

Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
 Data File : VH057098.D
 Acq On : 21 Sep 2015 18:40
 Operator : SY/FY
 Sample : G3747-02
 Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_H
 ClientSampleId :
 TRENCH-1-2

Integration Parameters: rteint.p

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.001
 Stop Thrs : 0

Filtering: 9
 Min Area: 250 Area counts
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 1

Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.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	4.070	340	346	359	rBV	476769	1526046	6.25%	0.347%
2	4.827	410	418	432	rBV2	964100	3695036	15.12%	0.840%
3	5.562	482	488	498	rBV	1101764	3089822	12.65%	0.702%
4	7.502	669	673	688	rVV	1552087	3856684	15.79%	0.876%
5	8.407	755	759	766	rVB2	185012	484472	1.98%	0.110%
6	8.563	770	774	778	rBV	116641	303246	1.24%	0.069%
7	8.911	802	807	812	rBV2	138103	448250	1.83%	0.102%
8	9.069	818	822	828	rVB	269144	682999	2.80%	0.155%
9	9.237	833	838	843	rBV	256034	585843	2.40%	0.133%
10	9.709	879	883	892	rVV	1548178	3902554	15.97%	0.887%
11	9.836	892	895	900	rVB2	109687	263663	1.08%	0.060%
12	10.172	921	927	935	rBV4	212742	762151	3.12%	0.173%
13	10.299	935	939	944	rBV2	150529	348008	1.42%	0.079%
14	10.456	948	954	958	rBV3	259480	819699	3.36%	0.186%
15	10.614	965	969	977	rBV2	1030479	2298062	9.41%	0.522%
16	10.730	977	980	982	rVV2	185982	423502	1.73%	0.096%
17	10.772	982	984	992	rVV2	312030	829223	3.39%	0.188%
18	10.889	992	995	999	rVB4	129993	311744	1.28%	0.071%
19	10.993	1000	1005	1008	rBV2	498514	1275246	5.22%	0.290%
20	11.067	1008	1012	1014	rVV2	989424	2157063	8.83%	0.490%
21	11.109	1014	1016	1019	rVV	958483	1580241	6.47%	0.359%
22	11.214	1019	1026	1030	rVV2	1479858	3427280	14.03%	0.779%
23	11.349	1034	1039	1045	rVV3	1148726	3370378	13.79%	0.766%
24	11.434	1045	1047	1050	rVV2	423305	944695	3.87%	0.215%
25	11.496	1050	1053	1056	rVV	1314739	2640248	10.81%	0.600%
26	11.549	1056	1058	1062	rVV3	712586	1960241	8.02%	0.445%
27	11.613	1062	1064	1067	rVV2	624128	1236213	5.06%	0.281%
28	11.675	1067	1070	1072	rVV2	522329	1224119	5.01%	0.278%
29	11.717	1072	1074	1076	rVV2	622653	1086761	4.45%	0.247%
30	11.759	1076	1078	1082	rVV3	498366	1082423	4.43%	0.246%
31	11.832	1082	1085	1087	rVV2	613327	1291723	5.29%	0.294%
32	11.896	1087	1091	1093	rVV2	1042601	2772906	11.35%	0.630%
33	11.938	1093	1095	1099	rVV	3173842	5416681	22.17%	1.231%
34	12.011	1099	1102	1107	rVV2	986401	2917244	11.94%	0.663%

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Integration Parameters: rteint.p

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.001

Stop Thrs : 0

Filtering: 9

Min Area: 250 Area counts

Max Peaks: 100

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35	12.096	1107	1110	1112	rVV2	1573684	3775832	15.45%	0.858%
36	12.139	1112	1114	1118	rVV3	4293482	8941587	36.60%	2.032%
37	12.224	1119	1122	1125	rVV4	1300365	3523673	14.42%	0.801%
38	12.266	1125	1126	1128	rVV2	999771	1695330	6.94%	0.385%
39	12.319	1128	1131	1133	rVV	2431346	5151953	21.09%	1.171%
40	12.362	1133	1135	1137	rVV2	2521519	4619368	18.91%	1.050%
41	12.394	1137	1138	1143	rVV2	3121250	6600256	27.01%	1.500%
42	12.479	1143	1146	1149	rVV2	3709643	7311680	29.93%	1.662%
43	12.543	1149	1152	1155	rVV	2763981	6224098	25.48%	1.414%
44	12.607	1155	1158	1160	rVV4	1149861	3313003	13.56%	0.753%
45	12.681	1162	1165	1166	rVV2	1656495	3879117	15.88%	0.882%
46	12.724	1166	1169	1172	rVV2	3068686	9145723	37.43%	2.078%
47	12.777	1172	1174	1178	rVV3	3387413	8220122	33.64%	1.868%
48	12.869	1180	1183	1188	rVV4	2886207	10581281	43.31%	2.405%
49	12.954	1188	1191	1194	rVV	2650720	6404680	26.21%	1.455%
50	13.016	1194	1197	1199	rVV2	2915112	7997962	32.74%	1.818%
51	13.080	1202	1203	1205	rVV	3592162	5908142	24.18%	1.343%
52	13.134	1205	1208	1212	rVV3	4486605	13234017	54.17%	3.007%
53	13.187	1212	1213	1214	rVV	2115625	2351913	9.63%	0.534%
54	13.219	1214	1216	1219	rVV2	5426044	12127612	49.64%	2.756%
55	13.272	1219	1221	1222	rVV2	3601518	6223512	25.47%	1.414%
56	13.304	1222	1224	1227	rVV2	5526101	12644673	51.75%	2.874%
57	13.358	1227	1229	1233	rVV3	2312755	7654419	31.33%	1.739%
58	13.420	1233	1235	1237	rVV2	4770062	8535352	34.93%	1.940%
59	13.473	1237	1240	1247	rVV2	4155380	16085592	65.84%	3.655%
60	13.567	1247	1249	1251	rVV	2684631	5285770	21.63%	1.201%
61	13.642	1252	1256	1259	rVV3	3963030	14307862	58.56%	3.251%
62	13.738	1263	1265	1267	rVV2	4166732	8566448	35.06%	1.947%
63	13.781	1267	1269	1273	rVV4	10778638	24432123	100.00%	5.552%
64	13.845	1273	1275	1278	rVV	4797296	10345710	42.34%	2.351%
65	13.898	1278	1280	1282	rVV2	2645551	5194478	21.26%	1.180%
66	13.940	1282	1284	1285	rVV2	1826551	3296534	13.49%	0.749%
67	14.003	1285	1290	1293	rVV2	3552101	13428304	54.96%	3.052%
68	14.086	1293	1298	1302	rVV6	6241182	22759471	93.15%	5.172%
69	14.148	1302	1304	1309	rVV3	4706564	13386378	54.79%	3.042%
70	14.231	1309	1312	1315	rVV3	4109354	10272651	42.05%	2.334%
71	14.283	1315	1317	1320	rVV5	2944085	6934181	28.38%	1.576%

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72	14.358	1320	1324	1325	rVV3	2518727	7525477	30.80%	1.710%
73	14.379	1325	1326	1328	rVV2	3435988	5628733	23.04%	1.279%
74	14.422	1328	1330	1334	rVV3	2738358	7720170	31.60%	1.754%
75	14.485	1334	1336	1337	rVV	2668765	4296078	17.58%	0.976%
76	14.517	1337	1339	1346	rVV4	2555072	9759872	39.95%	2.218%
77	14.656	1350	1352	1358	rVV3	3308280	8277169	33.88%	1.881%
78	14.750	1358	1361	1362	rVV	1335364	2849350	11.66%	0.648%
79	14.782	1362	1364	1370	rVV5	1257659	4877075	19.96%	1.108%
80	14.909	1374	1376	1381	rVV3	1228155	3238239	13.25%	0.736%
81	14.982	1381	1383	1385	rVV2	435191	910116	3.73%	0.207%
82	15.023	1385	1387	1390	rVV2	806837	1597852	6.54%	0.363%
83	15.076	1390	1392	1398	rVV5	639941	1683228	6.89%	0.383%
84	15.191	1398	1403	1406	rVV2	668266	1540045	6.30%	0.350%
85	15.245	1406	1408	1411	rVV3	149084	383167	1.57%	0.087%
86	15.309	1411	1414	1417	rVB	182818	379043	1.55%	0.086%

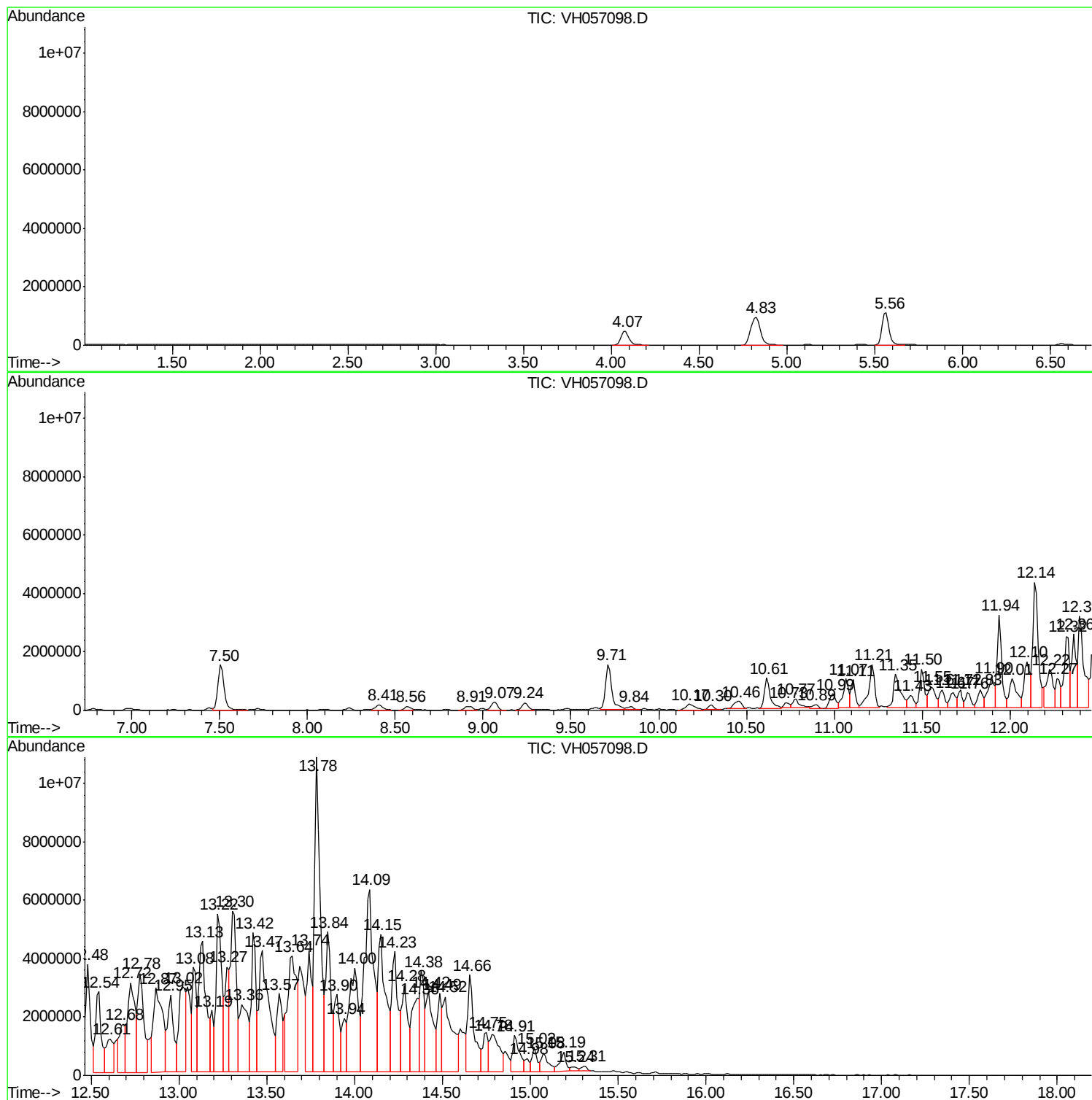
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Instrument :
MSVOA_H
ClientSampleId :
TRENCH-1-2

Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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TRENCH-1-2

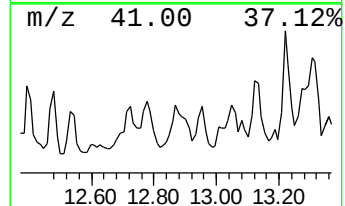
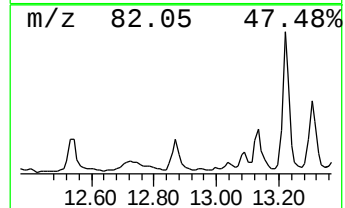
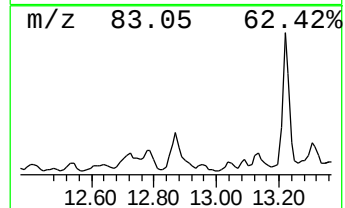
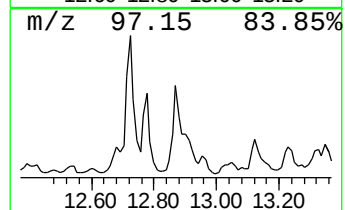
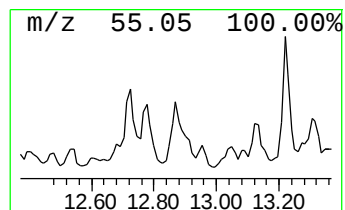
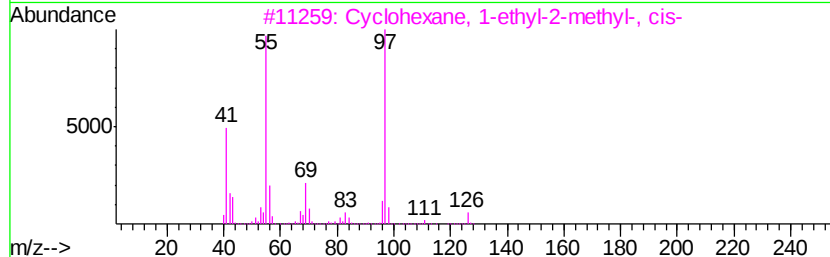
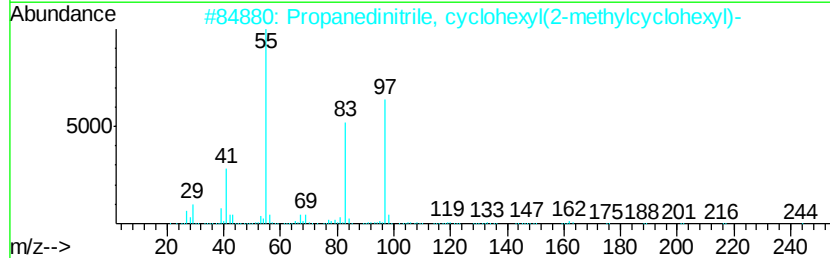
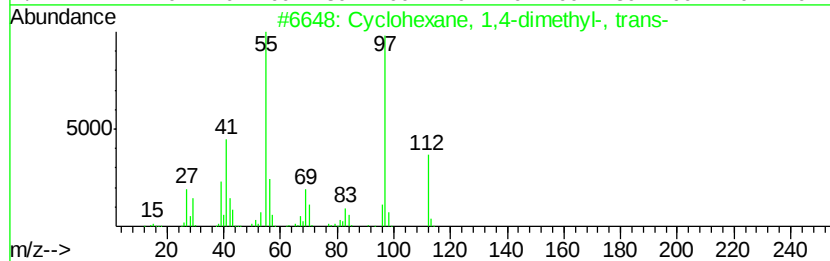
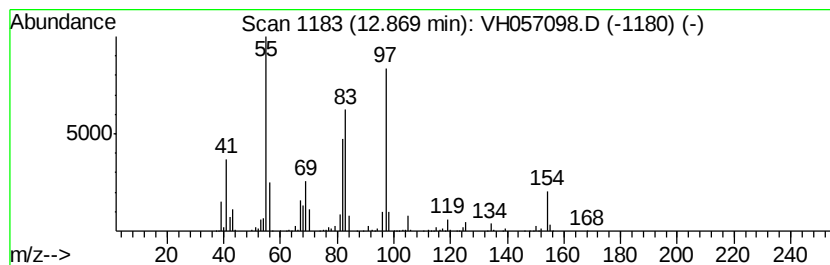
Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown12.87 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.87	72.36 ug/l	10581300	1,4-Dichlorobenzene-d4	12.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	47
2		Propanedinitrile, cyclohexyl(2-m...	244	C16H24N2	074764-55-9	47
3		Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-77-7	46
4		Cyclohexane, pentyl-	154	C11H22	004292-92-6	43
5		Cyclohexane, 1,4-dimethyl-	112	C8H16	000589-90-2	43



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Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 13 Sample Multiplier: 1

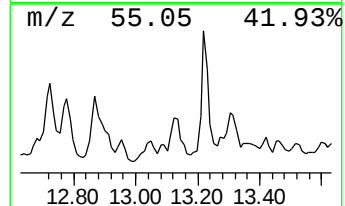
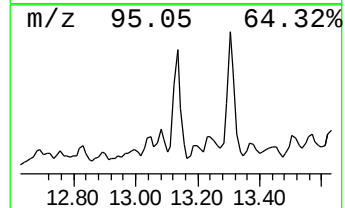
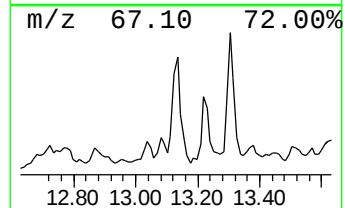
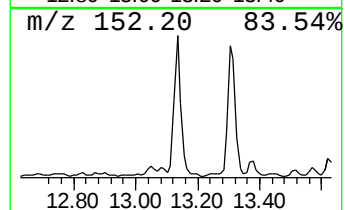
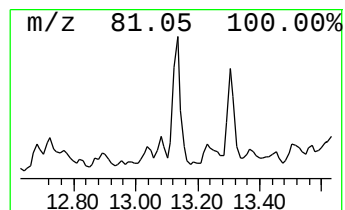
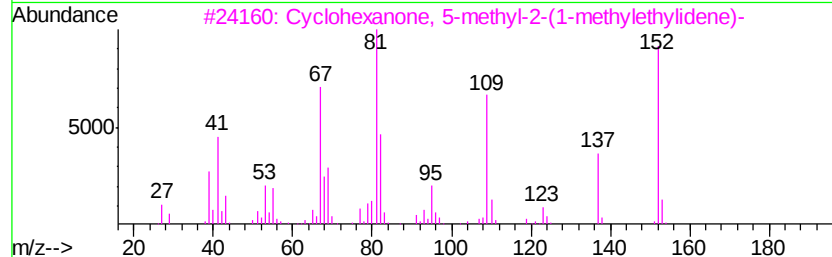
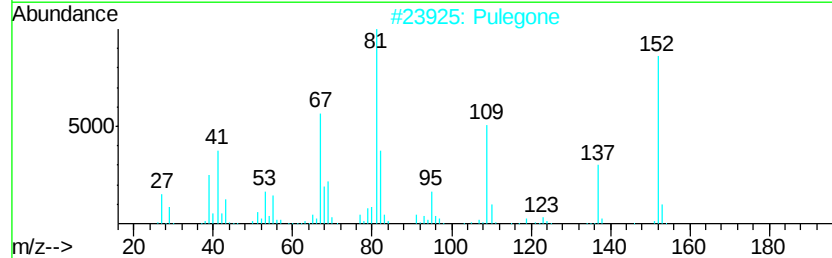
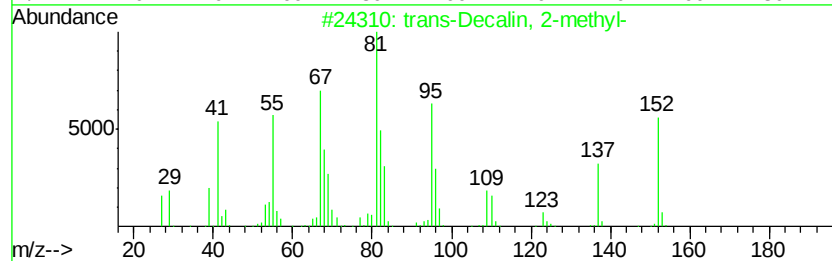
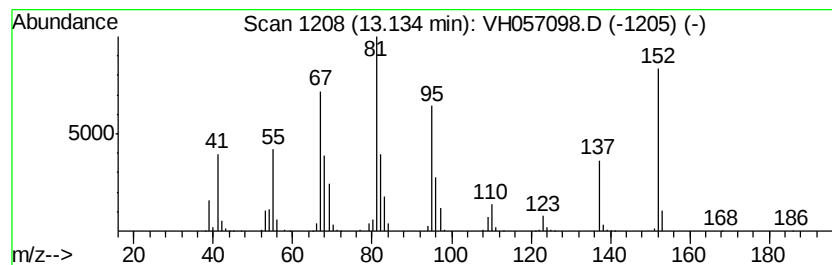
Instrument :
MSVOA_H
ClientSampleId :
TRENCH-1-2

Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 trans-Decalin, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.13	90.50 ug/l	13234000	1,4-Dichlorobenzene-d4	12.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	trans-Decalin, 2-methyl-	152 C11H20	1000152-47-3	94
2	Pulegone	152 C10H16O	000089-82-7	74
3	Cyclohexanone, 5-methyl-2-(1-methyl-2-propenyl)-	152 C10H16O	015932-80-6	74
4	Naphthalene, decahydro-2-methyl-	152 C11H20	002958-76-1	64
5	Bicyclo[4.1.0]heptan-3-one, 4,7,8-trimethyl-	152 C10H16O	004176-04-9	58



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ClientSampled :
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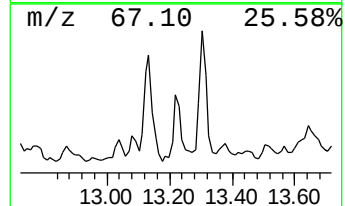
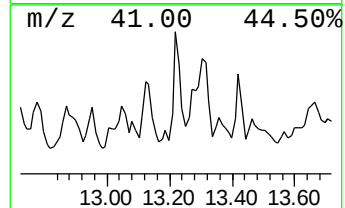
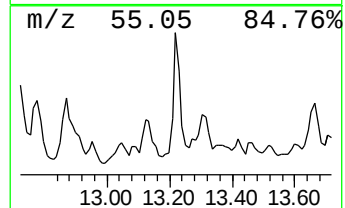
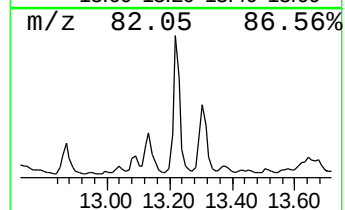
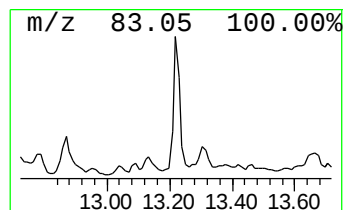
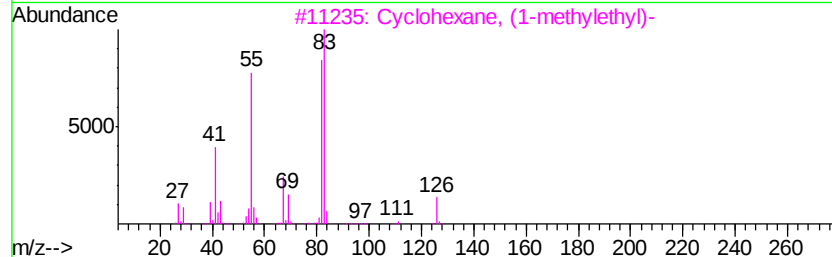
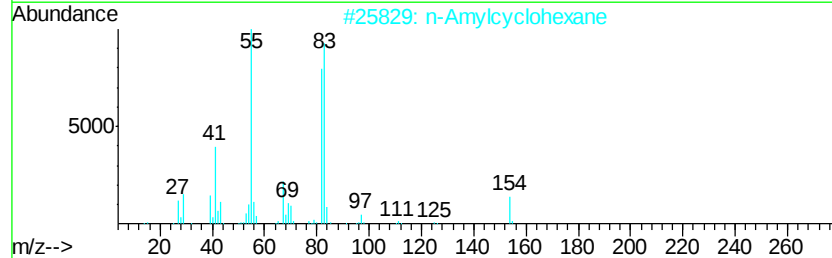
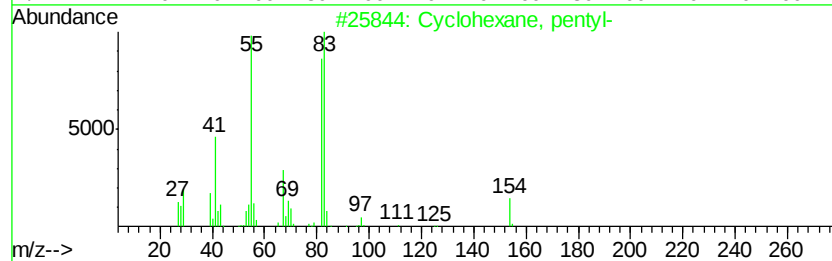
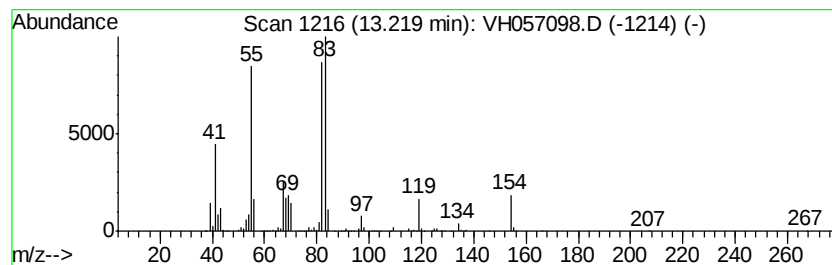
Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Cyclohexane, pentyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	82.93 ug/l	12127600	1,4-Dichlorobenzene-d4	12.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, pentyl-	154	C11H22	004292-92-6	90
2		n-Amylcyclohexane	154	C11H22	029949-27-7	83
3		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	74
4		Cyclohexane, undecyl-	238	C17H34	054105-66-7	72
5		Cyclohexane, octyl-	196	C14H28	001795-15-9	72



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Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 13 Sample Multiplier: 1

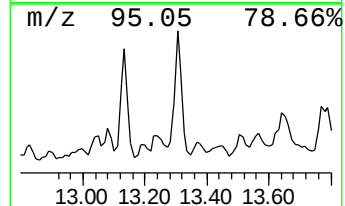
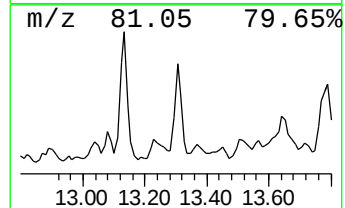
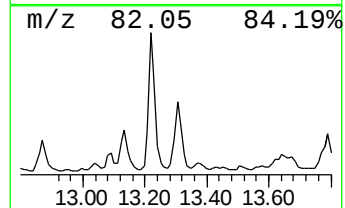
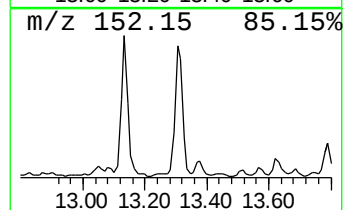
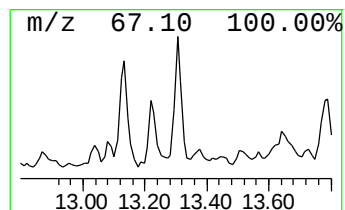
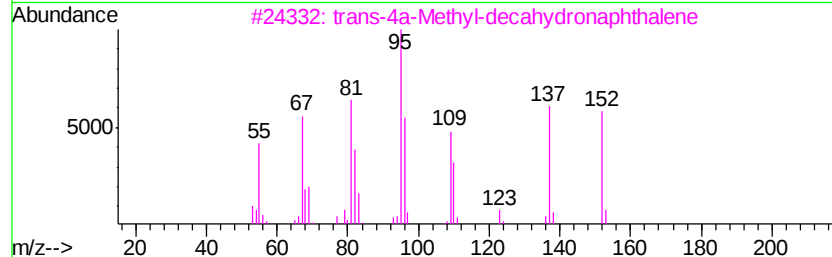
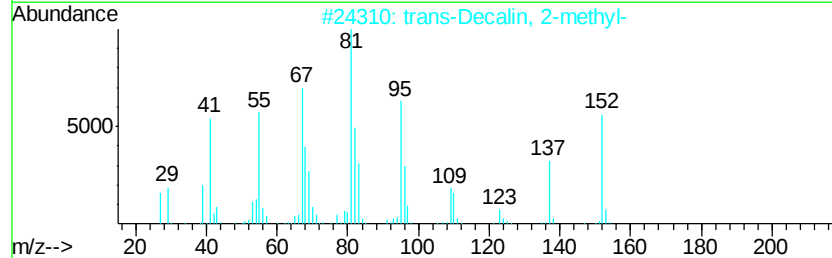
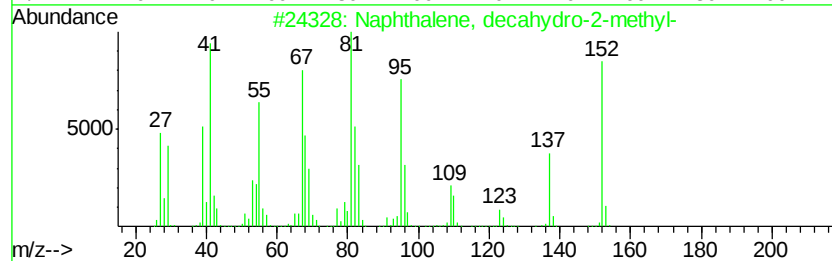
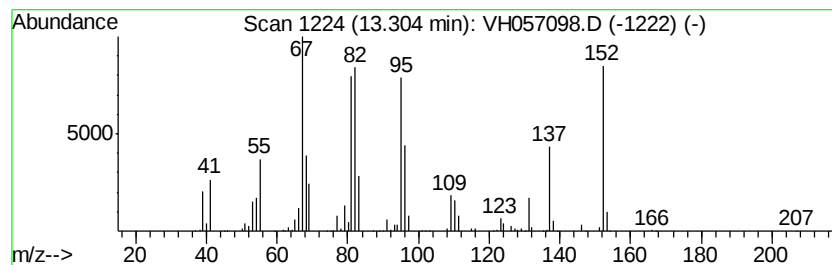
Instrument :
MSVOA_H
ClientSampleId :
TRENCH-1-2

Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Naphthalene, decahydro-2-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.30	86.47 ug/l	12644700	1,4-Dichlorobenzene-d4	12.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, decahydro-2-methyl-	152 C11H20	002958-76-1 93
2		trans-Decalin, 2-methyl-	152 C11H20	1000152-47-3 70
3		trans-4a-Methyl-decahydronaphtha...	152 C11H20	002547-27-5 64
4		Bicyclo[3.3.2]decan-9-one	152 C10H16O	028054-91-3 58
5		Cyclododecene, (Z)-	166 C12H22	001129-89-1 49



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057098.D
Acq On : 21 Sep 2015 18:40
Operator : SY/FY
Sample : G3747-02
Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 13 Sample Multiplier: 1

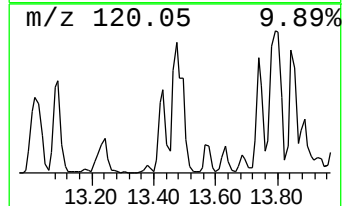
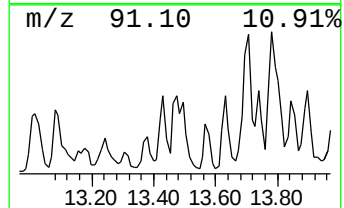
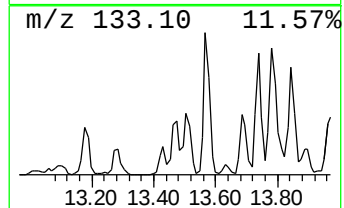
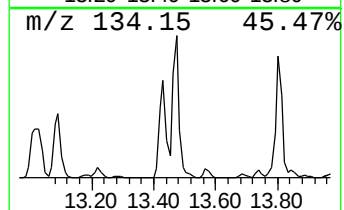
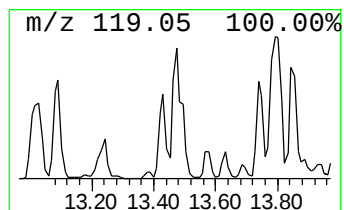
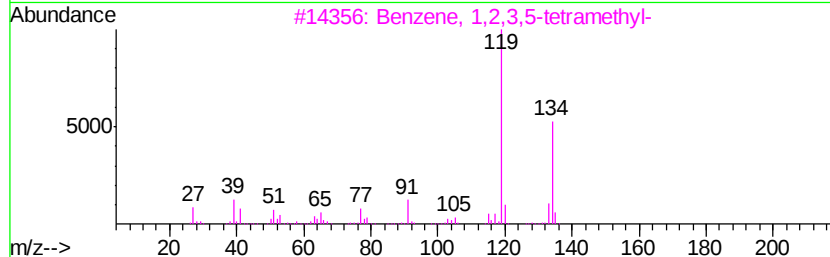
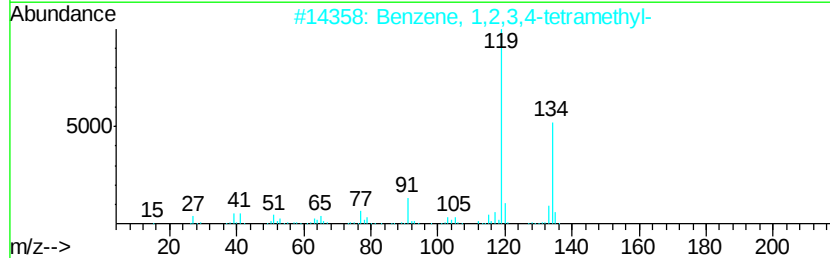
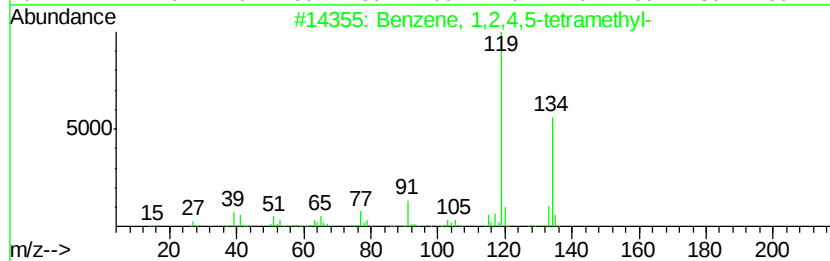
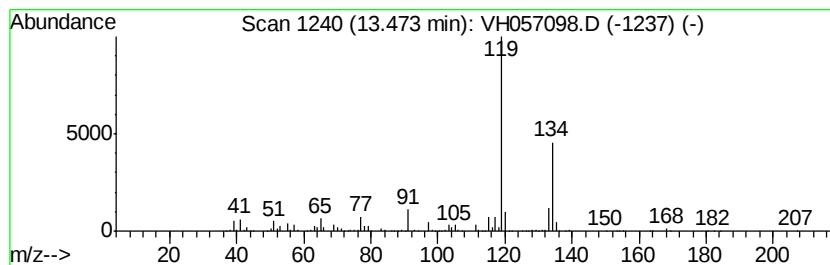
Instrument :
MSVOA_H
ClientSampleId :
TRENCH-1-2

Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	110.00 ug/l	16085600	1,4-Dichlorobenzene-d4	12.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2	95
2	Benzene, 1,2,3,4-tetramethyl-	134 C10H14	000488-23-3	95
3	Benzene, 1,2,3,5-tetramethyl-	134 C10H14	000527-53-7	95
4	Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9	94
5	Benzene, 1-methyl-4-(1-methyleth...	134 C10H14	000099-87-6	93



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057098.D
Acq On : 21 Sep 2015 18:40
Operator : SY/FY
Sample : G3747-02
Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 13 Sample Multiplier: 1

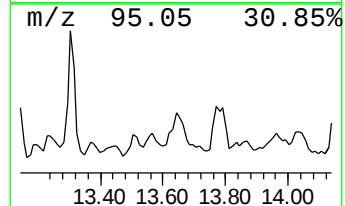
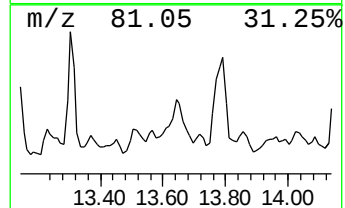
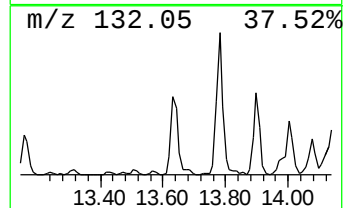
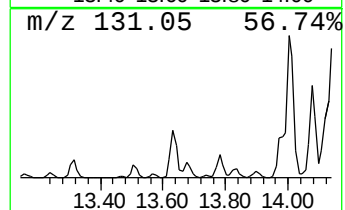
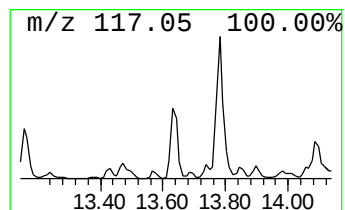
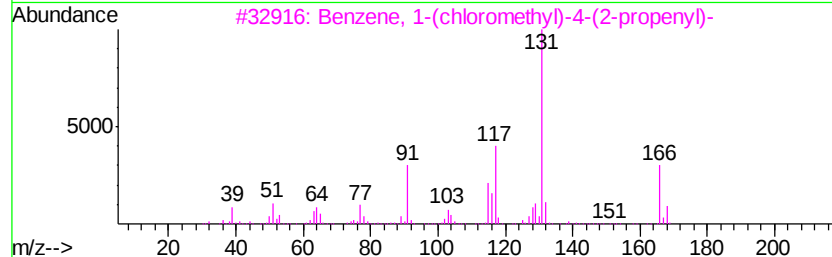
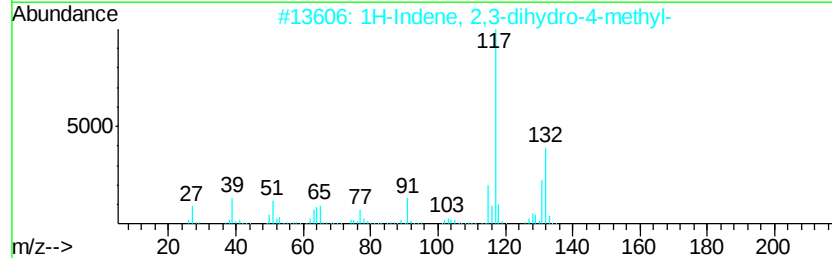
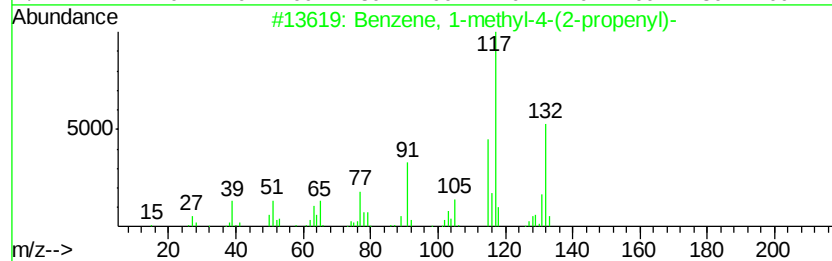
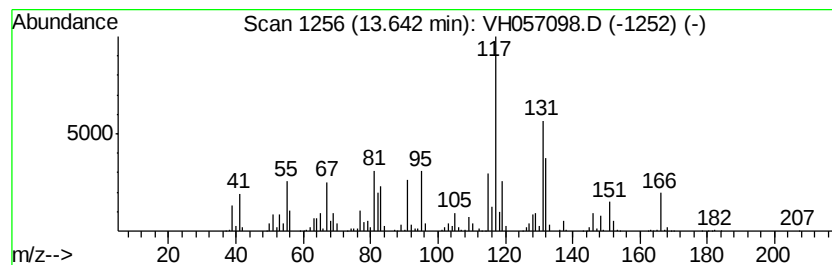
Instrument :
MSVOA_H
ClientSampleId :
TRENCH-1-2

Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-methyl-4-(2-prop... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.64	97.84 ug/l	14307900	1,4-Dichlorobenzene-d4	12.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	60
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	45
3	Benzene, 1-(chloromethyl)-4-(2-p...	166	C10H11Cl	036875-10-2	42
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	41
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	41



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057098.D
Acq On : 21 Sep 2015 18:40
Operator : SY/FY
Sample : G3747-02
Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_H
ClientSampleId :
TRENCH-1-2

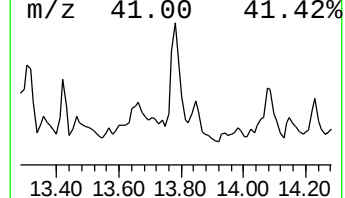
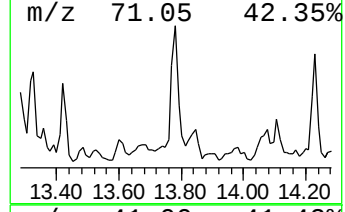
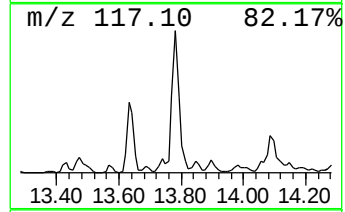
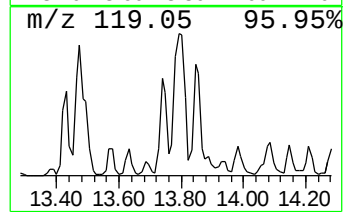
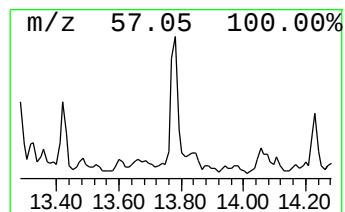
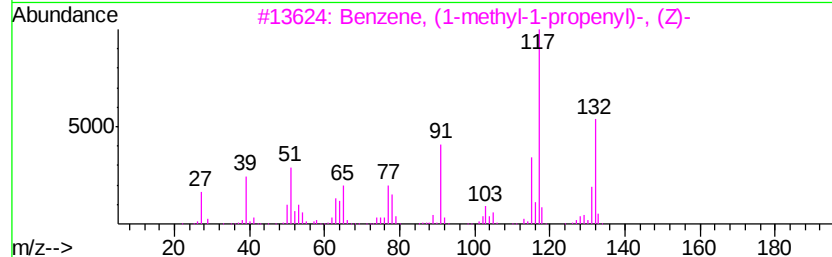
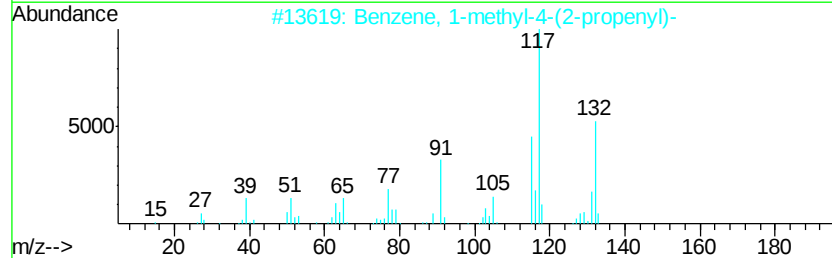
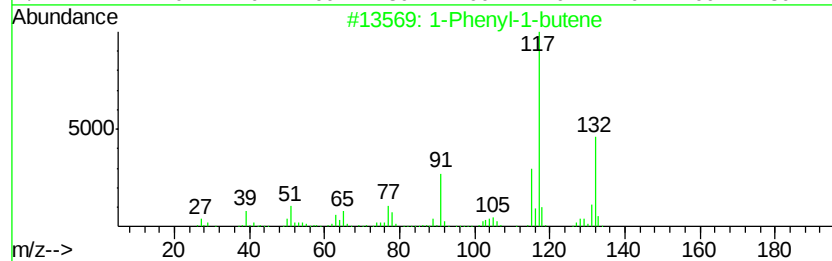
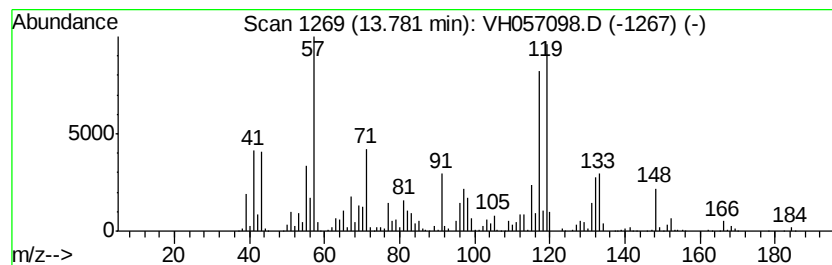
Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1-Phenyl-1-butene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.78	167.08 ug/l	24432100	1,4-Dichlorobenzene-d4	12.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	80
2		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	53
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	44
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	42
5		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	42



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057098.D
Acq On : 21 Sep 2015 18:40
Operator : SY/FY
Sample : G3747-02
Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_H
ClientSampleId :
TRENCH-1-2

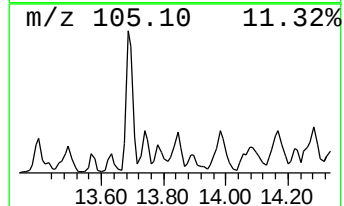
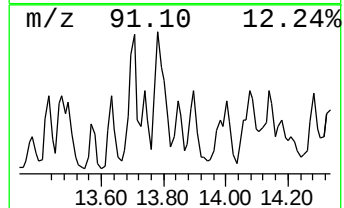
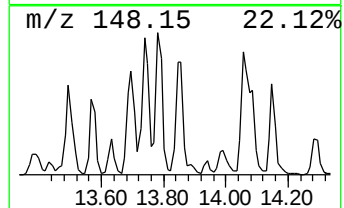
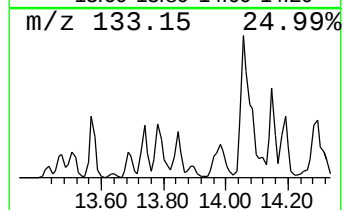
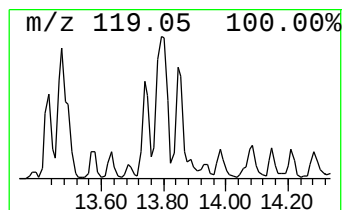
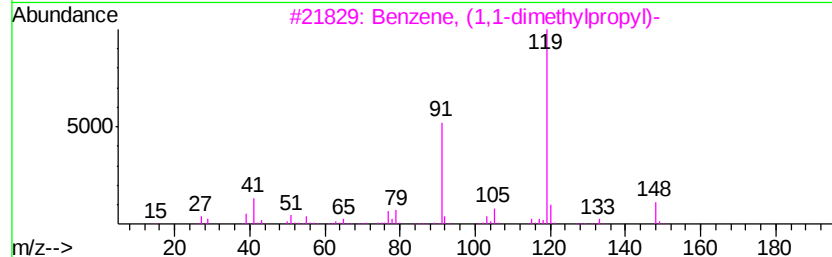
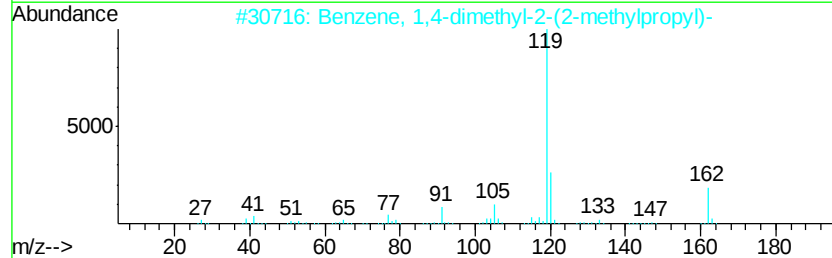
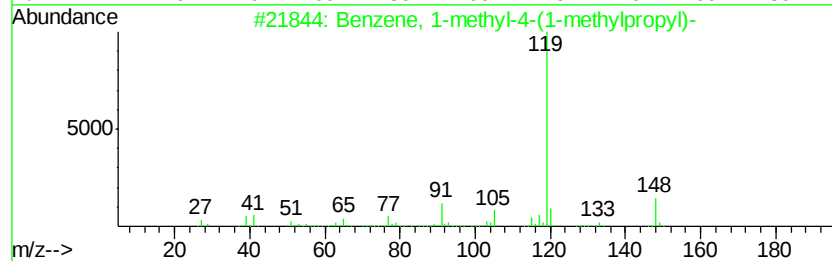
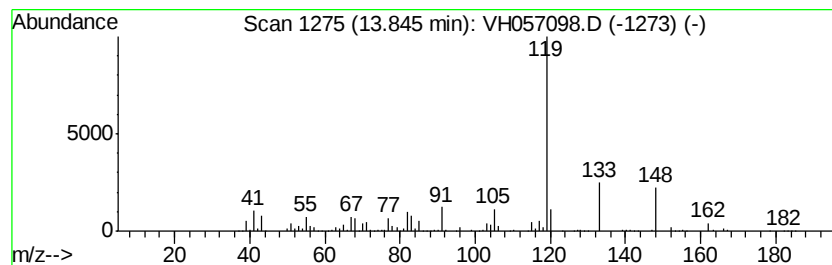
Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 1-methyl-4-(1-meth... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.84	70.75 ug/l	10345700	1,4-Dichlorobenzene-d4	12.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	81
2		Benzene, 1,4-dimethyl-2-(2-methy...	162	C12H18	055669-88-0	70
3		Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	64
4		Benzene, 1-ethyl-4-(2-methylprop...	162	C12H18	100319-40-2	58
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	58



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057098.D
Acq On : 21 Sep 2015 18:40
Operator : SY/FY
Sample : G3747-02
Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_H
ClientSampleId :
TRENCH-1-2

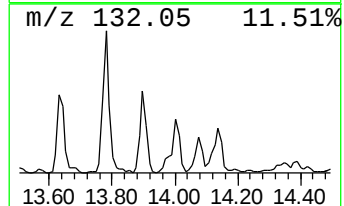
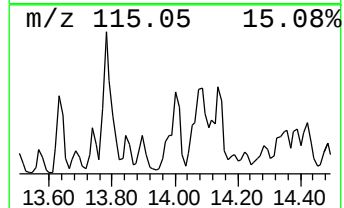
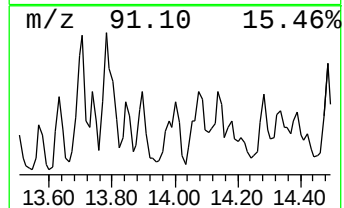
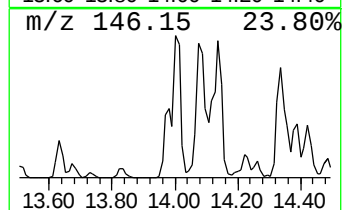
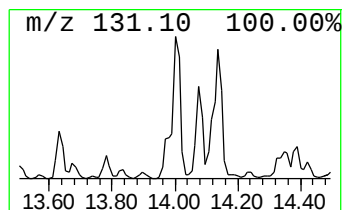
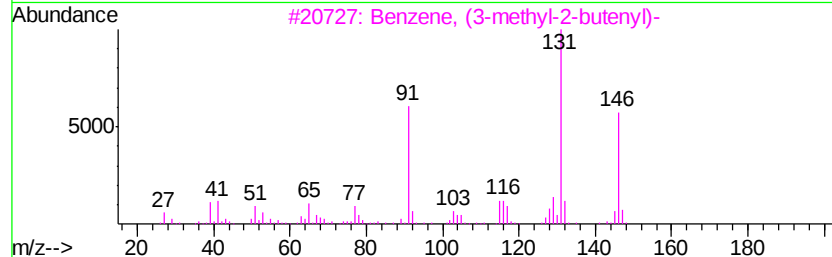
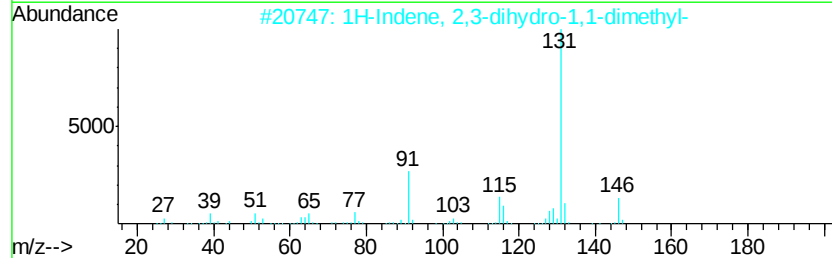
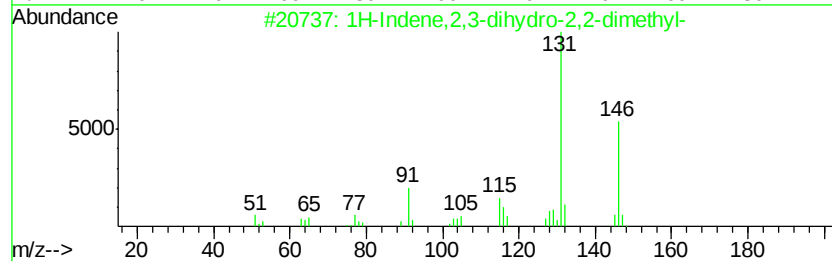
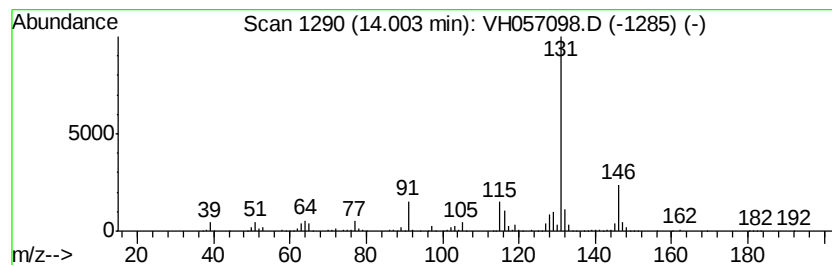
Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1H-Indene,2,3-dihydro-2,2-d... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.00	91.83 ug/l	13428300	1,4-Dichlorobenzene-d4	12.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene,2,3-dihydro-2,2-dimethyl-	146	C11H14	020836-11-7	94
2		1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	91
3		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	91
4		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
5		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057098.D
Acq On : 21 Sep 2015 18:40
Operator : SY/FY
Sample : G3747-02
Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_H
ClientSampleId :
TRENCH-1-2

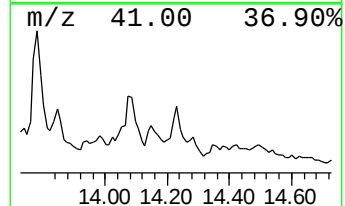
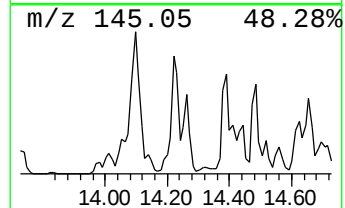
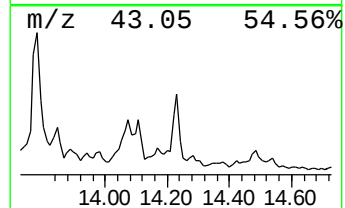
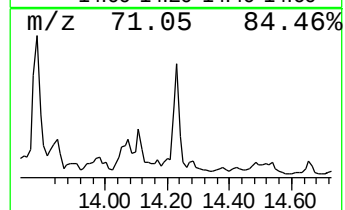
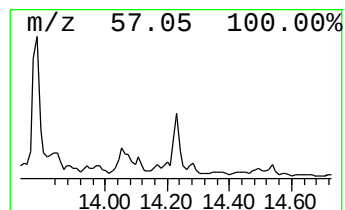
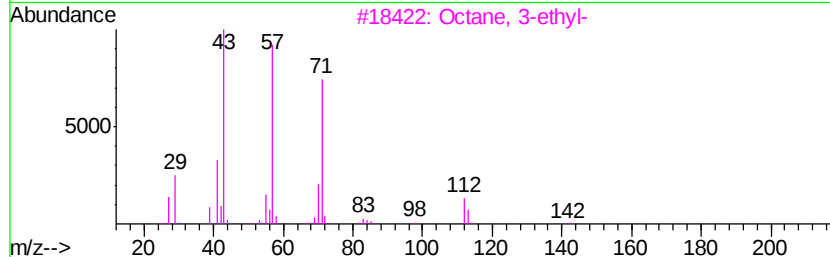
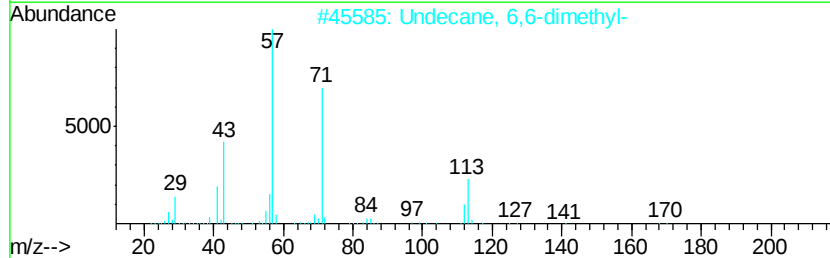
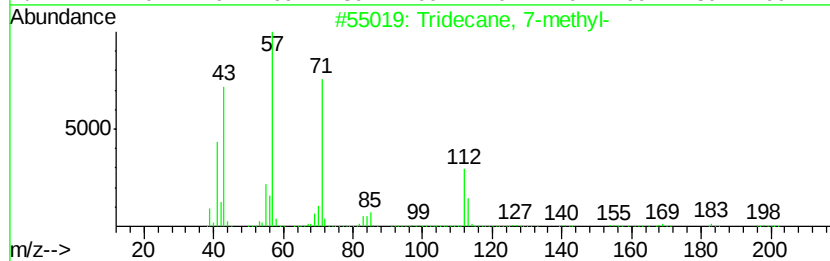
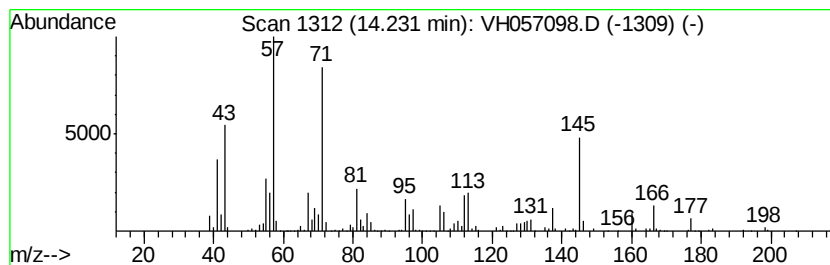
Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 unknown14.23 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.23	70.25 ug/l	10272700	1,4-Dichlorobenzene-d4	12.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 7-methyl-	198	C14H30	026730-14-3	46
2		Undecane, 6,6-dimethyl-	184	C13H28	017312-76-4	38
3		Octane, 3-ethyl-	142	C10H22	005881-17-4	35
4		Octane, 2-bromo-	192	C8H17Br	000557-35-7	35
5		Nonane, 4-methyl-5-propyl-	184	C13H28	062185-55-1	35



Data Path : W:\HPCHEM1\MSVOA_H\DATA\XH092115\
Data File : VH057098.D
Acq On : 21 Sep 2015 18:40
Operator : SY/FY
Sample : G3747-02
Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_H
ClientSampleId :
TRENCH-1-2

Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown12.87	12.87	72.4	ug/l	10581300	4	12.49	7311680	50.0
trans-Decalin, 2-...	13.13	90.5	ug/l	13234000	4	12.49	7311680	50.0
Cyclohexane, pentyl-	13.22	82.9	ug/l	12127600	4	12.49	7311680	50.0
Naphthalene, deca...	13.30	86.5	ug/l	12644700	4	12.49	7311680	50.0
Benzene, 1,2,4,5-...	13.47	110.0	ug/l	16085600	4	12.49	7311680	50.0
Benzene, 1-methyl...	13.64	97.8	ug/l	14307900	4	12.49	7311680	50.0
1-Phenyl-1-butene	13.78	167.1	ug/l	24432100	4	12.49	7311680	50.0
Benzene, 1-methyl...	13.84	70.8	ug/l	10345700	4	12.49	7311680	50.0
1H-Indene, 2,3-dih...	14.00	91.8	ug/l	13428300	4	12.49	7311680	50.0
unknown14.23	14.23	70.3	ug/l	10272700	4	12.49	7311680	50.0