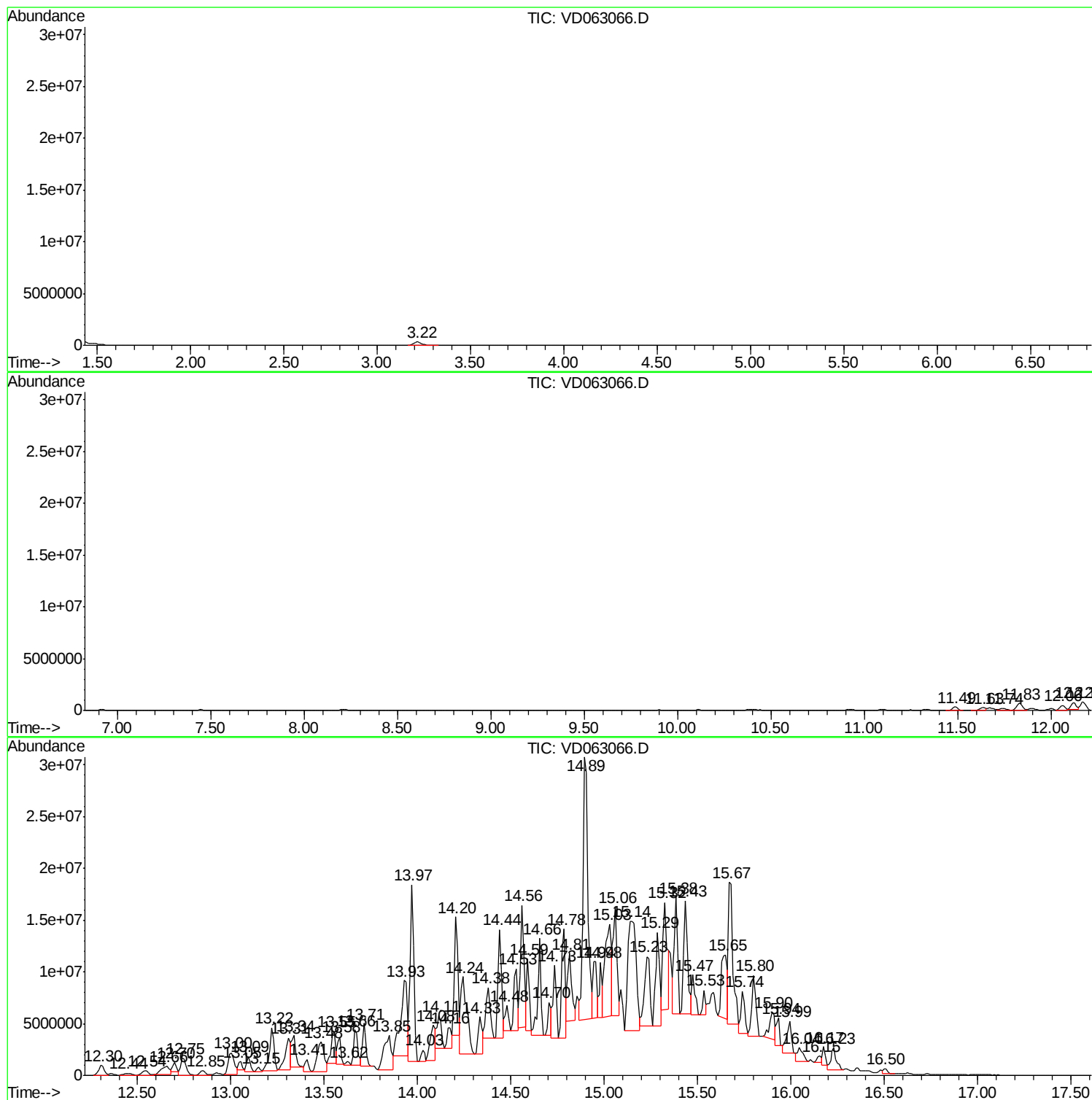
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061919S.M
Quant Title : SW846 8260

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



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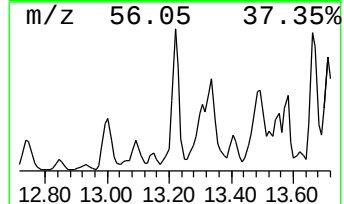
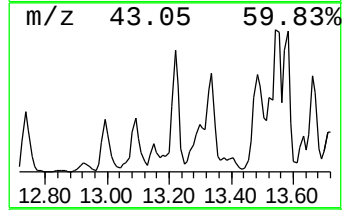
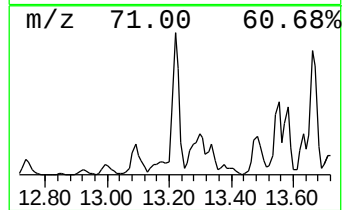
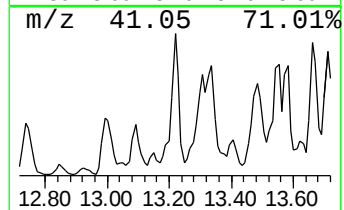
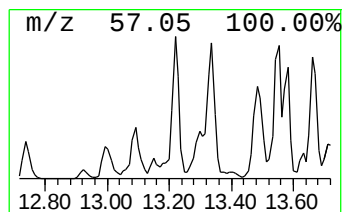
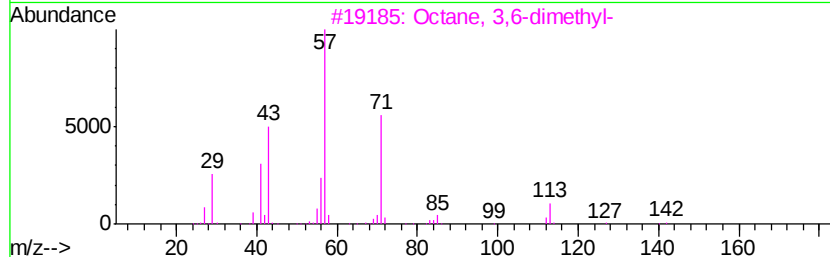
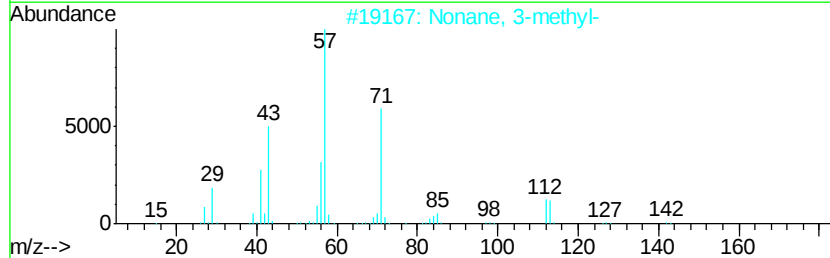
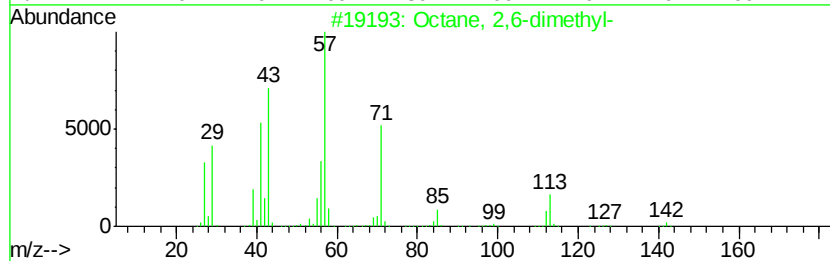
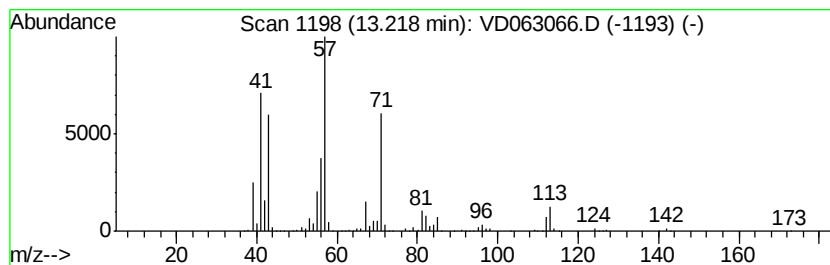
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061919S.M
Quant Title : SW846 8260

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

Peak Number 1 Octane, 2,6-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	183.50 ug/l	7025280	Chlorobenzene-d5	12.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	87
2		Nonane, 3-methyl-	142	C10H22	005911-04-6	81
3		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	74
4		Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	53
5		Octane, 1,1'-oxybis-	242	C16H34O	000629-82-3	53



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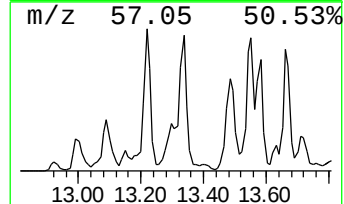
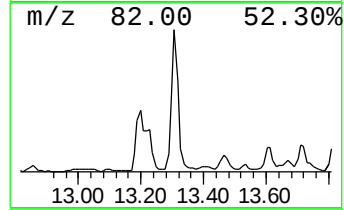
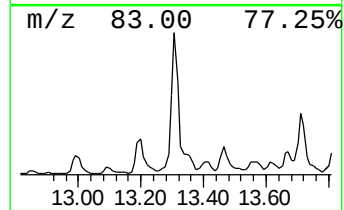
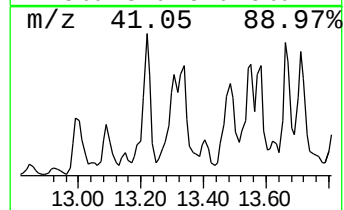
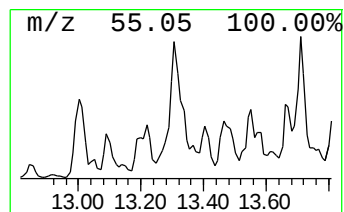
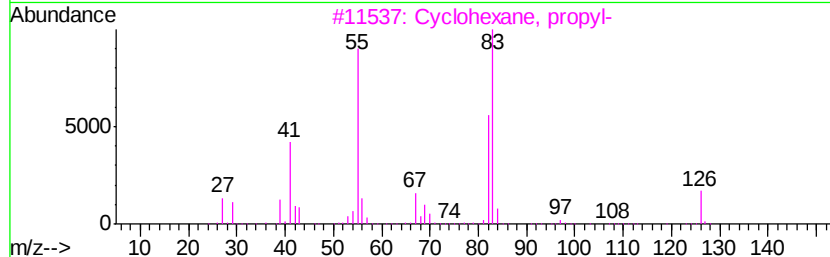
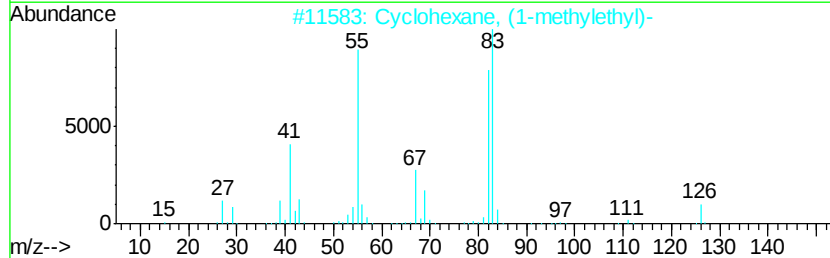
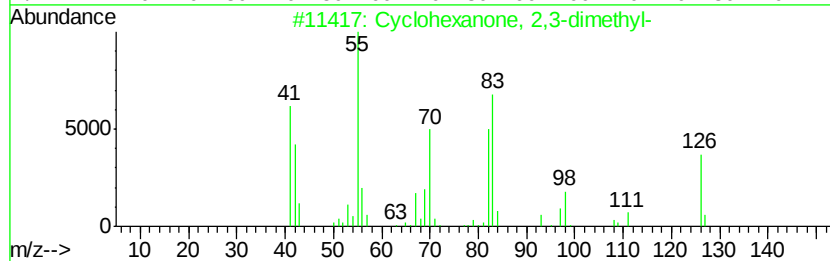
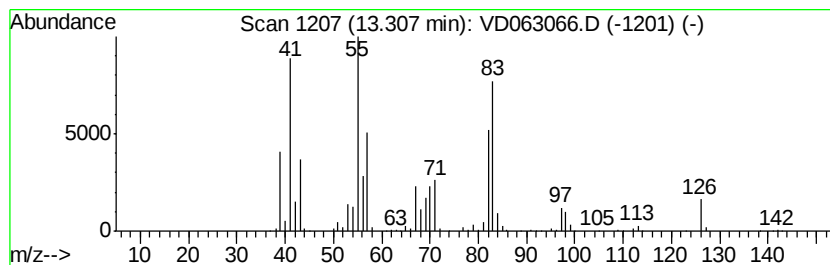
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061919S.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclohexanone, 2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.31	162.82 ug/l	6233600	Chlorobenzene-d5	12.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanone, 2,3-dimethyl-	126	C8H14O	013395-76-1	72
2		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	70
3		Cyclohexane, propyl-	126	C9H18	001678-92-8	58
4		Cyclohexane, methyl-	98	C7H14	000108-87-2	53
5		Cyclohexanol, 1-ethenyl-	126	C8H14O	001940-19-8	46



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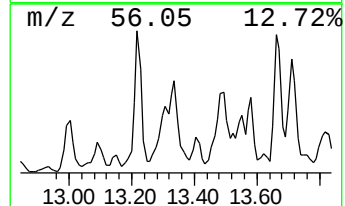
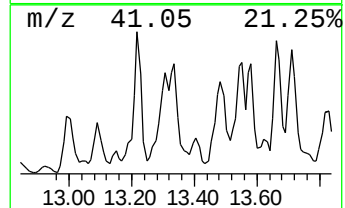
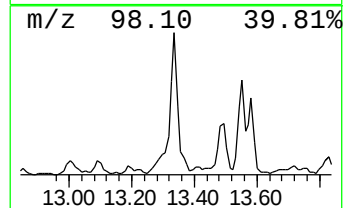
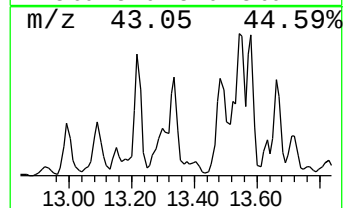
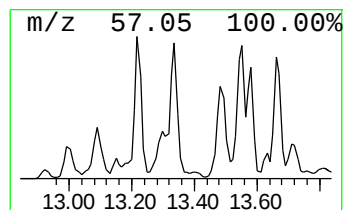
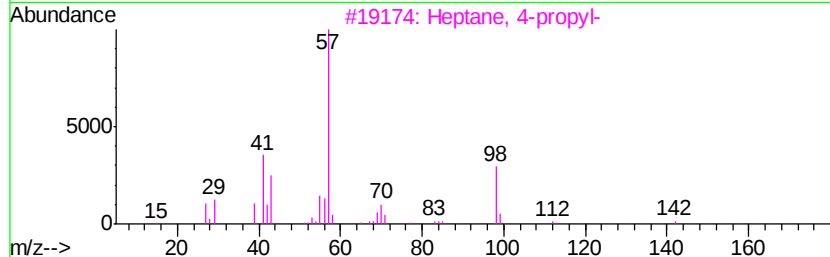
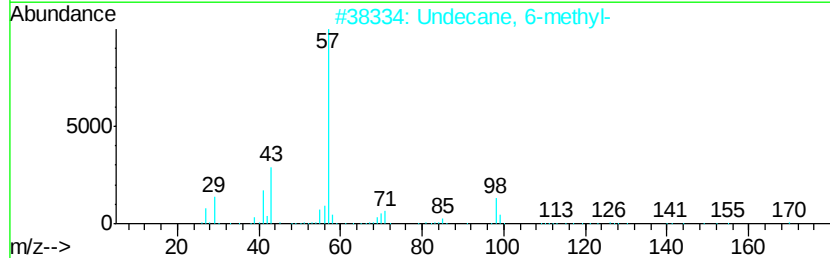
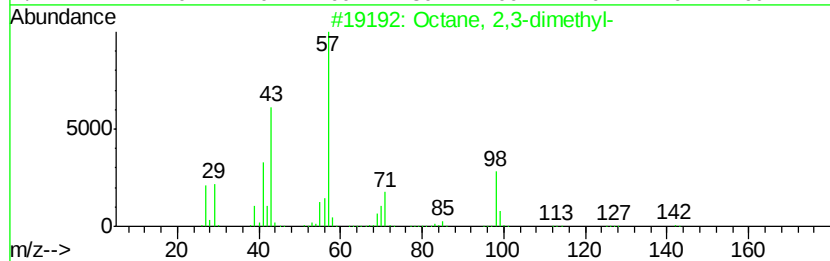
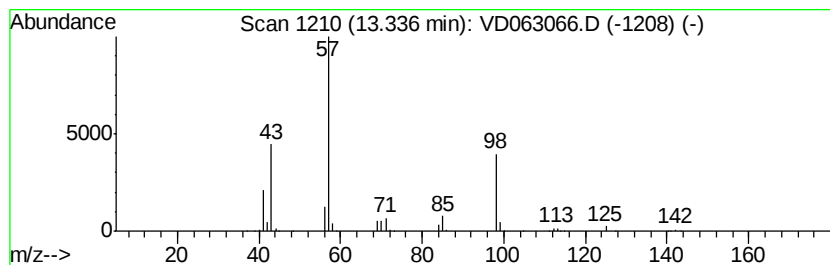
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061919S.M
Quant Title : SW846 8260

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

Peak Number 3 Octane, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	115.95 ug/l	4439210	Chlorobenzene-d5	12.36

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	59
2		Undecane, 6-methyl-	170	C12H26	017302-33-9	59
3		Heptane, 4-propyl-	142	C10H22	003178-29-8	50
4		Dodecane, 6-methyl-	184	C13H28	006044-71-9	37
5		Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	37



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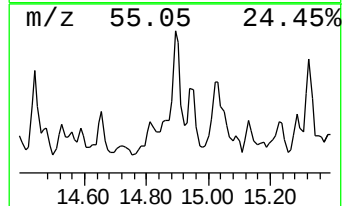
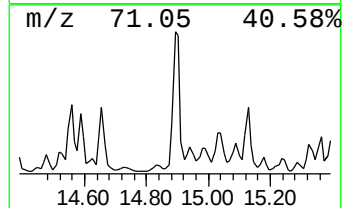
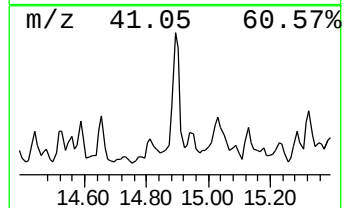
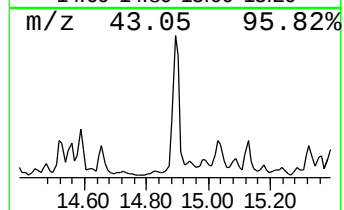
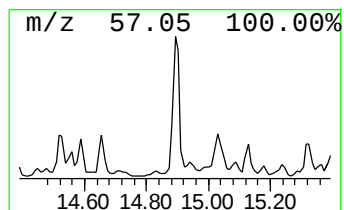
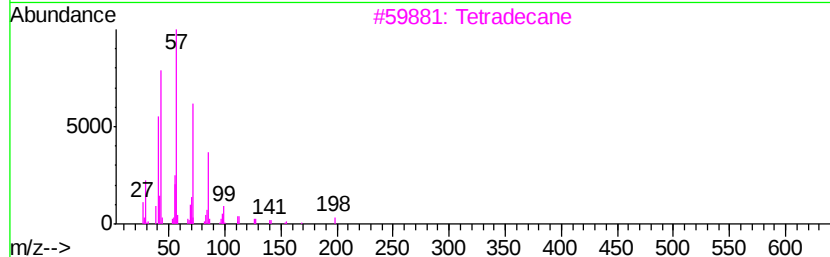
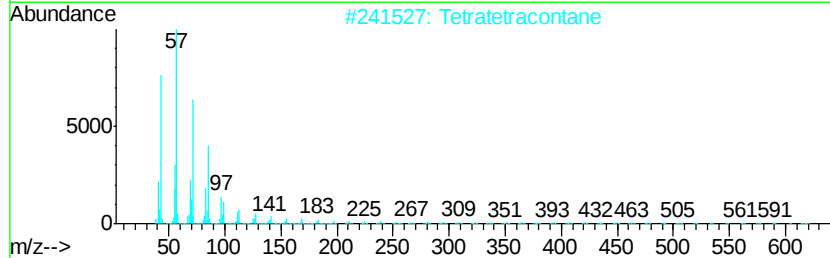
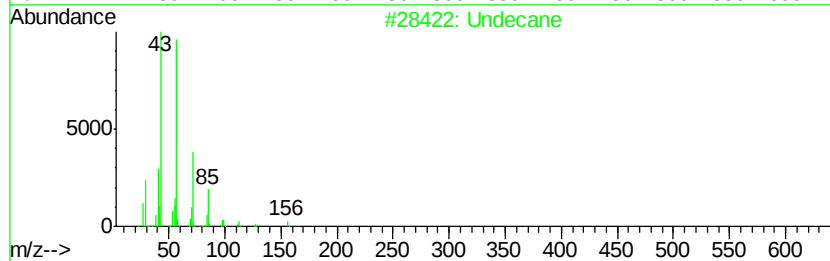
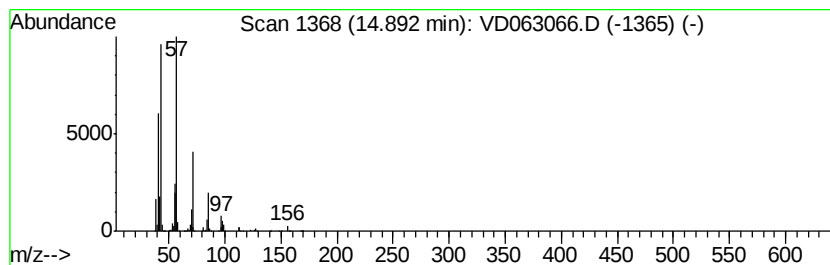
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Quant Title : SW846 8260

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

Peak Number 4 Undecane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.89	269.77 ug/l	45518800	1,4-Dichlorobenzene-d4	14.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Undecane	156	C11H24	001120-21-4	90
2		Tetratetracontane	619	C44H90	007098-22-8	78
3		Tetradecane	198	C14H30	000629-59-4	78
4		Undecane, 4,7-dimethyl-	184	C13H28	017301-32-5	72
5		Octane, 2,7-dimethyl-	142	C10H22	001072-16-8	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD071019\
Data File : VD063066.D
Acq On : 10 Jul 2019 18:59
Operator : VA/AP
Sample : K3726-03
Misc : 5.48g/5ml/MSVOA D/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
PX-3 070919

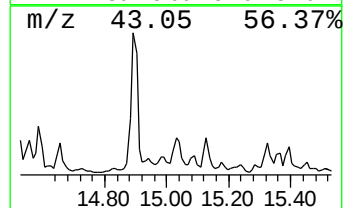
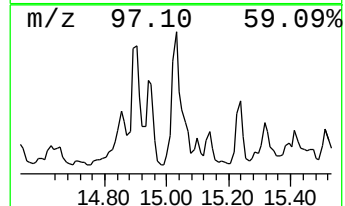
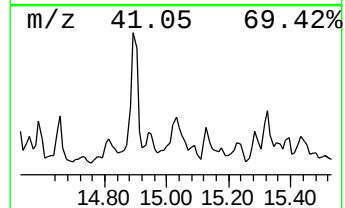
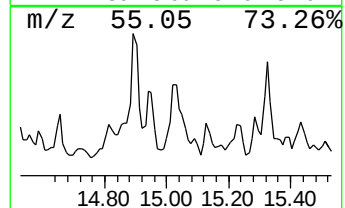
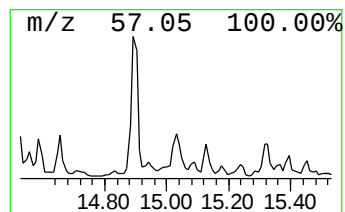
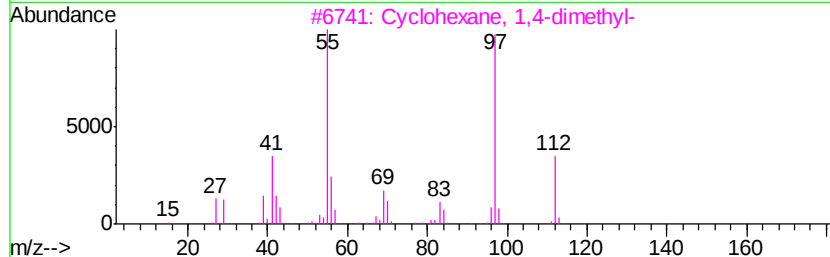
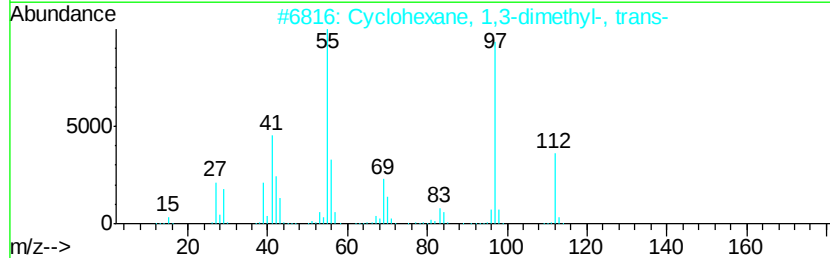
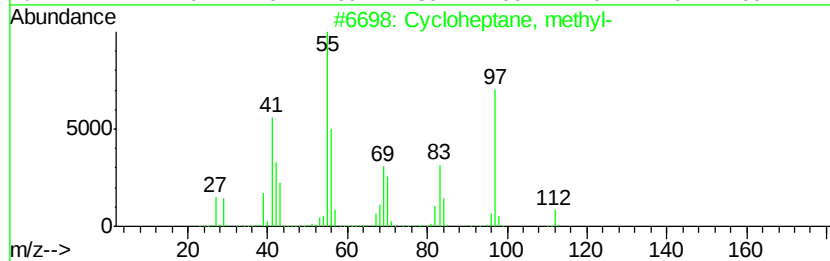
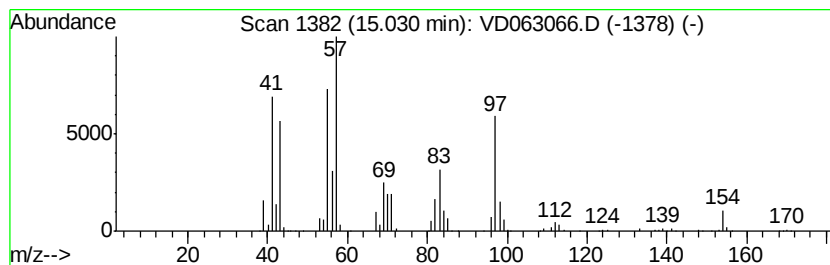
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Quant Title : SW846 8260

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

Peak Number 5 Cycloheptane, methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.03	124.40 ug/l	20989900	1,4-Dichlorobenzene-d4	14.52

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cycloheptane, methyl-	112	C8H16	004126-78-7	64
2		Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	49
3		Cyclohexane, 1,4-dimethyl-	112	C8H16	000589-90-2	46
4		Cyclopropane, 1,1,2,3-tetramethyl-	98	C7H14	074752-93-5	45
5		Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD071019\
Data File : VD063066.D
Acq On : 10 Jul 2019 18:59
Operator : VA/AP
Sample : K3726-03
Misc : 5.48g/5ml/MSVOA D/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
PX-3 070919

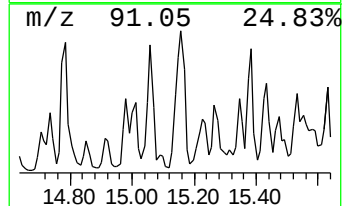
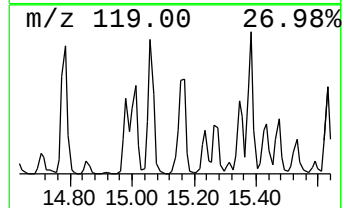
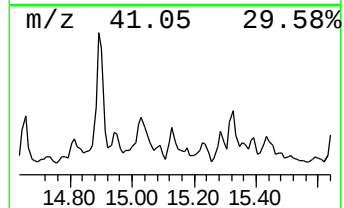
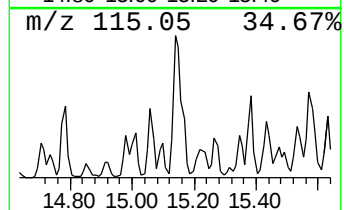
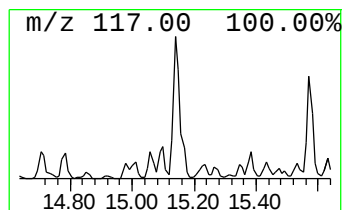
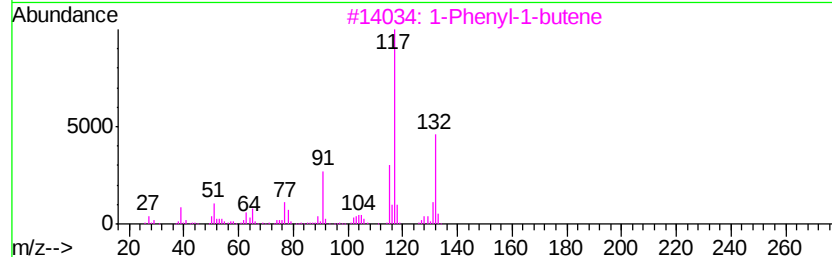
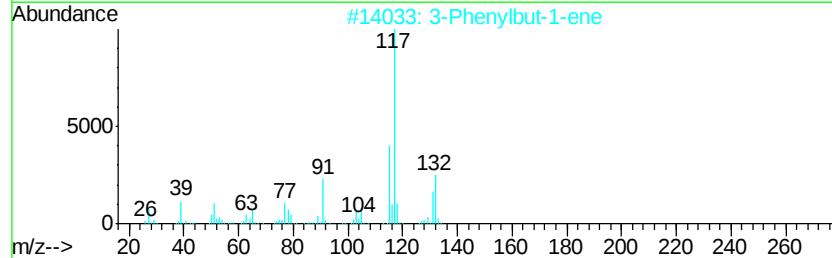
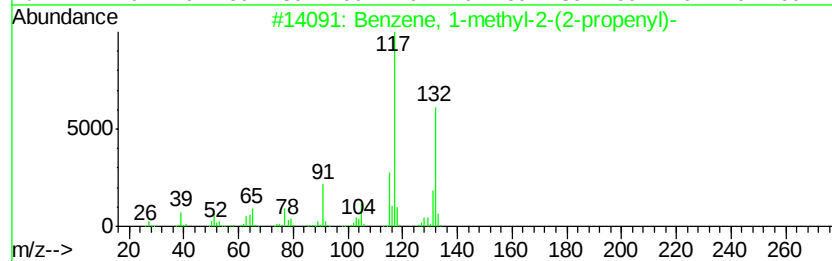
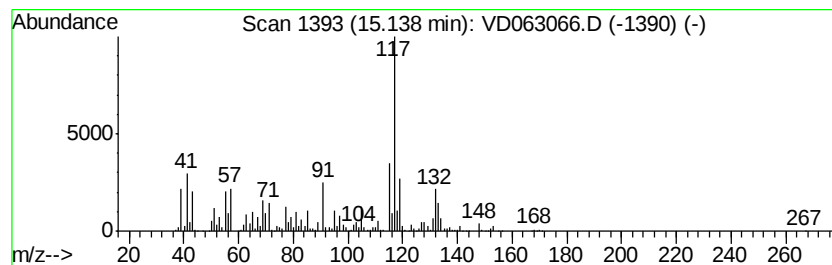
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-methyl-2-(2-prop... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.14	194.19 ug/l	32765000	1,4-Dichlorobenzene-d4	14.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	70
4		Indan, 1-methyl-	132	C10H12	000767-58-8	64
5		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD071019\
Data File : VD063066.D
Acq On : 10 Jul 2019 18:59
Operator : VA/AP
Sample : K3726-03
Misc : 5.48g/5ml/MSVOA D/SOIL
ALS Vial : 12 Sample Multiplier: 1

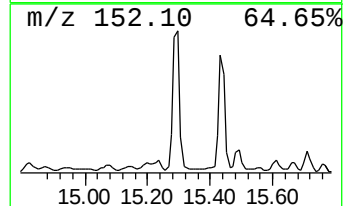
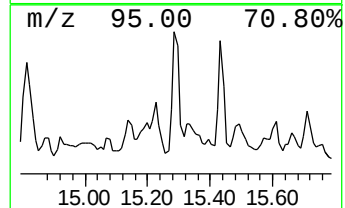
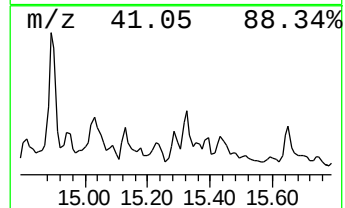
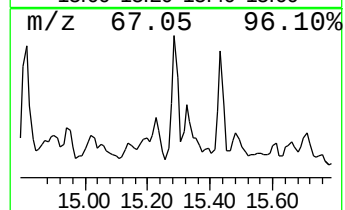
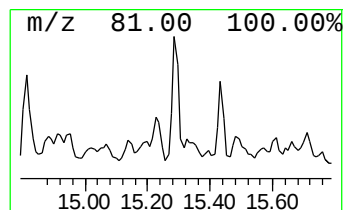
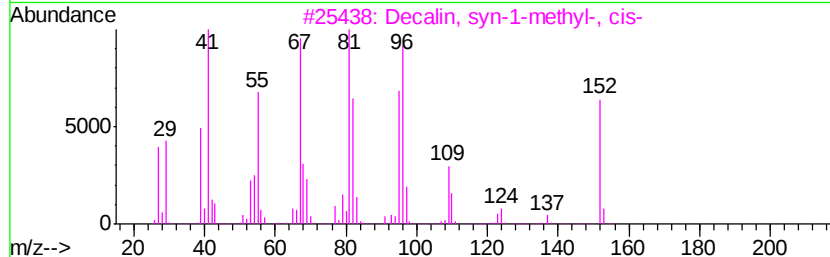
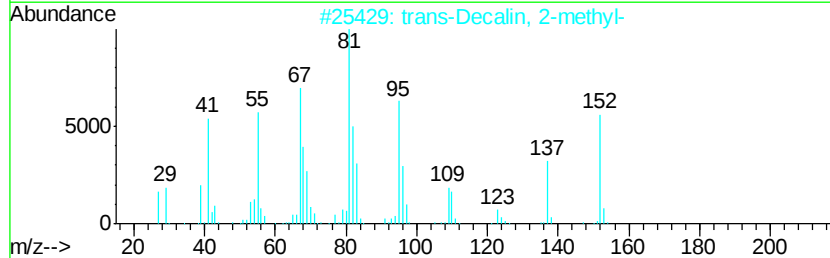
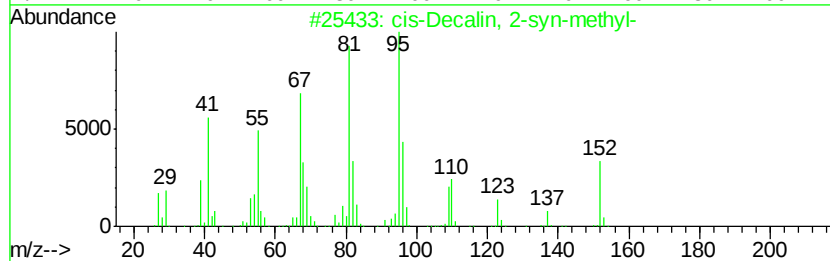
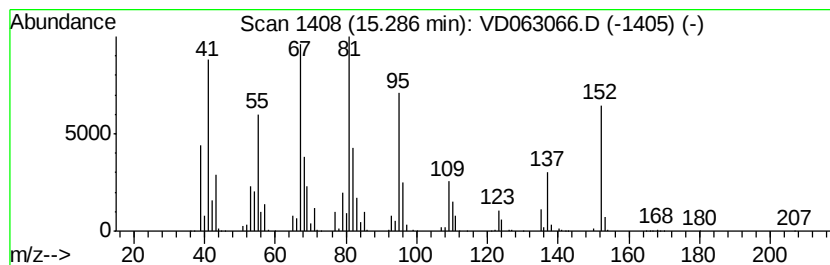
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ClientSampled :
PX-3 070919

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061919S.M
Quant Title : SW846 8260

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

Peak Number 7 cis-Decalin, 2-syn-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.29	94.83 ug/l	15999900	1,4-Dichlorobenzene-d4	14.52
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		cis-Decalin, 2-syn-methyl-	152 C11H20	1000155-85-6 91
2		trans-Decalin, 2-methyl-	152 C11H20	1000152-47-3 90
3		Decalin, syn-1-methyl-, cis-	152 C11H20	1000158-89-1 86
4		Spiro[5.5]undecane	152 C11H20	000180-43-8 86
5		7-Hexadecyne	222 C16H30	074685-28-2 72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD071019\
Data File : VD063066.D
Acq On : 10 Jul 2019 18:59
Operator : VA/AP
Sample : K3726-03
Misc : 5.48g/5ml/MSVOA D/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
PX-3 070919

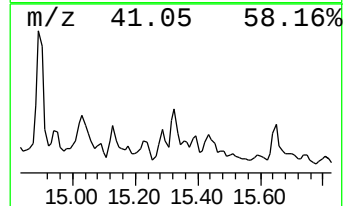
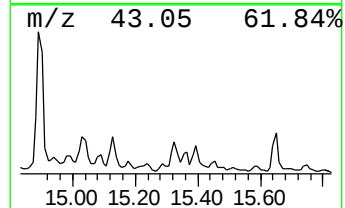
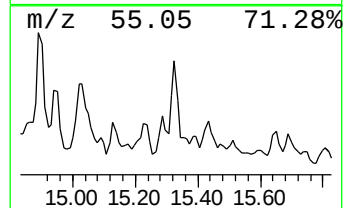
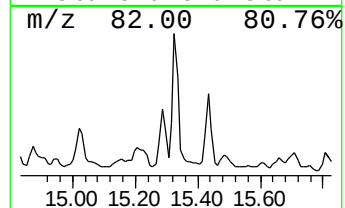
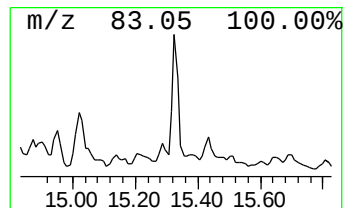
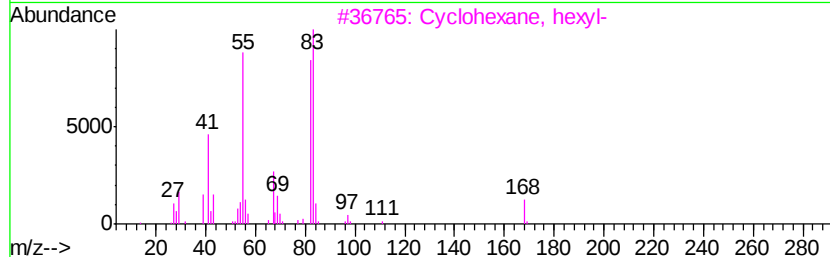
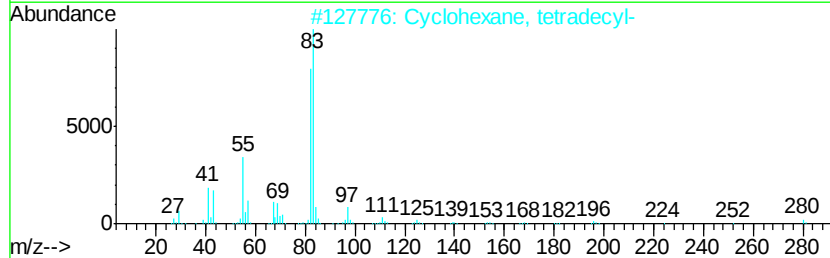
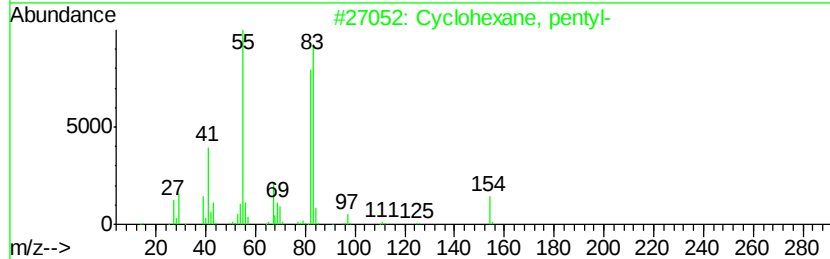
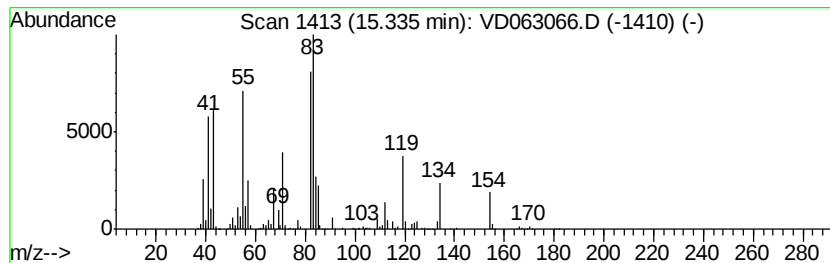
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Quant Title : SW846 8260

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

Peak Number 8 Cyclohexane, pentyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.33	101.73 ug/l	17164900	1,4-Dichlorobenzene-d4	14.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, pentyl-	154	C11H22	004292-92-6	70
2		Cyclohexane, tetradecyl-	280	C20H40	001795-18-2	53
3		Cyclohexane, hexyl-	168	C12H24	004292-75-5	50
4		Cyclohexane, butyl-	140	C10H20	001678-93-9	50
5		Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD071019\
Data File : VD063066.D
Acq On : 10 Jul 2019 18:59
Operator : VA/AP
Sample : K3726-03
Misc : 5.48g/5ml/MSVOA D/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampled :
PX-3 070919

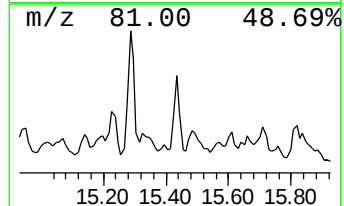
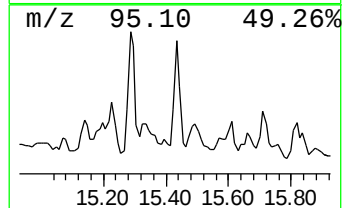
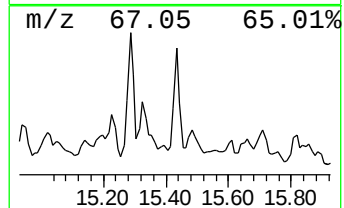
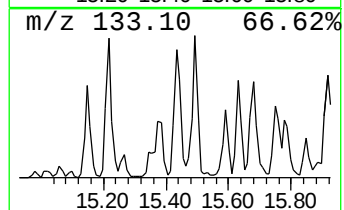
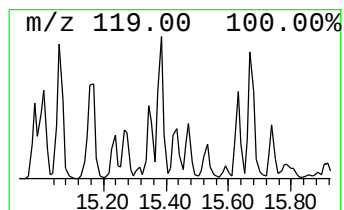
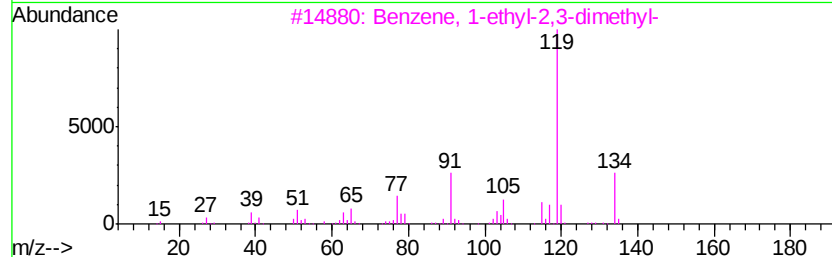
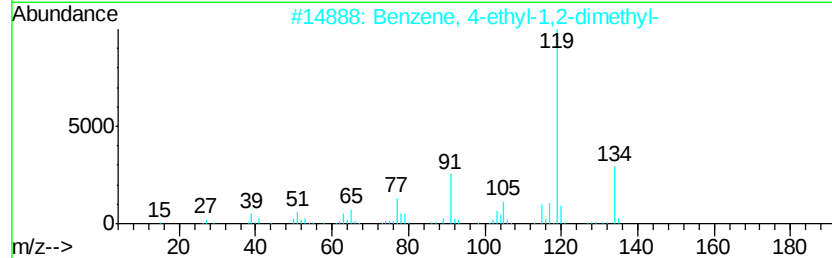
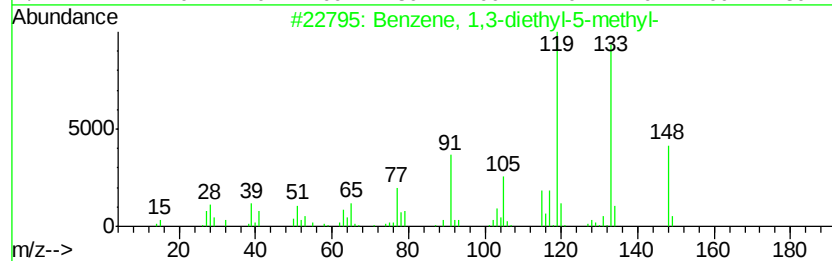
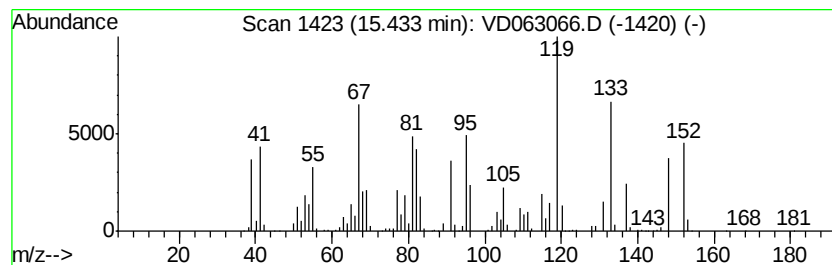
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,3-diethyl-5-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.43	96.79 ug/l	16331500	1,4-Dichlorobenzene-d4	14.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	59
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	25
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	25
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	25
5		Spiro[5.5]undecane	152	C11H20	000180-43-8	18



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD071019\
Data File : VD063066.D
Acq On : 10 Jul 2019 18:59
Operator : VA/AP
Sample : K3726-03
Misc : 5.48g/5ml/MSVOA D/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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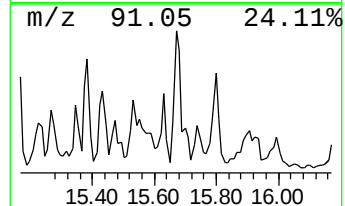
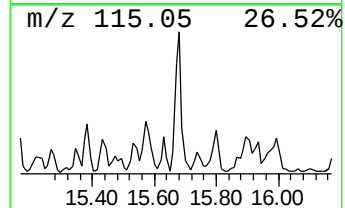
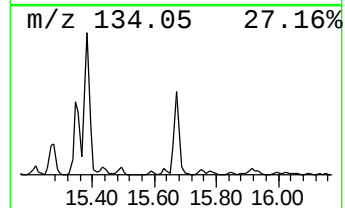
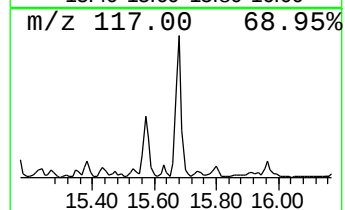
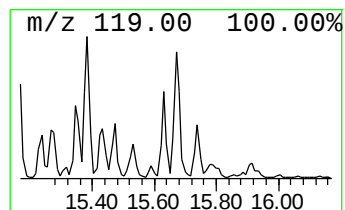
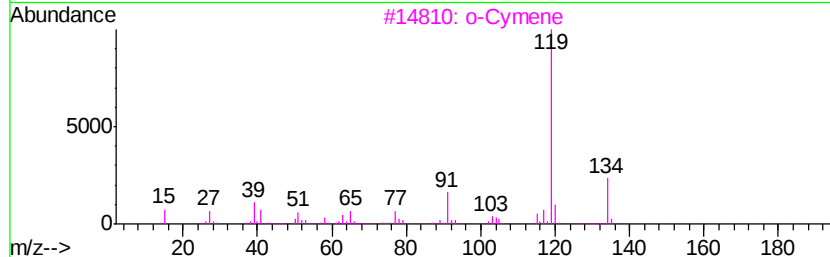
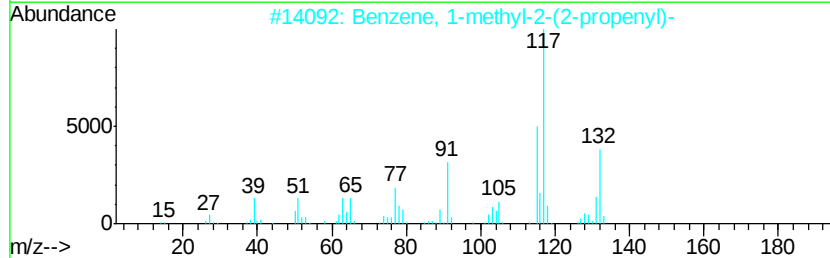
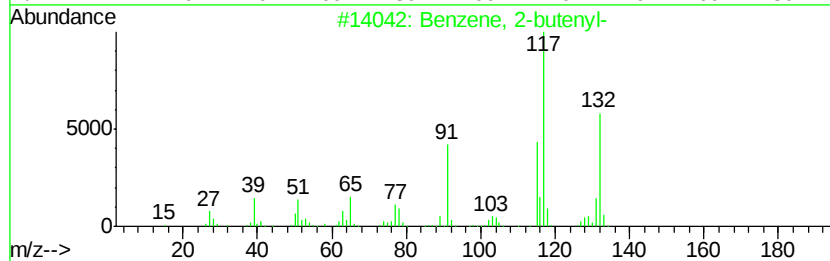
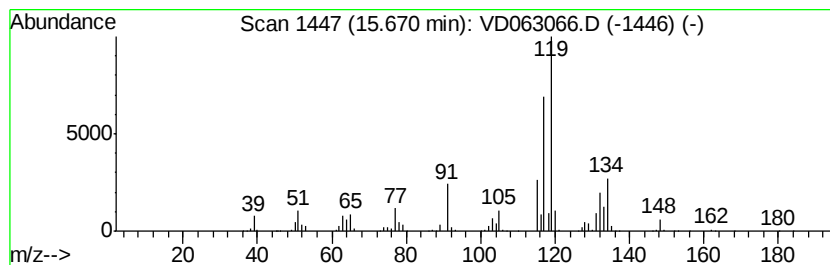
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-butenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.67	131.18 ug/l	22133600	1,4-Dichlorobenzene-d4	14.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	86
2		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	84
3		o-Cymene	134	C10H14	000527-84-4	55
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	55
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD071019\
Data File : VD063066.D
Acq On : 10 Jul 2019 18:59
Operator : VA/AP
Sample : K3726-03
Misc : 5.48g/5ml/MSVOA_D/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
PX-3_070919

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061919S.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
Octane, 2,6-dimet...	13.22	183.5	ug/l	7025280	3	12.36	1914220	50.0
Cyclohexanone, 2,...	13.31	162.8	ug/l	6233600	3	12.36	1914220	50.0
Octane, 2,3-dimet...	13.34	116.0	ug/l	4439210	3	12.36	1914220	50.0
Undecane	14.89	269.8	ug/l	45518800	4	14.52	8436460	50.0
Cycloheptane, met...	15.03	124.4	ug/l	20989900	4	14.52	8436460	50.0
Benzene, 1-methyl...	15.14	194.2	ug/l	32765000	4	14.52	8436460	50.0
cis-Decalin, 2-sy...	15.29	94.8	ug/l	15999900	4	14.52	8436460	50.0
Cyclohexane, pentyl-	15.33	101.7	ug/l	17164900	4	14.52	8436460	50.0
Benzene, 1,3-diet...	15.43	96.8	ug/l	16331500	4	14.52	8436460	50.0
Benzene, 2-butenyl-	15.67	131.2	ug/l	22133600	4	14.52	8436460	50.0