

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD102418\
 Data File : VD060180.D
 Acq On : 24 Oct 2018 14:25
 Operator : JC/SY
 Sample : J5622-03
 Misc : 4.91g/5ml/MSVOA D/SOIL
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 KG-V23608

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\82D100118S.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	3.161	159	166	172	rBV	36350	125034	1.84%	0.089%
2	3.722	218	223	245	rVB	297764	903818	13.32%	0.644%
3	5.651	411	419	429	rBV2	55312	213935	3.15%	0.153%
4	6.320	481	487	492	rBV	845141	2272326	33.50%	1.620%
5	6.409	492	496	513	rVB	1099387	2918126	43.02%	2.081%
6	6.783	528	534	547	rBV	663555	1728302	25.48%	1.232%
7	7.442	595	601	618	rBV	1663437	3962775	58.42%	2.826%
8	8.032	653	661	670	rBV5	28959	111985	1.65%	0.080%
9	9.450	796	805	812	rBV	1602324	3757165	55.39%	2.679%
10	9.548	812	815	821	rVB	66762	145579	2.15%	0.104%
11	9.784	836	839	848	rVB	44241	108663	1.60%	0.077%
12	9.951	848	856	863	rVB2	73223	212544	3.13%	0.152%
13	10.109	863	872	876	rBV3	53533	138720	2.05%	0.099%
14	10.188	876	880	884	rBV	136570	331896	4.89%	0.237%
15	10.306	889	892	898	rVB2	38625	90044	1.33%	0.064%
16	10.443	903	906	910	rVB2	48964	90001	1.33%	0.064%
17	10.581	915	920	922	rBV	80305	172348	2.54%	0.123%
18	10.630	922	925	929	rVB4	43768	100722	1.48%	0.072%
19	10.699	929	932	936	rVB	190861	347209	5.12%	0.248%
20	10.778	936	940	944	rBV	264401	572775	8.44%	0.408%
21	10.837	944	946	951	rVB3	59980	126915	1.87%	0.090%
22	10.926	951	955	958	rBV3	50246	107776	1.59%	0.077%
23	10.985	958	961	963	rBV	100565	158962	2.34%	0.113%
24	11.044	963	967	970	rBV3	198127	381740	5.63%	0.272%
25	11.211	976	984	987	rBV2	171966	407815	6.01%	0.291%
26	11.280	987	991	996	rVB	1589015	2446308	36.07%	1.744%
27	11.368	996	1000	1004	rVB3	68953	181688	2.68%	0.130%
28	11.447	1004	1008	1012	rBV4	129968	302476	4.46%	0.216%
29	11.526	1012	1016	1017	rBV2	189473	343059	5.06%	0.245%
30	11.595	1021	1023	1025	rBV	319633	478170	7.05%	0.341%
31	11.624	1025	1026	1034	rVB2	359196	860883	12.69%	0.614%
32	11.742	1034	1038	1042	rBV2	102927	244428	3.60%	0.174%
33	11.801	1042	1044	1048	rVB2	45640	89330	1.32%	0.064%
34	11.890	1048	1053	1056	rBV3	479720	1195427	17.62%	0.852%

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35	11.929	1056	1057	1059	rVB2	111816	106345	1.57%	0.076%
36	11.979	1059	1062	1064	rVB3	153879	303904	4.48%	0.217%
37	12.018	1064	1066	1068	rVB	79698	88378	1.30%	0.063%
38	12.097	1068	1074	1077	rBV4	523840	1370625	20.21%	0.977%
39	12.185	1077	1083	1087	rBV4	911262	2237807	32.99%	1.596%
40	12.274	1089	1092	1095	rVB2	335482	598660	8.83%	0.427%
41	12.343	1095	1099	1103	rBV3	484455	1192605	17.58%	0.850%
42	12.421	1103	1107	1111	rVB2	990464	1805387	26.62%	1.287%
43	12.480	1111	1113	1115	rBV2	311834	442344	6.52%	0.315%
44	12.530	1115	1118	1119	rBV3	175651	252901	3.73%	0.180%
45	12.579	1119	1123	1130	rVB3	1010739	2059505	30.36%	1.469%
46	12.687	1130	1134	1138	rBV3	477665	1192437	17.58%	0.850%
47	12.786	1138	1144	1145	rBV3	779081	2197245	32.39%	1.567%
48	12.825	1145	1148	1151	rVB2	1361785	2433021	35.87%	1.735%
49	12.884	1151	1154	1156	rBV3	131454	197313	2.91%	0.141%
50	12.923	1156	1158	1159	rVB	186332	197821	2.92%	0.141%
51	12.973	1159	1163	1165	rBV4	525720	1334897	19.68%	0.952%
52	13.012	1165	1167	1169	rVB2	121238	91279	1.35%	0.065%
53	13.051	1169	1171	1174	rVB	742733	911915	13.44%	0.650%
54	13.100	1174	1176	1180	rVB3	638710	1247748	18.40%	0.890%
55	13.179	1181	1184	1185	rBV2	350518	527140	7.77%	0.376%
56	13.219	1185	1188	1192	rBV4	864282	1671495	24.64%	1.192%
57	13.287	1192	1195	1201	rBV4	1773883	3238848	47.75%	2.309%
58	13.366	1201	1203	1207	rVB3	869024	1293979	19.08%	0.923%
59	13.425	1207	1209	1212	rVB3	219280	279827	4.13%	0.200%
60	13.494	1212	1216	1220	rVB4	568239	1088322	16.04%	0.776%
61	13.553	1220	1222	1227	rVB4	834585	1531759	22.58%	1.092%
62	13.622	1227	1229	1230	rBV2	314986	393624	5.80%	0.281%
63	13.661	1230	1233	1236	rBV4	1456253	2699375	39.80%	1.925%
64	13.730	1237	1240	1243	rBV2	2108679	2684118	39.57%	1.914%
65	13.789	1243	1246	1249	rVB3	909769	1498769	22.10%	1.069%
66	13.868	1250	1254	1256	rBV2	1242475	2516768	37.10%	1.795%
67	13.898	1256	1257	1260	rVB	1432339	1561511	23.02%	1.113%
68	13.986	1262	1266	1271	rVB3	1392725	2900374	42.76%	2.068%
69	14.075	1271	1275	1277	rVB4	1033665	1928644	28.43%	1.375%
70	14.134	1277	1281	1282	rBV3	1677749	2873395	42.36%	2.049%
71	14.163	1282	1284	1289	rVB3	2157782	4803605	70.82%	3.425%

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72	14.232	1289	1291	1292	rBV2	311625	371186	5.47%	0.265%
73	14.272	1292	1295	1298	rVB3	1982794	2911244	42.92%	2.076%
74	14.321	1298	1300	1302	rBV2	668737	1217110	17.94%	0.868%
75	14.419	1308	1310	1314	rBV3	789597	1885733	27.80%	1.345%
76	14.488	1314	1317	1318	rVV2	2446146	3635689	53.60%	2.592%
77	14.518	1318	1320	1324	rVV2	3120971	6783002	100.00%	4.837%
78	14.586	1324	1327	1330	rVV2	2733194	4756347	70.12%	3.392%
79	14.655	1330	1334	1336	rVV4	1559254	3313019	48.84%	2.362%
80	14.754	1340	1344	1345	rVV3	1346601	3033205	44.72%	2.163%
81	14.842	1348	1353	1355	rVV2	1813276	4086435	60.25%	2.914%
82	14.892	1355	1358	1359	rVV3	1393057	2499271	36.85%	1.782%
83	14.911	1359	1360	1362	rVV2	1284649	1530098	22.56%	1.091%
84	14.990	1365	1368	1370	rVV2	2928151	4894465	72.16%	3.490%
85	15.019	1370	1371	1373	rVV	1246829	1432420	21.12%	1.021%
86	15.069	1373	1376	1378	rVV3	2257393	3610266	53.23%	2.574%
87	15.138	1381	1383	1387	rVB3	631618	1473370	21.72%	1.051%
88	15.197	1387	1389	1391	rBV	1010681	1376003	20.29%	0.981%
89	15.265	1393	1396	1398	rVB3	965941	1618368	23.86%	1.154%
90	15.325	1398	1402	1403	rBV4	1242409	2174222	32.05%	1.550%
91	15.423	1409	1412	1415	rVV3	721779	1085808	16.01%	0.774%
92	15.472	1415	1417	1420	rVV	890139	1303488	19.22%	0.929%
93	15.580	1424	1428	1430	rVB	1354924	2312017	34.09%	1.649%
94	15.620	1430	1432	1436	rVB4	382324	620941	9.15%	0.443%
95	15.718	1440	1442	1445	rVB	588611	889562	13.11%	0.634%
96	15.777	1445	1448	1453	rVB3	386792	925422	13.64%	0.660%
97	15.895	1457	1460	1465	rVB3	383772	851321	12.55%	0.607%
98	16.102	1478	1481	1487	rVB5	125581	269970	3.98%	0.193%

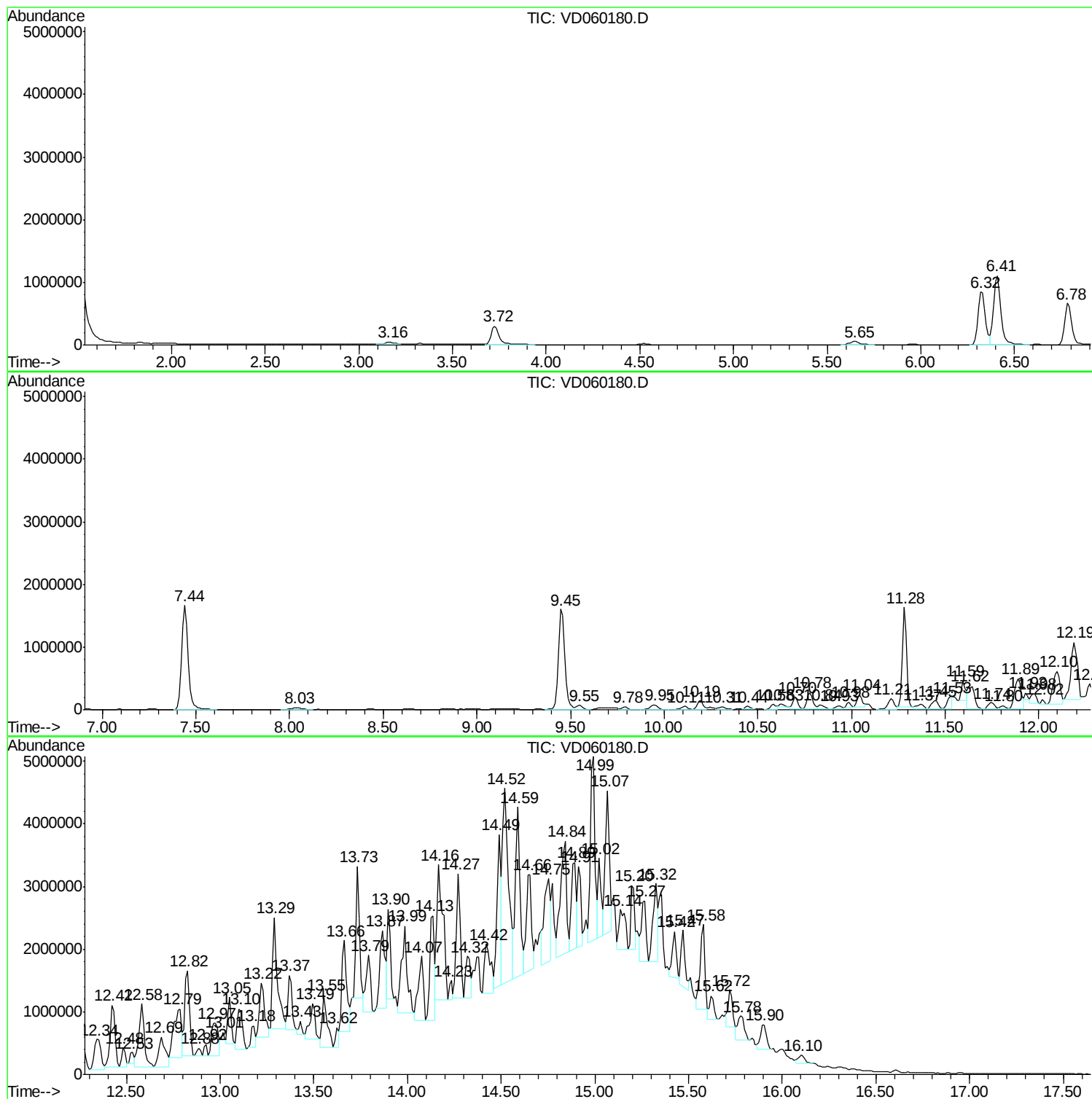
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Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D100118S.M
Quant Title : SW846 8260

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



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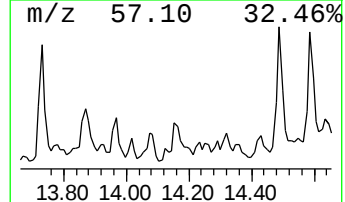
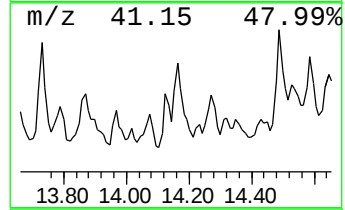
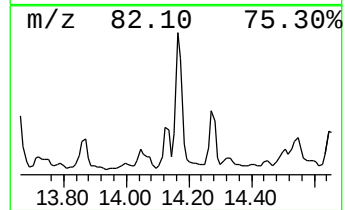
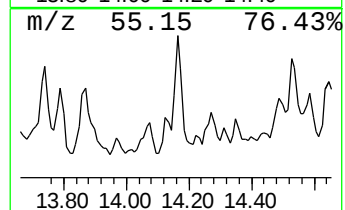
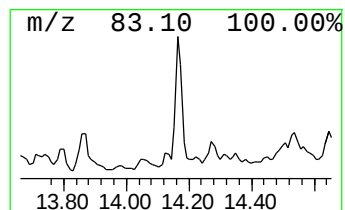
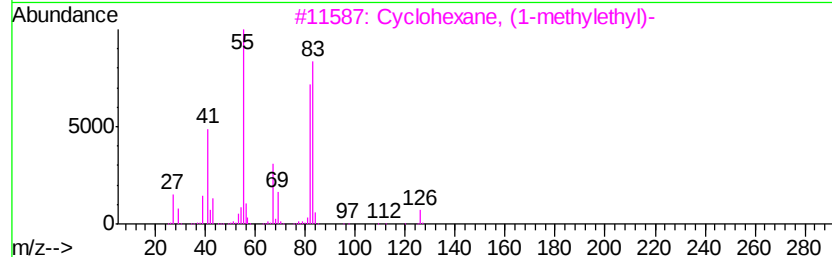
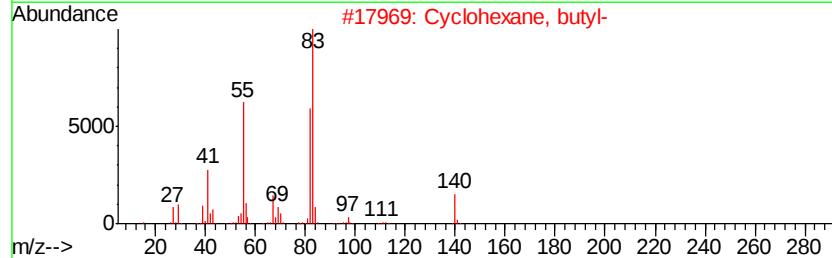
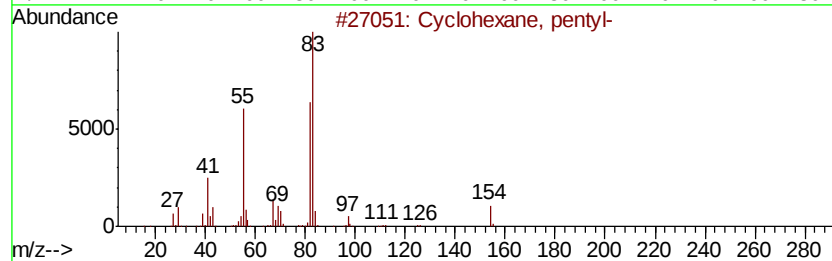
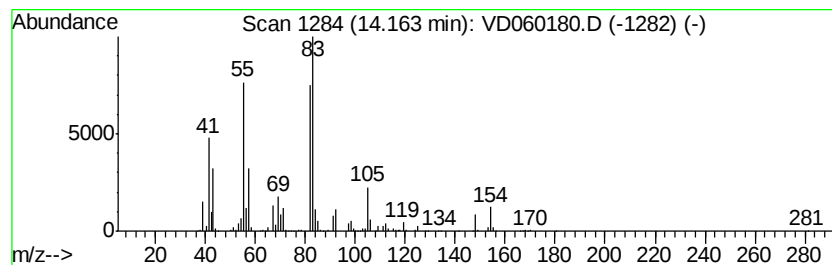
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D100118S.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclohexane, pentyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.16	185.61 ug/l	4803610	1,4-Dichlorobenzene-d4	13.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, pentyl-	154	C11H22	004292-92-6	81
2		Cyclohexane, butyl-	140	C10H20	001678-93-9	59
3		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	59
4		Cyclohexane, octyl-	196	C14H28	001795-15-9	59
5		Cyclohexane, propyl-	126	C9H18	001678-92-8	59



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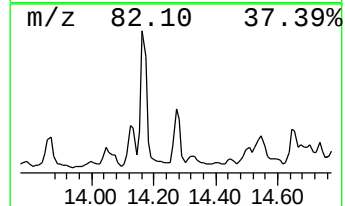
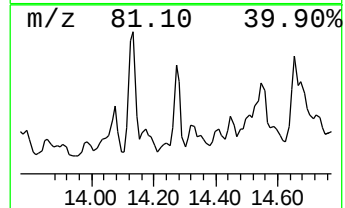
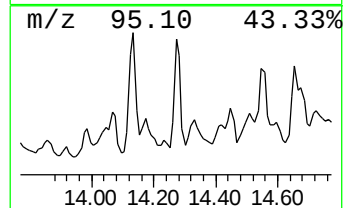
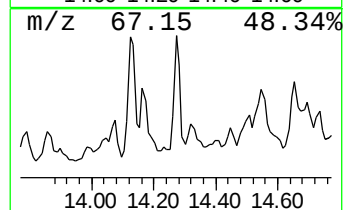
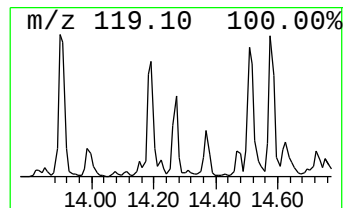
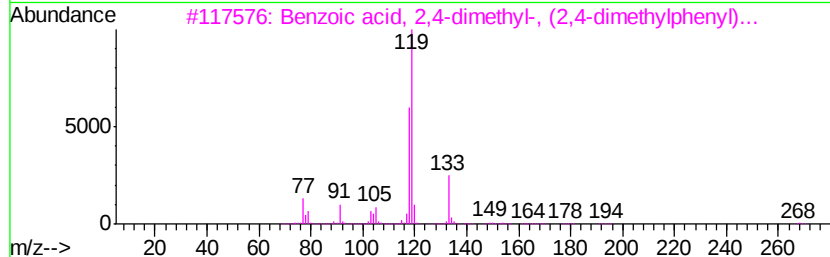
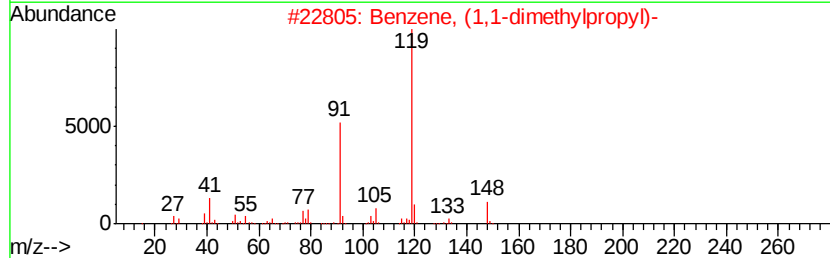
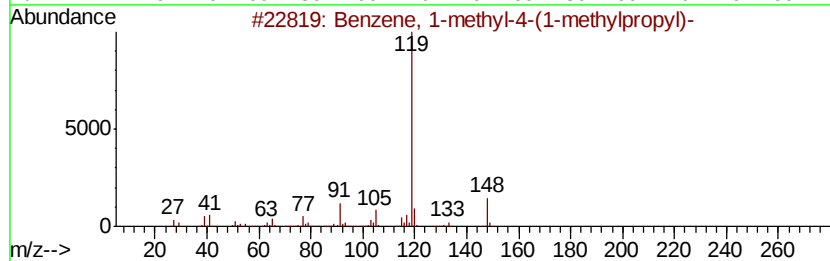
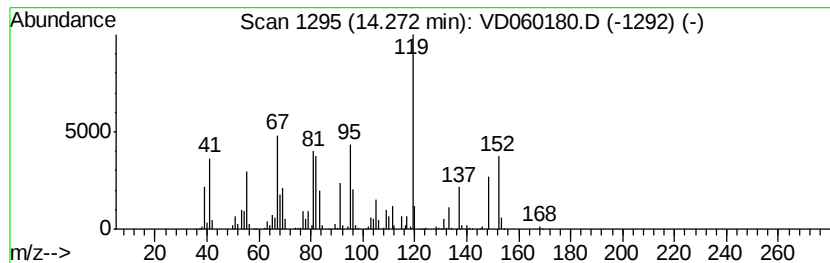
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown14.27 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	112.49 ug/l	2911240	1,4-Dichlorobenzene-d4	13.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	42
2		Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	27
3		Benzoic acid, 2,4-dimethyl-, (2,...	268	C18H20O2	055000-43-6	27
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	22
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	22



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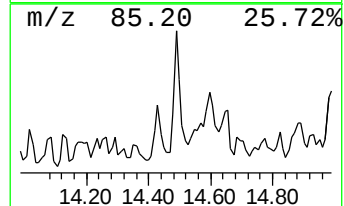
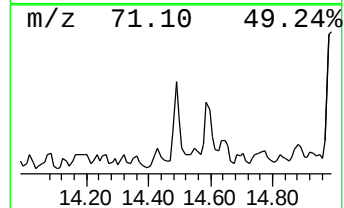
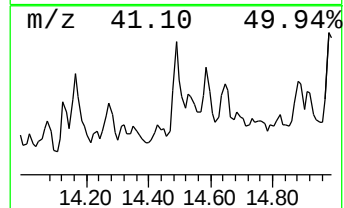
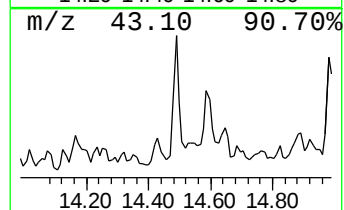
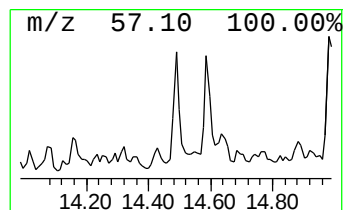
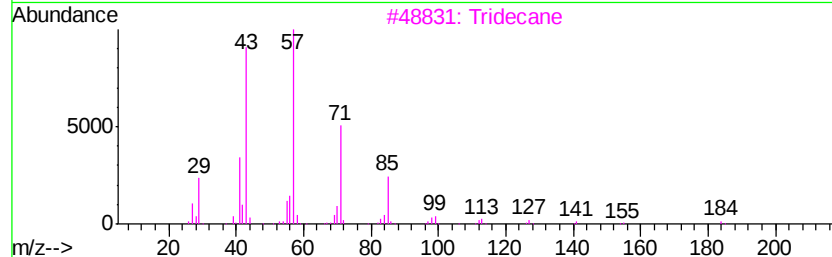
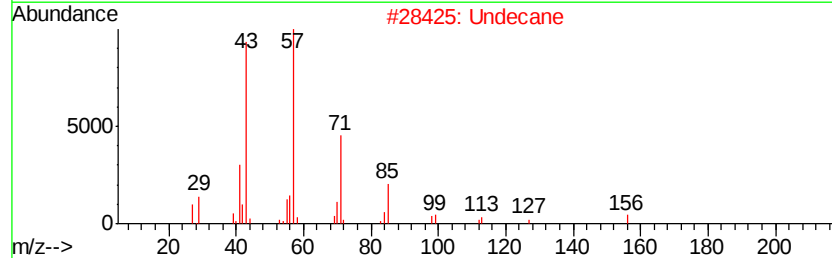
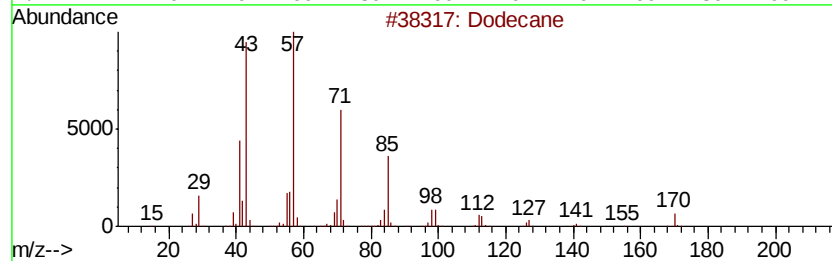
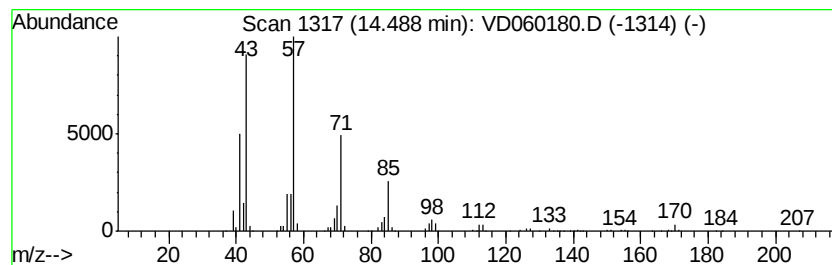
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Quant Title : SW846 8260

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

Peak Number 3 Dodecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.49	140.49 ug/l	3635690	1,4-Dichlorobenzene-d4	13.38

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane		170	C12H26	000112-40-3	95
2	Undecane		156	C11H24	001120-21-4	86
3	Tridecane		184	C13H28	000629-50-5	81
4	Undecane, 3,5-dimethyl-		184	C13H28	017312-81-1	80
5	Tetradecane		198	C14H30	000629-59-4	78



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD102418\
Data File : VD060180.D
Acq On : 24 Oct 2018 14:25
Operator : JC/SY
Sample : J5622-03
Misc : 4.91a/5ml/MSVOA D/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
KG-V23608

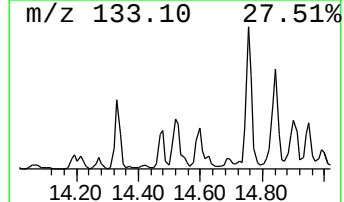
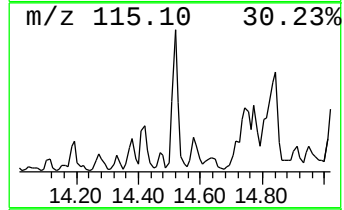
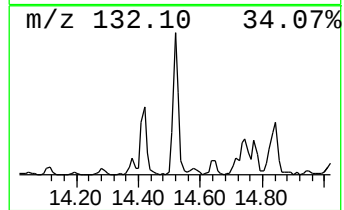
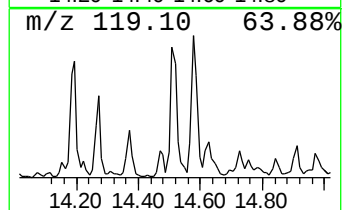
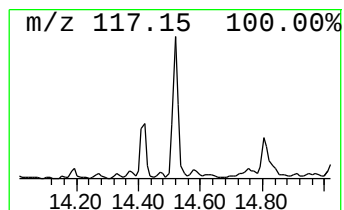
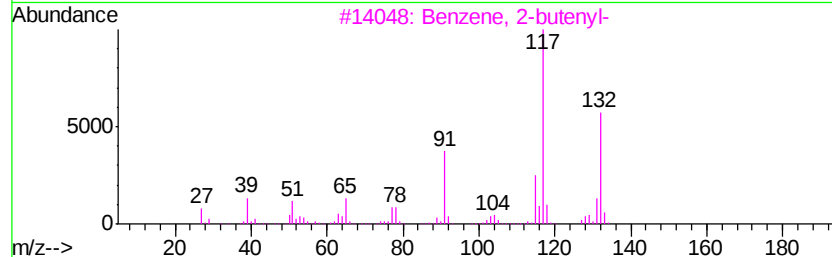
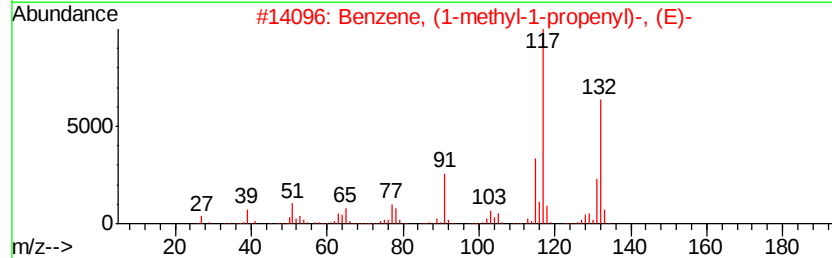
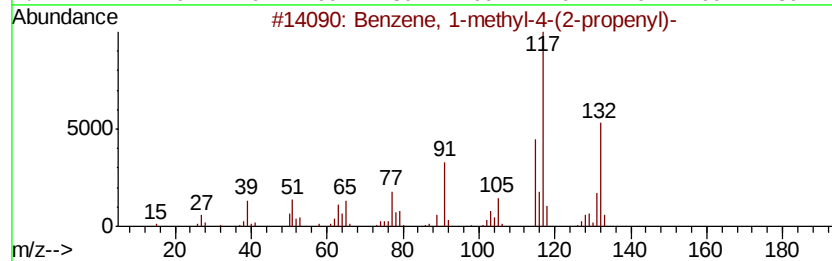
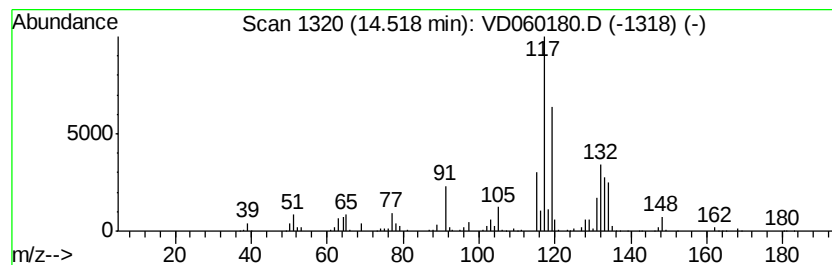
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.52	262.10 ug/l	6783000	1,4-Dichlorobenzene-d4	13.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	86
2		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	64
3		Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	60
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD102418\
Data File : VD060180.D
Acq On : 24 Oct 2018 14:25
Operator : JC/SY
Sample : J5622-03
Misc : 4.91a/5ml/MSVOA D/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
KG-V23608

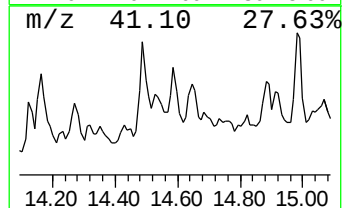
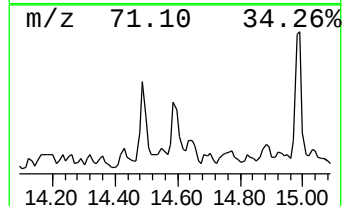
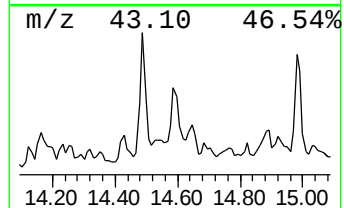
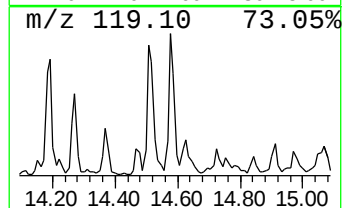
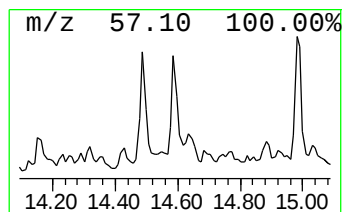
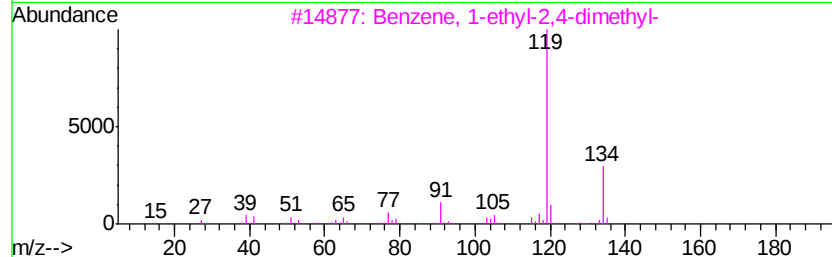
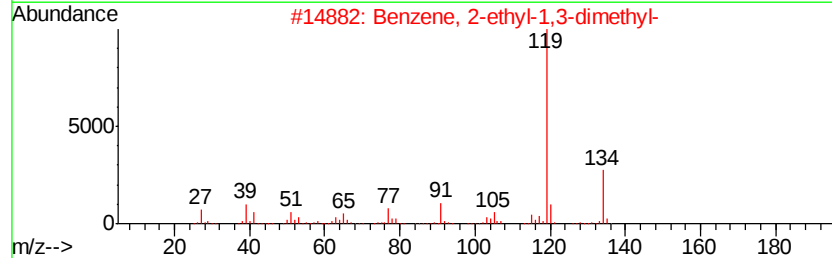
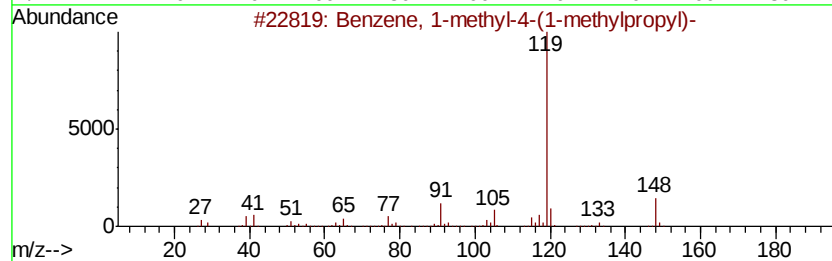
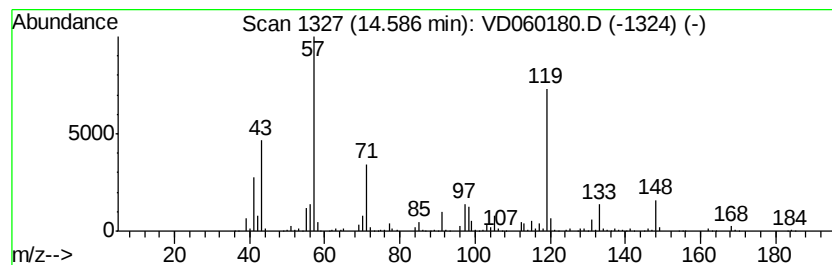
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-methyl-4-(1-meth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	183.79 ug/l	4756350	1,4-Dichlorobenzene-d4	13.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	83
2		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	38
3		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	38
4		Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	38
5		Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD102418\
Data File : VD060180.D
Acq On : 24 Oct 2018 14:25
Operator : JC/SY
Sample : J5622-03
Misc : 4.91a/5ml/MSVOA D/SOIL
ALS Vial : 12 Sample Multiplier: 1

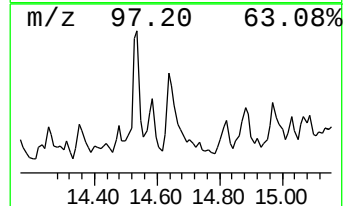
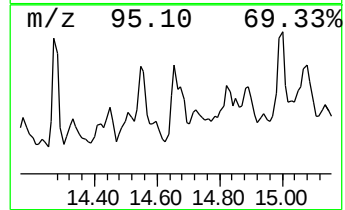
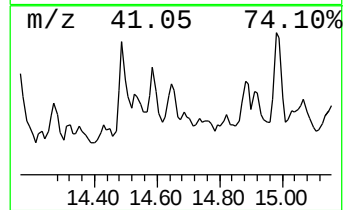
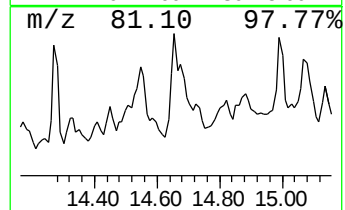
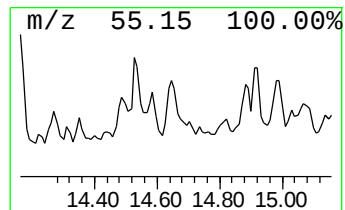
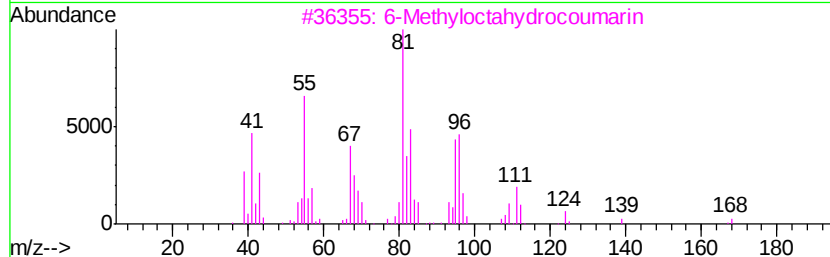
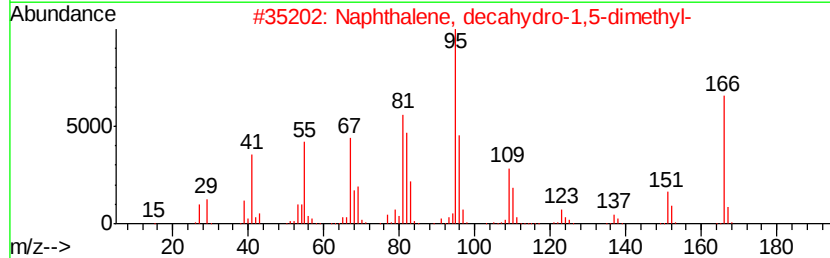
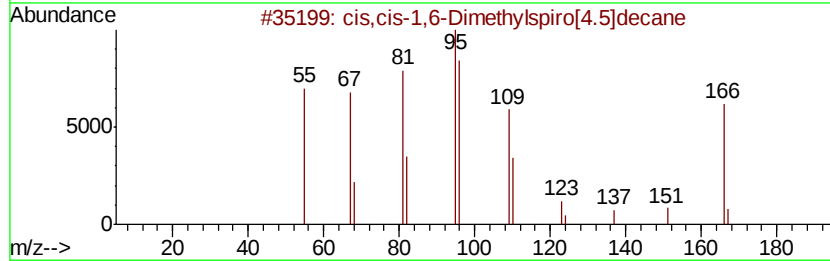
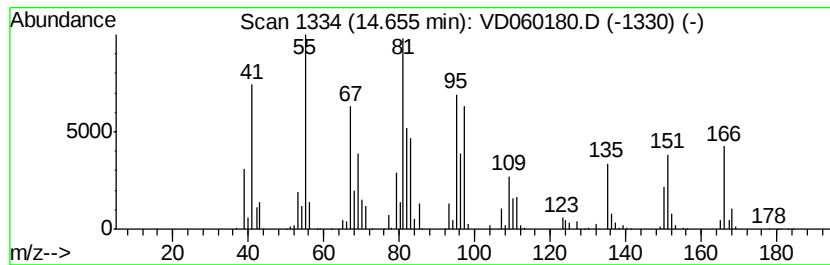
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ClientSampleId :
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Quant Title : SW846 8260

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

Peak Number 6 unknown14.66 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.66	128.02 ug/l	3313020	1,4-Dichlorobenzene-d4	13.38
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		cis,cis-1,6-Dimethylspiro[4.5]de...	166 C12H22	1000111-72-4 49
2		Naphthalene, decahydro-1,5-dimet...	166 C12H22	066552-62-3 46
3		6-Methyloctahydrocoumarin	168 C10H16O2	080648-29-9 45
4		Naphthalene, decahydro-2,6-dimet...	166 C12H22	001618-22-0 43
5		trans,trans-1,6-Dimethylspiro[4....	166 C12H22	1000111-72-1 42



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD102418\
Data File : VD060180.D
Acq On : 24 Oct 2018 14:25
Operator : JC/SY
Sample : J5622-03
Misc : 4.91a/5ml/MSVOA D/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
KG-V23608

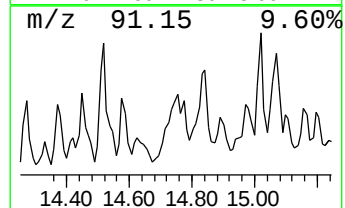
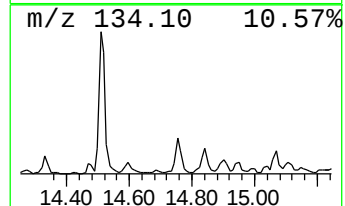
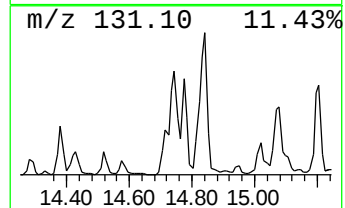
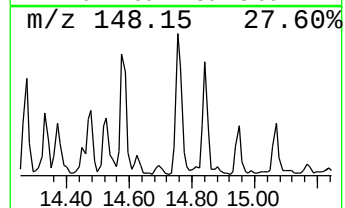
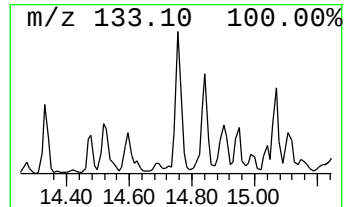
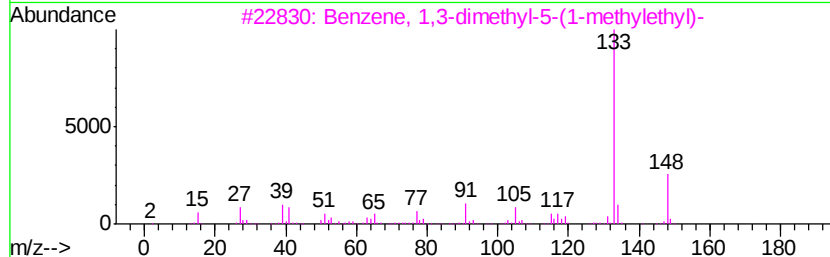
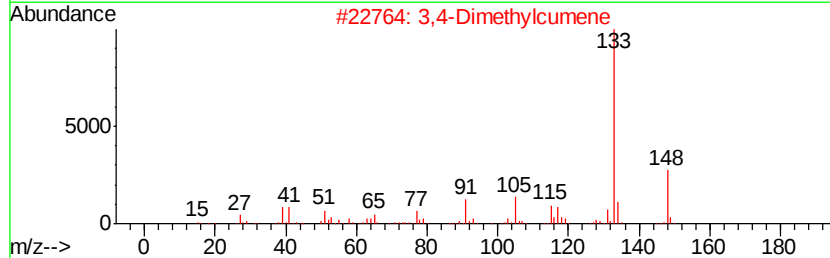
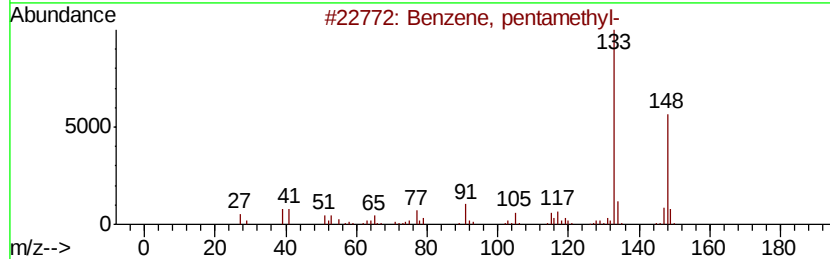
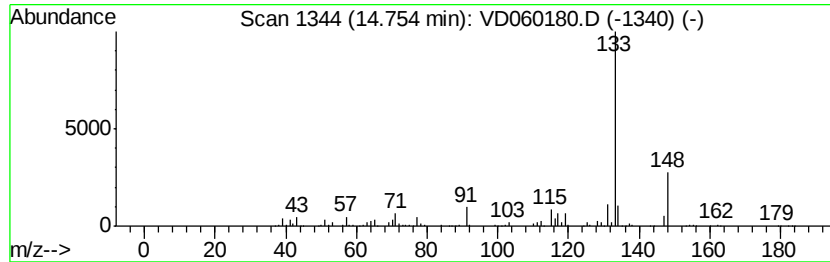
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, pentamethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.75	117.21 ug/l	3033210	1,4-Dichlorobenzene-d4	13.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, pentamethyl-	148	C11H16	000700-12-9	90
2		3,4-Dimethylcumene	148	C11H16	1000370-34-1	86
3		Benzene, 1,3-dimethyl-5-(1-methv...	148	C11H16	004706-90-5	86
4		Benzene, 1-ethyl-2,4,5-trimethyl-	148	C11H16	017851-27-3	83
5		Benzene, 2,4-dimethyl-1-(1-methy...	148	C11H16	004706-89-2	80



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD102418\
Data File : VD060180.D
Acq On : 24 Oct 2018 14:25
Operator : JC/SY
Sample : J5622-03
Misc : 4.91a/5ml/MSVOA D/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
KG-V23608

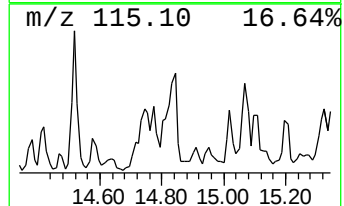
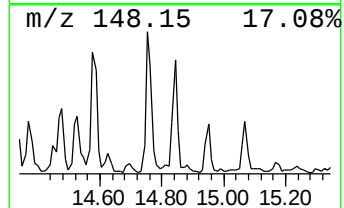
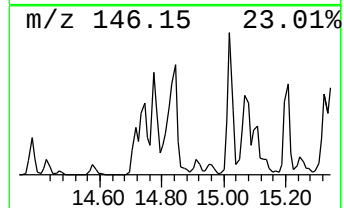
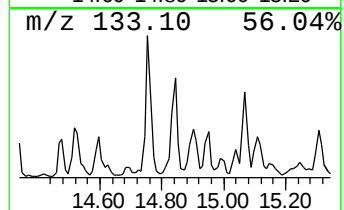
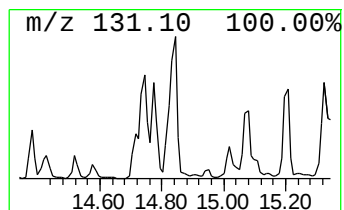
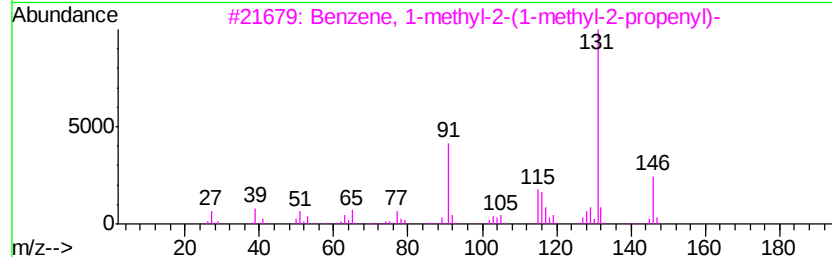
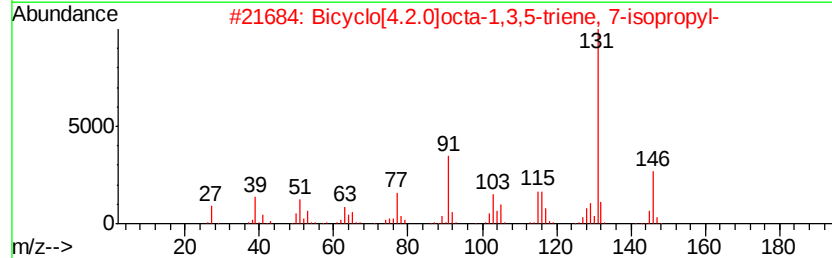
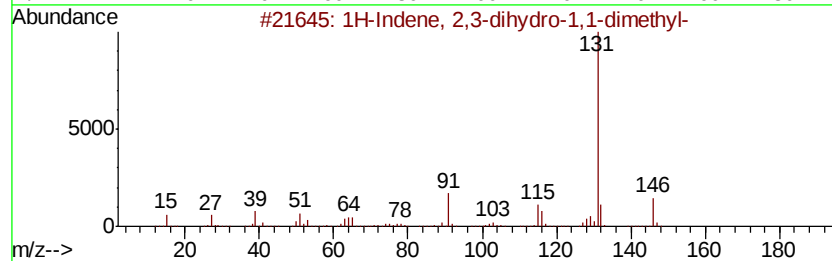
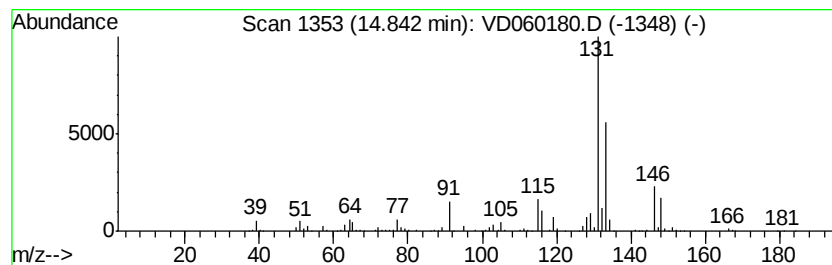
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Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-1,1-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.84	157.90 ug/l	4086440	1,4-Dichlorobenzene-d4	13.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	64
2		Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	62
3		Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	58
4		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	58
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD102418\
Data File : VD060180.D
Acq On : 24 Oct 2018 14:25
Operator : JC/SY
Sample : J5622-03
Misc : 4.91a/5ml/MSVOA D/SOIL
ALS Vial : 12 Sample Multiplier: 1

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ClientSampleId :
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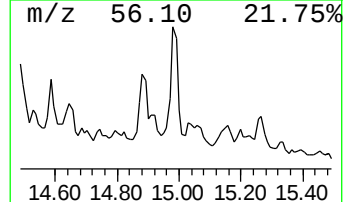
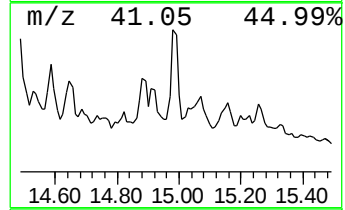
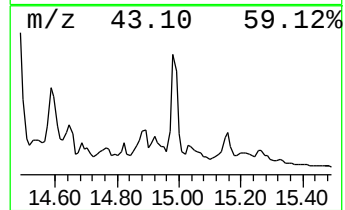
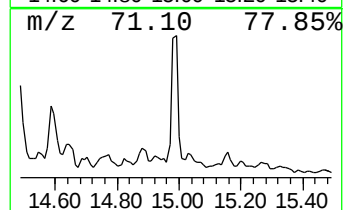
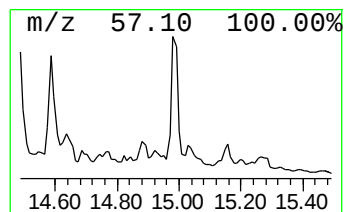
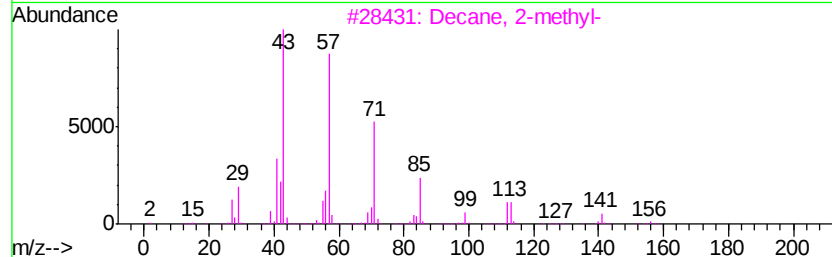
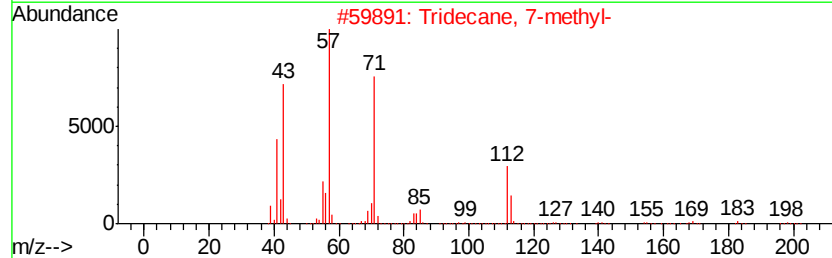
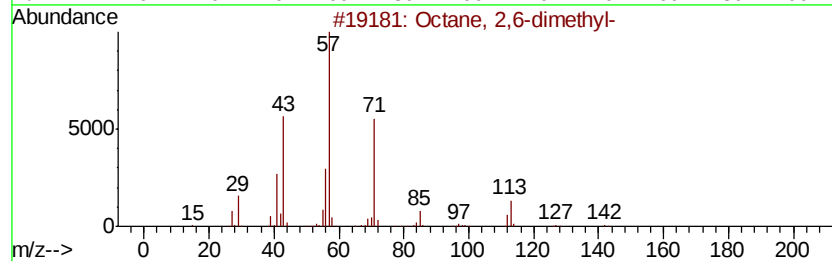
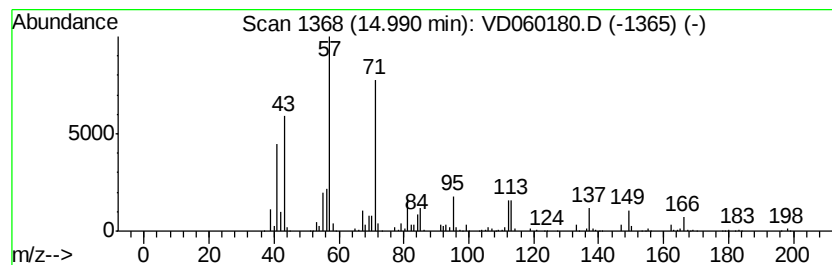
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.99	189.13 ug/l	4894470	1,4-Dichlorobenzene-d4	13.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64
2		Tridecane, 7-methyl-	198	C14H30	026730-14-3	64
3		Decane, 2-methyl-	156	C11H24	006975-98-0	59
4		Nonane, 3-methyl-	142	C10H22	005911-04-6	59
5		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD102418\
Data File : VD060180.D
Acq On : 24 Oct 2018 14:25
Operator : JC/SY
Sample : J5622-03
Misc : 4.91a/5ml/MSVOA D/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampled :
KG-V23608

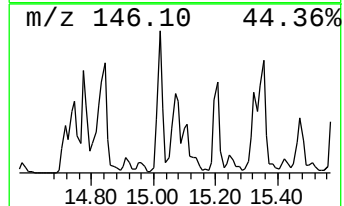
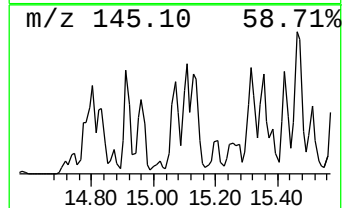
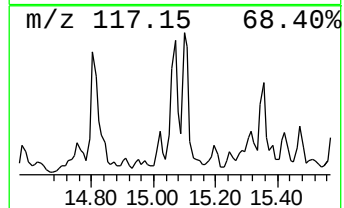
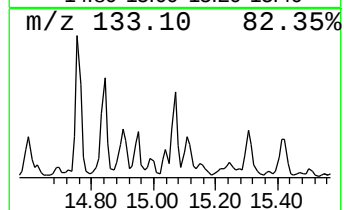
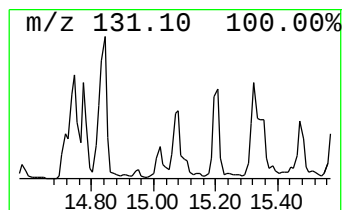
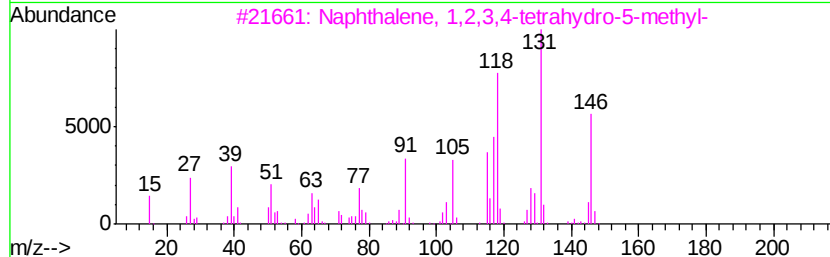
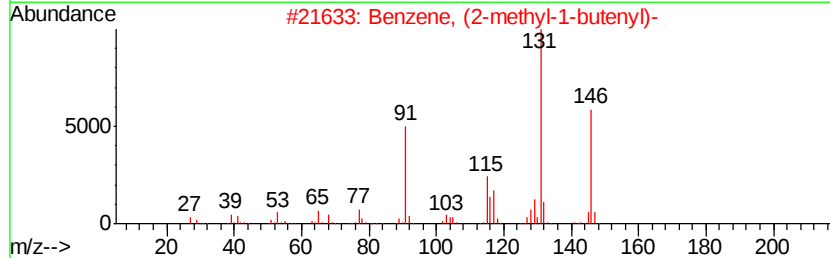
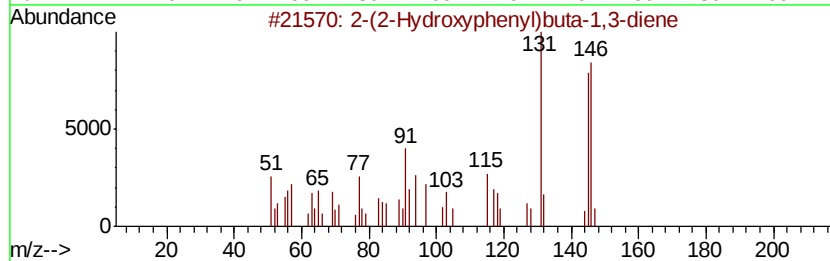
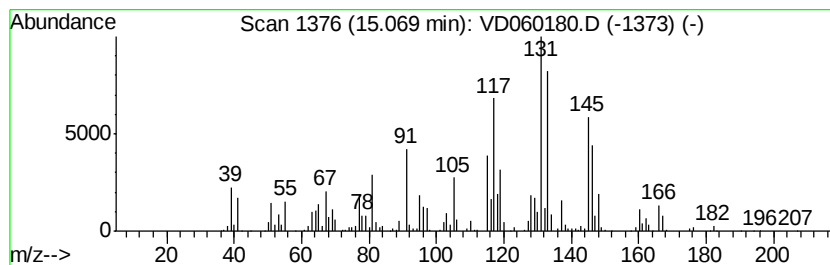
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D100118S.M
Quant Title : SW846 8260

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

Peak Number 10 2-(2-Hydroxyphenyl)buta-1,3... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	139.50 ug/l	3610270	1,4-Dichlorobenzene-d4	13.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-(2-Hydroxyphenyl)buta-1,3-diene	146	C10H10O	038865-47-3	58
2		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	46
3		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	46
4		Benzene, (3-chloro-1-methyl-1-pr...	166	C10H11Cl	1000327-43-7	44
5		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	42



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD102418\
Data File : VD060180.D
Acq On : 24 Oct 2018 14:25
Operator : JC/SY
Sample : J5622-03
Misc : 4.91g/5ml/MSVOA_D/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
KG-V23608

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D100118S.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, pentyl-	14.16	185.6	ug/l	4803610	4	13.38	1293980	50.0
unknown14.27	14.27	112.5	ug/l	2911240	4	13.38	1293980	50.0
Dodecane	14.49	140.5	ug/l	3635690	4	13.38	1293980	50.0
Benzene, 1-methyl...	14.52	262.1	ug/l	6783000	4	13.38	1293980	50.0
Benzene, 1-methyl...	14.59	183.8	ug/l	4756350	4	13.38	1293980	50.0
unknown14.66	14.66	128.0	ug/l	3313020	4	13.38	1293980	50.0
Benzene, pentamet...	14.75	117.2	ug/l	3033210	4	13.38	1293980	50.0
1H-Indene, 2,3-di...	14.84	157.9	ug/l	4086440	4	13.38	1293980	50.0
Octane, 2,6-dimet...	14.99	189.1	ug/l	4894470	4	13.38	1293980	50.0
2-(2-Hydroxypheny...	15.07	139.5	ug/l	3610270	4	13.38	1293980	50.0