

Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
 Data File : VH057094.D
 Acq On : 21 Sep 2015 16:08
 Operator : SY/FY
 Sample : G3747-02
 Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
 ALS Vial : 9 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	4.056	332	345	359	rBV	919767	2988052	13.48%	1.249%
2	4.583	390	395	404	rVB3	69078	239210	1.08%	0.100%
3	4.814	407	417	429	rBV2	1541769	6210439	28.03%	2.597%
4	5.541	481	486	498	rBV	1818001	5020289	22.66%	2.099%
5	6.971	615	622	631	rVB2	68860	262513	1.18%	0.110%
6	7.435	661	666	667	rBV	122250	228113	1.03%	0.095%
7	7.499	667	672	676	rVV	2872926	7012297	31.65%	2.932%
8	7.562	676	678	688	rVV	473424	1230232	5.55%	0.514%
9	8.227	734	741	746	rBV2	98980	232878	1.05%	0.097%
10	8.406	753	758	765	rVB	193237	533917	2.41%	0.223%
11	8.563	769	773	777	rBV	137353	362446	1.64%	0.152%
12	8.918	800	807	811	rBV2	157920	493263	2.23%	0.206%
13	9.055	816	820	827	rVB	291846	754347	3.40%	0.315%
14	9.233	832	837	841	rBV	292407	649959	2.93%	0.272%
15	9.630	866	875	878	rBV5	80586	267273	1.21%	0.112%
16	9.704	878	882	890	rVV	2369894	5704945	25.75%	2.385%
17	9.832	890	894	900	rVB	966014	2036709	9.19%	0.852%
18	10.062	912	916	922	rVB	750247	1471986	6.64%	0.615%
19	10.166	922	926	934	rVB3	226173	718443	3.24%	0.300%
20	10.294	934	938	944	rBV2	149248	373176	1.68%	0.156%
21	10.451	947	953	957	rBV3	269150	851371	3.84%	0.356%
22	10.610	964	968	976	rBV3	988454	2926530	13.21%	1.224%
23	10.727	976	979	981	rBV	166652	344002	1.55%	0.144%
24	10.769	981	983	991	rVB2	268505	626009	2.83%	0.262%
25	10.895	991	995	998	rVB2	131360	296312	1.34%	0.124%
26	10.991	999	1004	1007	rBV2	485942	1221352	5.51%	0.511%
27	11.064	1007	1011	1013	rBV	843454	1519611	6.86%	0.635%
28	11.105	1013	1015	1018	rVB	815477	1160027	5.24%	0.485%
29	11.211	1018	1025	1029	rVB2	1341582	3042944	13.73%	1.272%
30	11.349	1033	1038	1044	rBV2	2032982	4745181	21.41%	1.984%
31	11.432	1044	1046	1049	rVB2	317932	515845	2.33%	0.216%
32	11.496	1049	1052	1054	rBV	1164147	1826151	8.24%	0.764%
33	11.549	1054	1057	1061	rVB3	429507	927643	4.19%	0.388%
34	11.611	1061	1063	1064	rBV2	272126	364197	1.64%	0.152%

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ClientSampleId :
TRENCH-1-2

Integration Parameters: rteint.p

Integrator: RTE
Smoothing : OFF
Sampling : 1
Start Thrs: 0.001
Stop Thrs : 0
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Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
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35	11.653	1064	1067	1071	rVB	661187	1392297	6.28%	0.582%
36	11.717	1071	1073	1075	rVB	401334	534597	2.41%	0.224%
37	11.758	1075	1077	1080	rVB2	574862	926382	4.18%	0.387%
38	11.832	1080	1084	1086	rBV2	347702	724504	3.27%	0.303%
39	11.896	1086	1090	1091	rBV2	544093	1010883	4.56%	0.423%
40	11.938	1091	1094	1098	rVB	2150226	3502284	15.81%	1.464%
41	12.010	1098	1101	1106	rBV2	561691	1277293	5.76%	0.534%
42	12.091	1106	1109	1111	rBV2	1024602	2080487	9.39%	0.870%
43	12.144	1111	1114	1117	rVB2	4030571	6666313	30.08%	2.787%
44	12.227	1117	1122	1124	rVB4	672729	1316973	5.94%	0.551%
45	12.270	1124	1126	1128	rBV2	571119	733978	3.31%	0.307%
46	12.321	1128	1131	1133	rBV	1370495	2029964	9.16%	0.849%
47	12.362	1133	1135	1136	rBV	974596	1117618	5.04%	0.467%
48	12.394	1136	1138	1143	rVB2	1586964	2488583	11.23%	1.040%
49	12.479	1143	1146	1149	rBV	2902700	4378552	19.76%	1.831%
50	12.532	1149	1151	1155	rVB2	1956831	3078931	13.89%	1.287%
51	12.596	1155	1157	1159	rBV3	235300	413816	1.87%	0.173%
52	12.669	1159	1164	1166	rBV3	766780	2278654	10.28%	0.953%
53	12.722	1166	1169	1172	rVV3	1272441	2877091	12.98%	1.203%
54	12.775	1172	1174	1178	rVB3	1946790	3351117	15.12%	1.401%
55	12.871	1180	1183	1187	rBV4	1380921	4050225	18.28%	1.693%
56	12.935	1187	1189	1194	rVB2	2471492	5101573	23.02%	2.133%
57	13.030	1194	1198	1201	rBV3	1735086	5654778	25.52%	2.364%
58	13.082	1201	1203	1205	rVV	2120052	2798226	12.63%	1.170%
59	13.125	1205	1207	1211	rVB3	2587094	4708824	21.25%	1.969%
60	13.221	1214	1216	1219	rBV3	3464056	6236889	28.15%	2.608%
61	13.273	1219	1221	1222	rVV2	761278	1254670	5.66%	0.525%
62	13.305	1222	1224	1227	rVB2	3640337	5655267	25.52%	2.364%
63	13.369	1227	1230	1233	rVB4	510912	1089109	4.91%	0.455%
64	13.423	1233	1235	1237	rBV	2550050	3347672	15.11%	1.400%
65	13.465	1237	1239	1246	rVB2	3114716	7524923	33.96%	3.146%
66	13.571	1246	1249	1251	rBV	1459040	2303052	10.39%	0.963%
67	13.633	1252	1255	1258	rBV2	2312586	4575217	20.65%	1.913%
68	13.686	1258	1260	1263	rVV3	1171829	2063210	9.31%	0.863%
69	13.738	1263	1265	1267	rVV	1830746	2483710	11.21%	1.038%
70	13.781	1267	1269	1273	rVV2	6195975	11650639	52.58%	4.871%
71	13.845	1273	1275	1277	rVB	2343999	2913620	13.15%	1.218%

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Integration Parameters: rteint.p

Integrator: RTE
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72	13.898	1277	1280	1282	rVB	1604135	2611298	11.78%	1.092%
73	13.940	1282	1284	1285	rBV2	203248	336873	1.52%	0.141%
74	14.003	1285	1290	1293	rBV2	2302858	5705421	25.75%	2.385%
75	14.084	1293	1298	1302	rBV5	3524569	9843986	44.42%	4.116%
76	14.146	1302	1304	1309	rVB4	2705876	4559003	20.57%	1.906%
77	14.230	1309	1312	1315	rBV3	1124641	1791114	8.08%	0.749%
78	14.281	1315	1317	1320	rVB4	1081193	1633524	7.37%	0.683%
79	14.375	1320	1326	1329	rBV	12319536	22158921	100.00%	9.265%
80	14.481	1335	1336	1338	rBV	756033	1017368	4.59%	0.425%
81	14.513	1338	1339	1346	rVB3	1130251	1501356	6.78%	0.628%
82	14.657	1350	1353	1359	rVB3	2432547	4501260	20.31%	1.882%
83	14.740	1359	1361	1363	rBV	614676	990438	4.47%	0.414%
84	14.781	1363	1365	1371	rVB4	566605	1569586	7.08%	0.656%
85	14.908	1375	1377	1382	rVB2	788937	1355120	6.12%	0.567%
86	15.025	1385	1388	1390	rVB	455993	651684	2.94%	0.272%
87	15.079	1390	1393	1399	rVB5	417894	971679	4.39%	0.406%
88	15.186	1400	1403	1407	rVB	2177658	3328771	15.02%	1.392%
89	15.313	1412	1415	1422	rVB	1183023	1900119	8.57%	0.794%

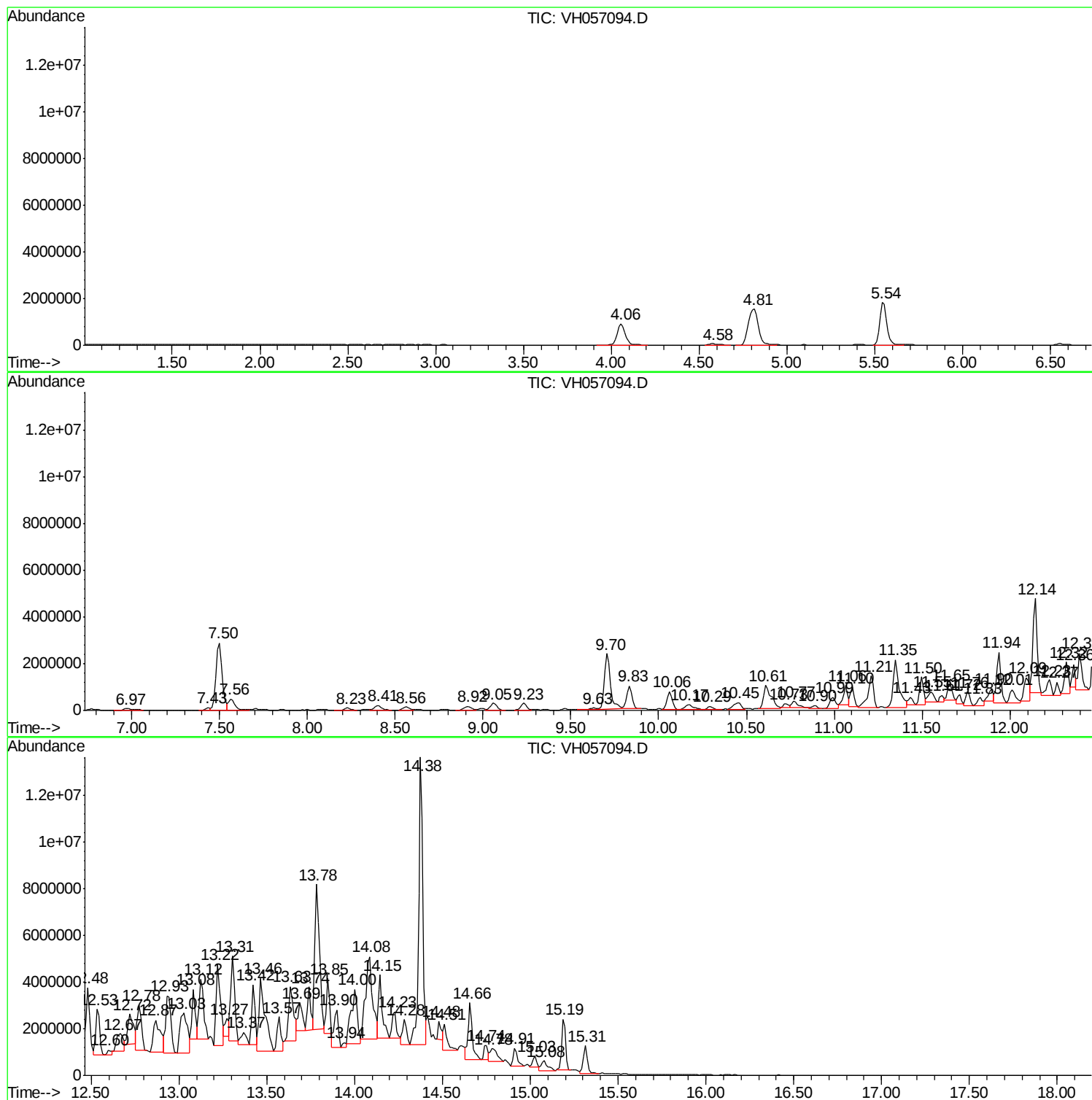
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Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

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



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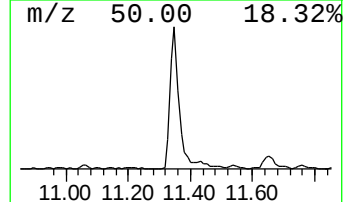
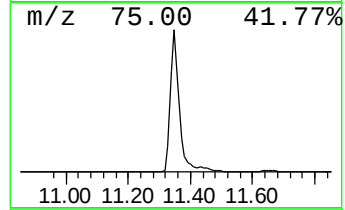
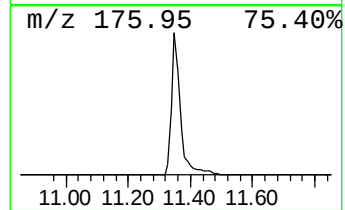
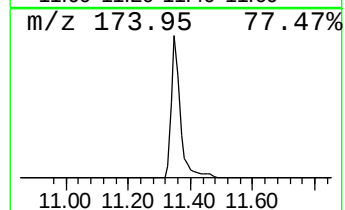
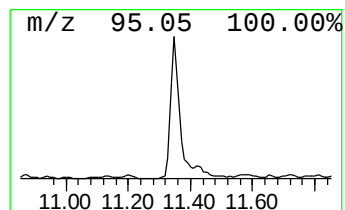
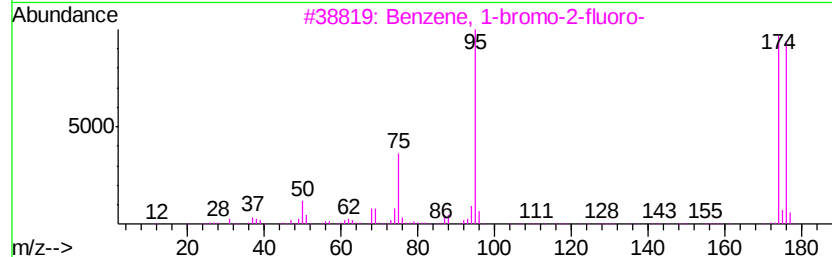
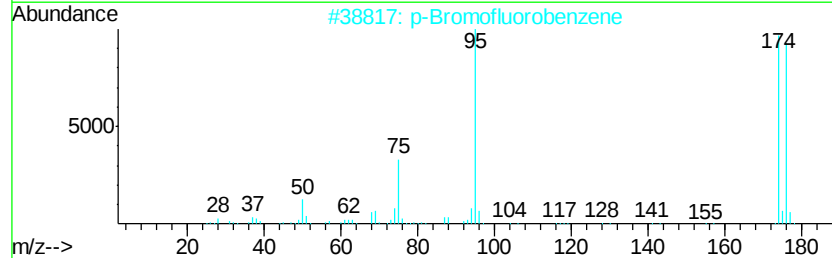
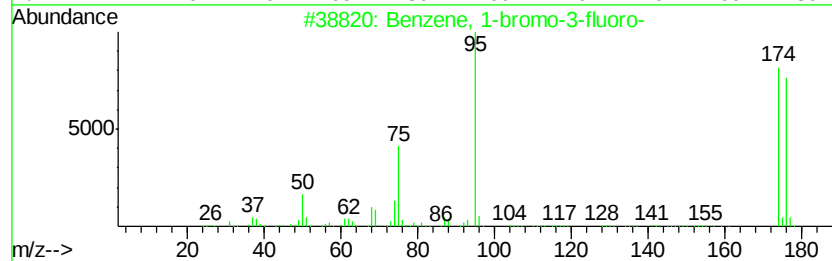
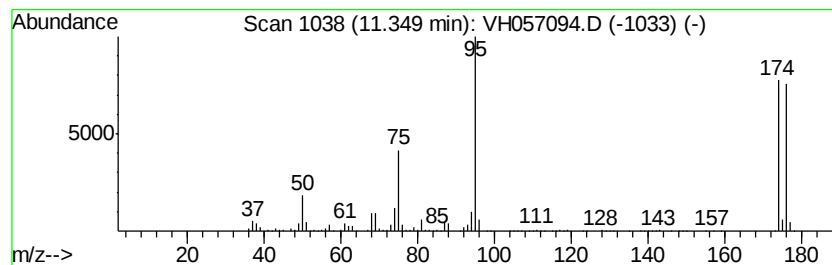
Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-bromo-3-fluoro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.35	54.19 ug/l	4745180	1,4-Dichlorobenzene-d4	12.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-bromo-3-fluoro-	174	C6H4BrF	001073-06-9	98
2		p-Bromofluorobenzene	174	C6H4BrF	000460-00-4	97
3		Benzene, 1-bromo-2-fluoro-	174	C6H4BrF	001072-85-1	96
4		4(1H)-Pyrimidinone, 5-bromo-	174	C4H3BrN2O	019808-30-1	37
5		Benzaldehyde, 2,6-difluoro-3,4-d...	174	C7H4F2O3	152434-98-5	14



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Instrument :
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ClientSampleId :
TRENCH-1-2

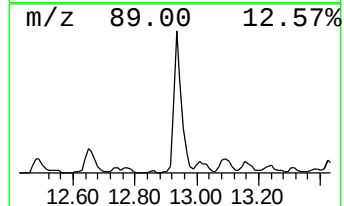
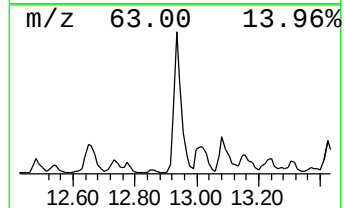
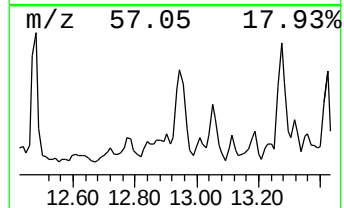
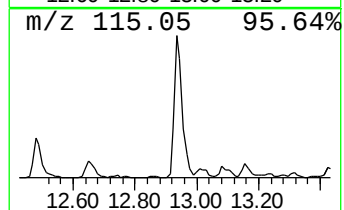
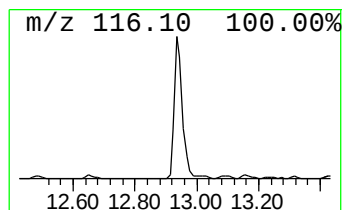
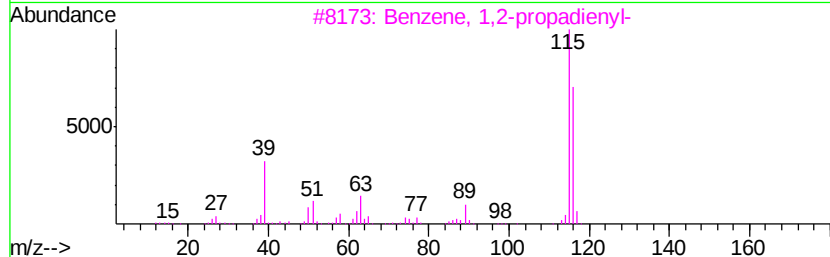
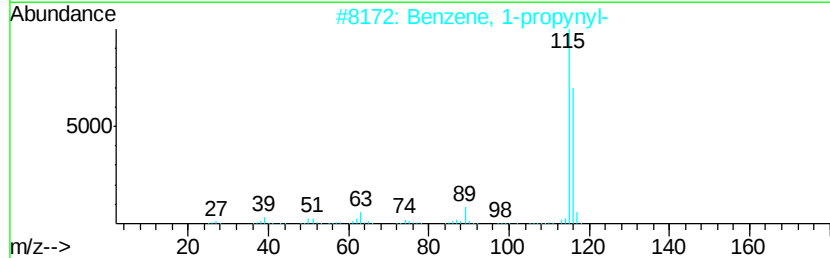
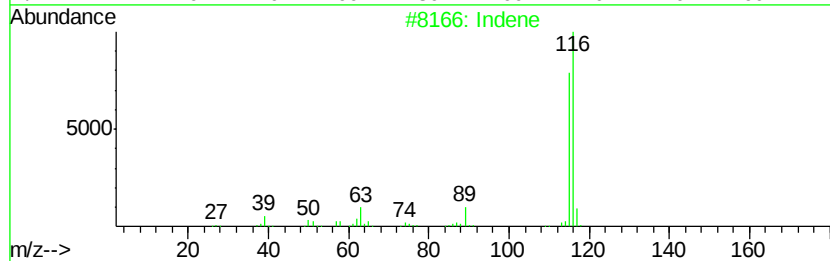
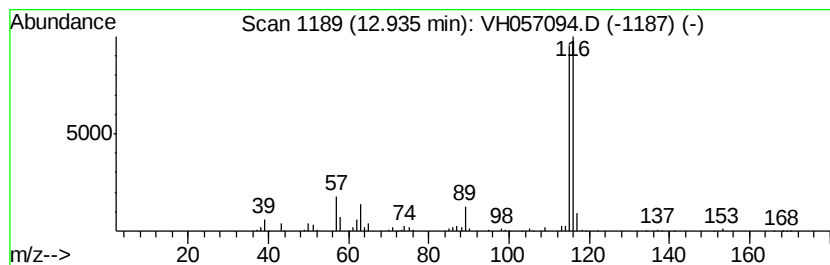
Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

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

Peak Number 2 Indene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.93	58.26 ug/l	5101570	1,4-Dichlorobenzene-d4	12.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indene	116	C9H8	000095-13-6	97
2		Benzene, 1-propynyl-	116	C9H8	000673-32-5	90
3		Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	90
4		Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	83
5		Tetracyclo[2.2.1.0(2,6).0(3,5)]h...	116	C9H8	081787-74-8	43



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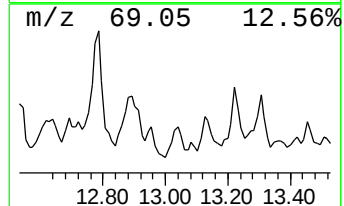
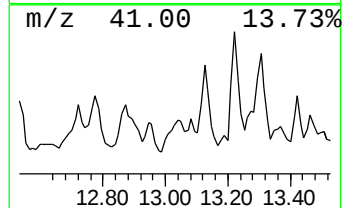
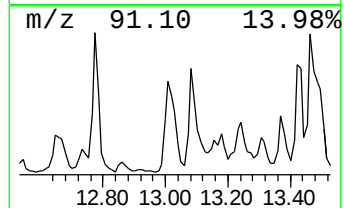
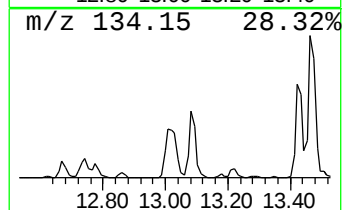
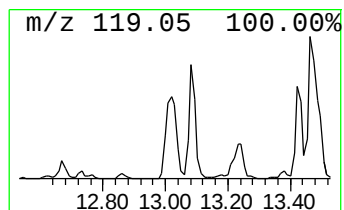
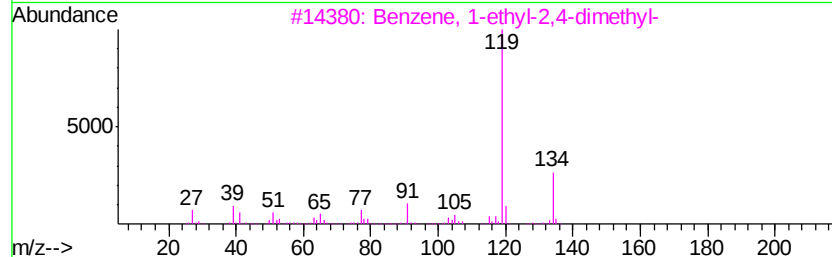
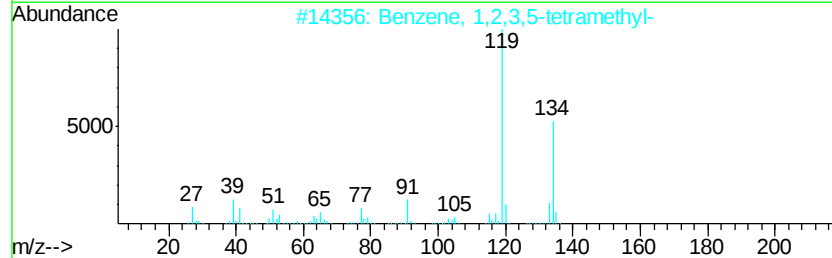
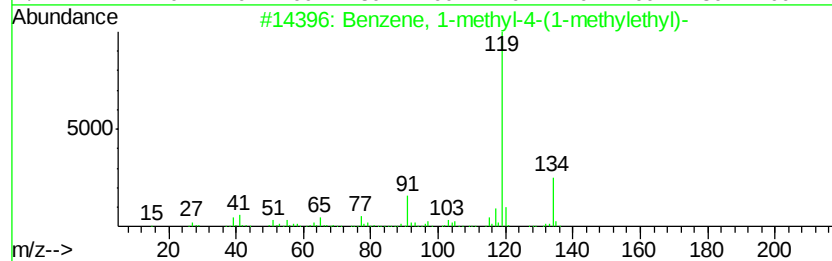
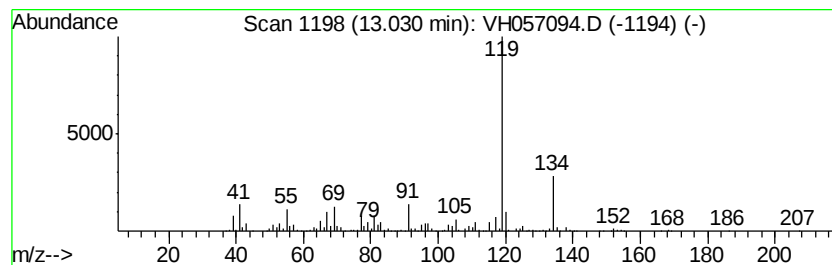
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-methyl-4-(1-meth... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.03	64.57 ug/l	5654780	1,4-Dichlorobenzene-d4	12.48
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1-methyl-4-(1-methyleth...	134 C10H14	000099-87-6	87
2	Benzene, 1,2,3,5-tetramethyl-	134 C10H14	000527-53-7	81
3	Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9	81
4	Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4	81
5	Benzene, 1-methyl-2-(1-methyleth...	134 C10H14	000527-84-4	81



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Acq On : 21 Sep 2015 16:08
Operator : SY/FY
Sample : G3747-02
Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_H
ClientSampleId :
TRENCH-1-2

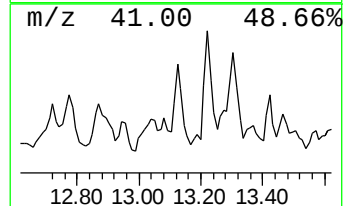
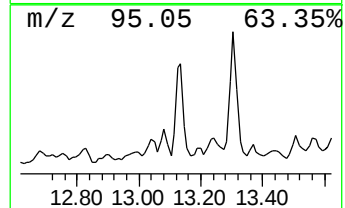
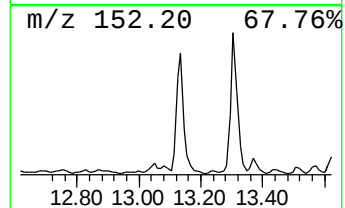
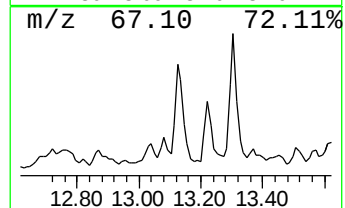
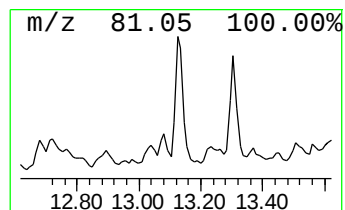
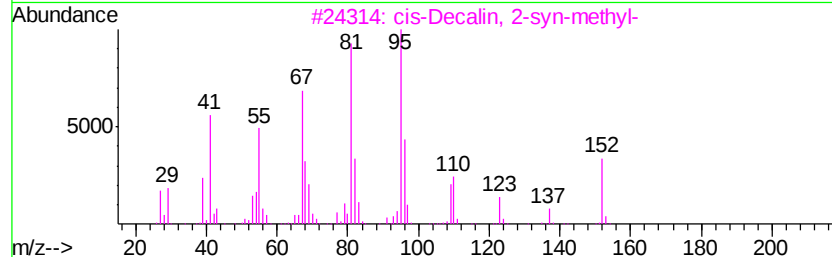
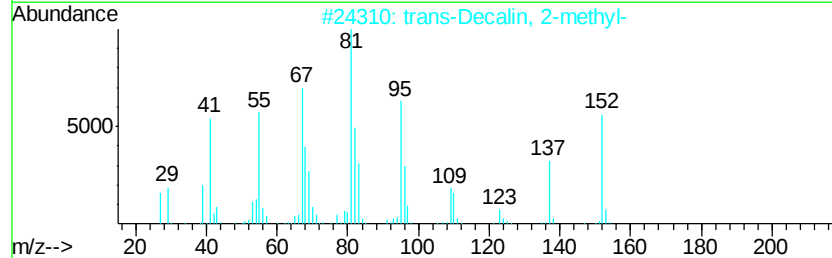
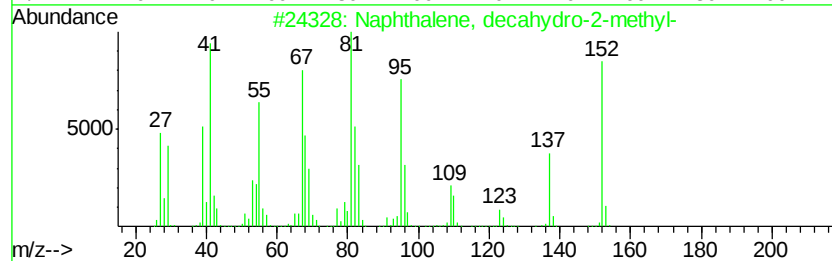
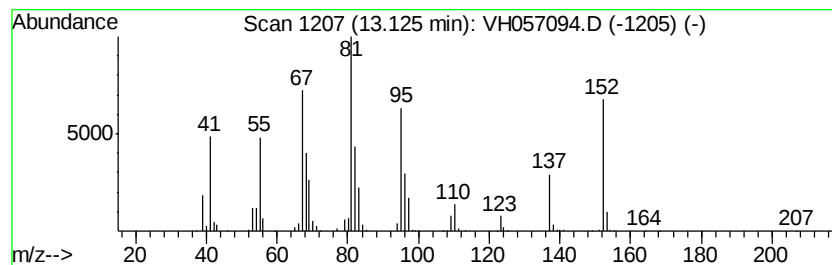
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Quant Title : SW846 8260

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

Peak Number 4 Naphthalene, decahydro-2-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	53.77 ug/l	4708820	1,4-Dichlorobenzene-d4	12.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	94
2		trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	94
3		cis-Decalin, 2-syn-methyl-	152	C11H20	1000155-85-6	72
4		Cyclohexanone, 5-methyl-2-(1-met...	152	C10H16O	015932-80-6	68
5		6a-Methyl-hexahdropentalene-1,6...	152	C9H12O2	092485-86-4	64



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057094.D
Acq On : 21 Sep 2015 16:08
Operator : SY/FY
Sample : G3747-02
Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_H
ClientSampleId :
TRENCH-1-2

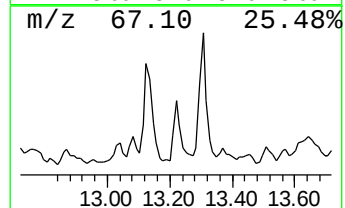
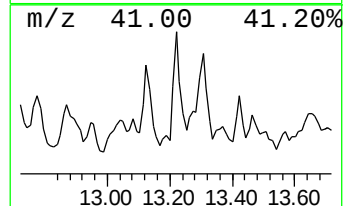
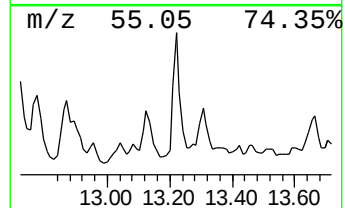
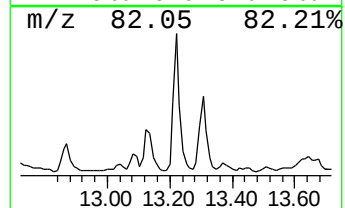
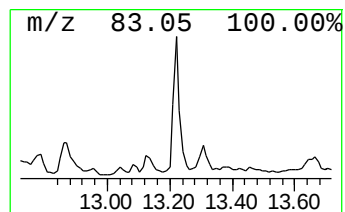
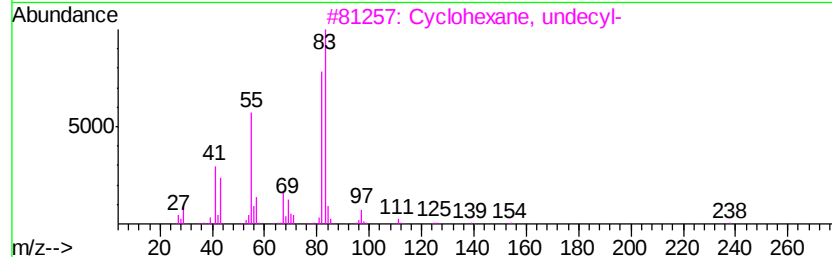
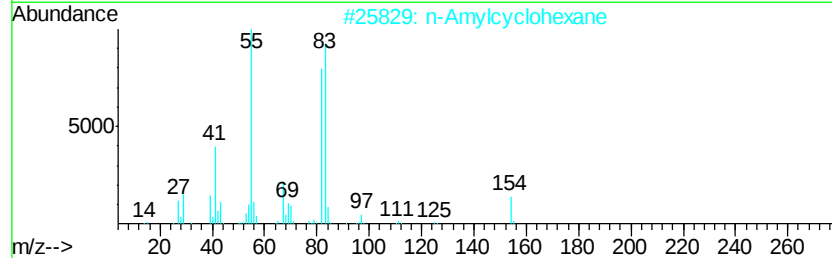
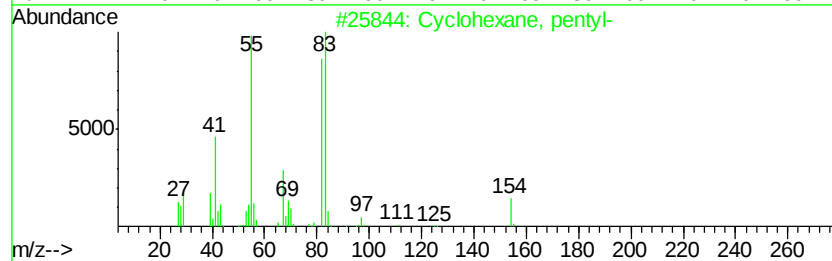
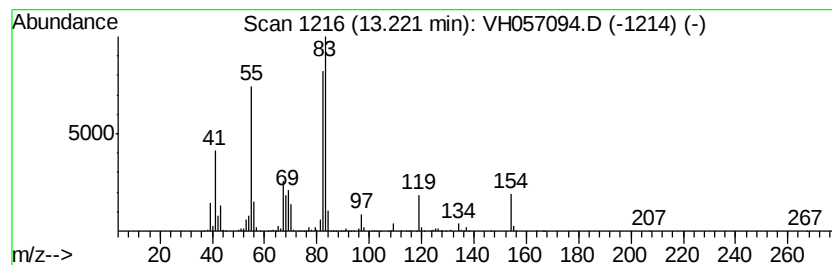
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Quant Title : SW846 8260

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

Peak Number 5 Cyclohexane, pentyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	71.22 ug/l	6236890	1,4-Dichlorobenzene-d4	12.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, pentyl-	154	C11H22	004292-92-6	90
2		n-Amylcyclohexane	154	C11H22	029949-27-7	87
3		Cyclohexane, undecyl-	238	C17H34	054105-66-7	72
4		Cyclohexane, octyl-	196	C14H28	001795-15-9	72
5		n-Pentadecylcyclohexane	294	C21H42	006006-95-7	72



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057094.D
Acq On : 21 Sep 2015 16:08
Operator : SY/FY
Sample : G3747-02
Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 9 Sample Multiplier: 1

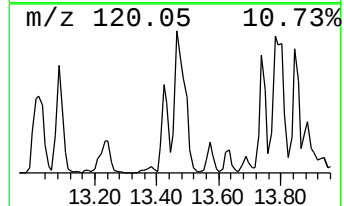
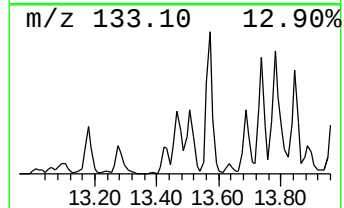
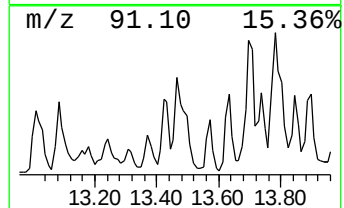
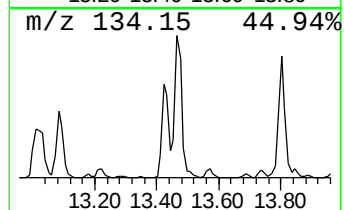
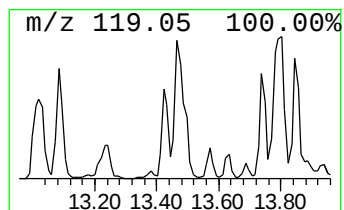
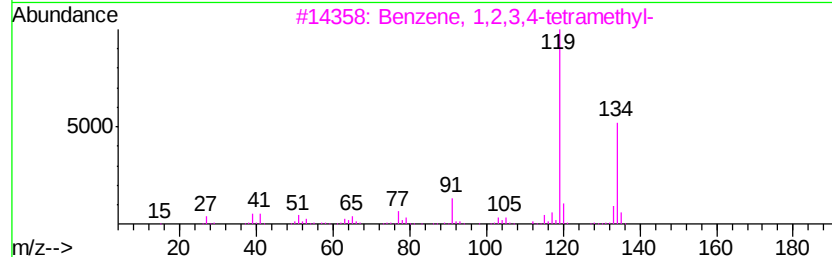
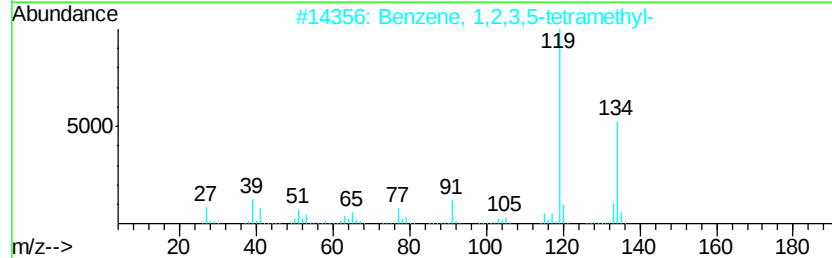
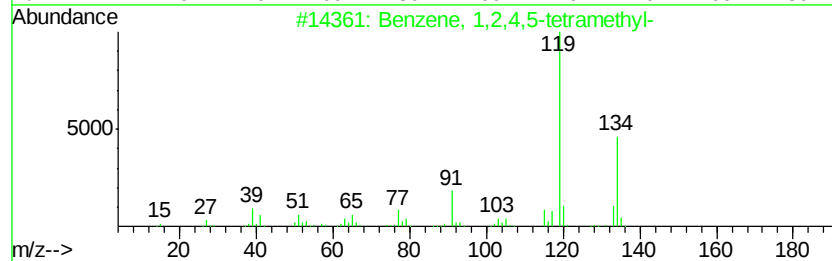
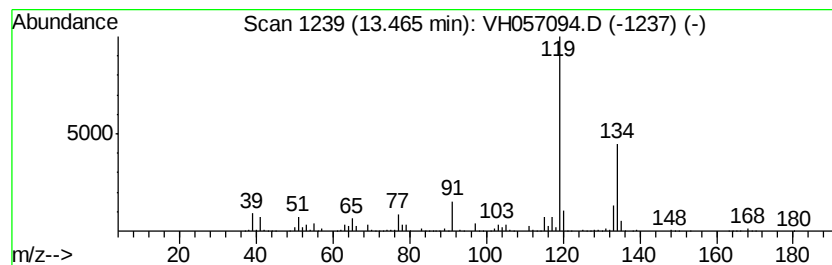
Instrument :
MSVOA_H
ClientSampled :
TRENCH-1-2

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 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.46	85.93 ug/l	7524920	1,4-Dichlorobenzene-d4	12.48	
Hit# of	5	Tentative ID	MW MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057094.D
Acq On : 21 Sep 2015 16:08
Operator : SY/FY
Sample : G3747-02
Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_H
ClientSampleId :
TRENCH-1-2

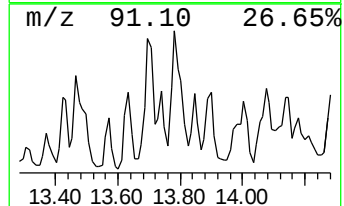
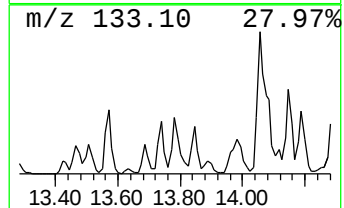
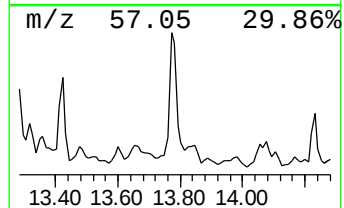
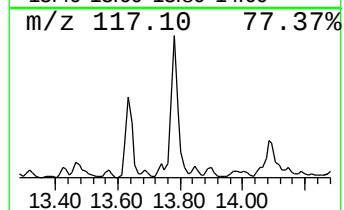
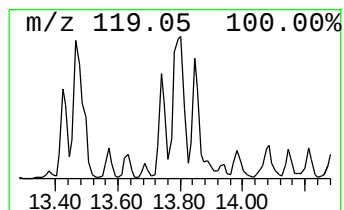
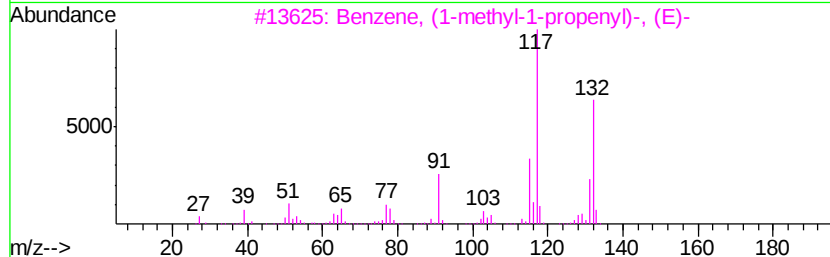
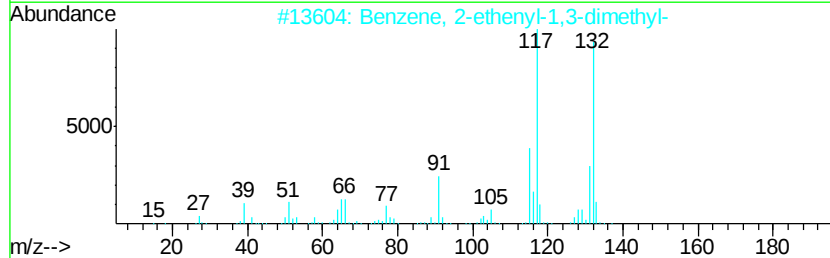
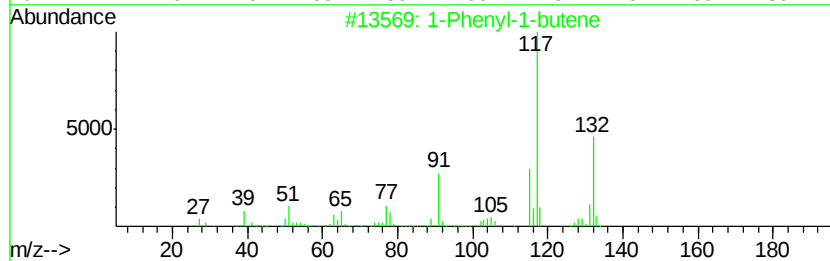
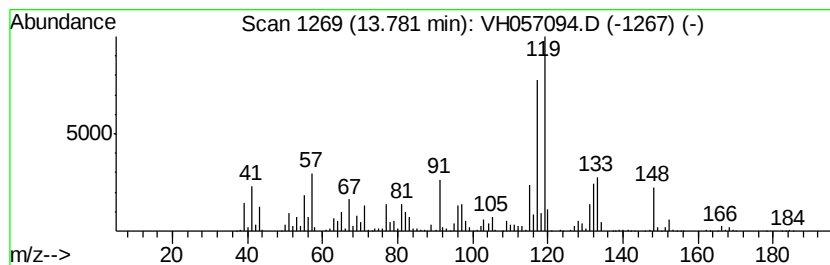
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Quant Title : SW846 8260

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

Peak Number 8 1-Phenyl-1-butene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.78	133.04 ug/l	11650600	1,4-Dichlorobenzene-d4	12.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	50
2		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	50
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	50
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	50
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	50



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057094.D
Acq On : 21 Sep 2015 16:08
Operator : SY/FY
Sample : G3747-02
Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_H
ClientSampleId :
TRENCH-1-2

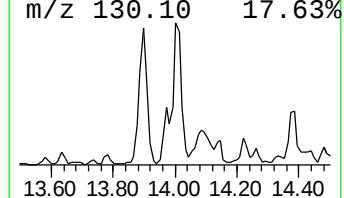
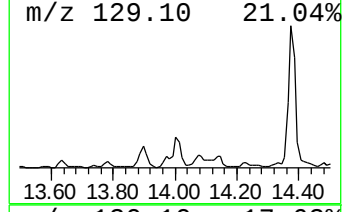
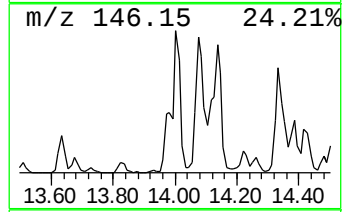
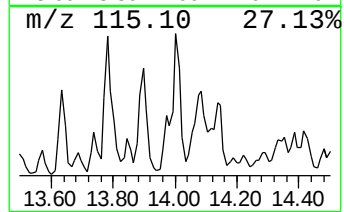
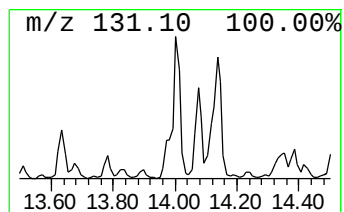
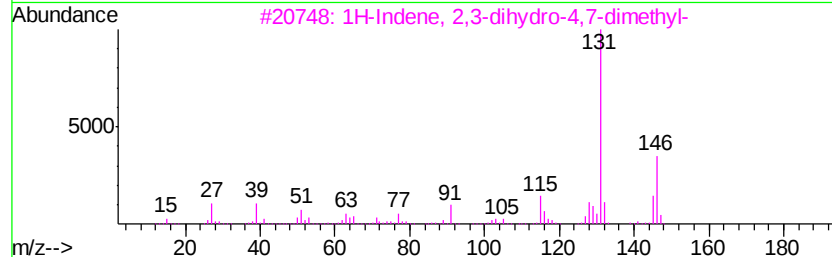
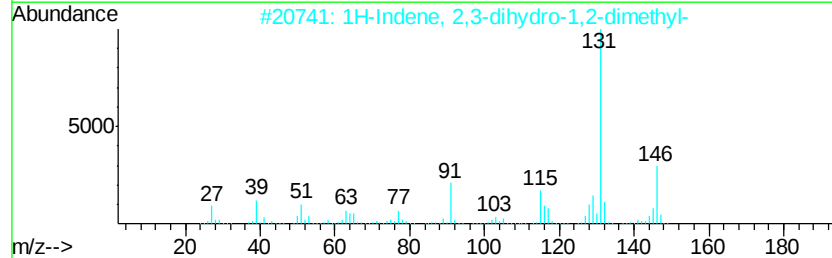
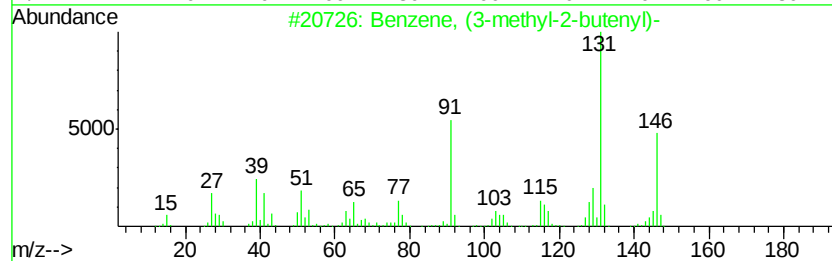
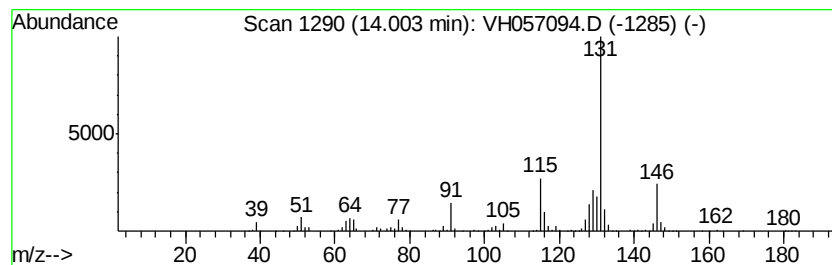
Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, (3-methyl-2-butenyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.00	65.15 ug/l	5705420	1,4-Dichlorobenzene-d4	12.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	90
2		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	87
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	86
4		.alpha.,.beta.,.beta.-Trimethyls...	146	C11H14	000769-57-3	83
5		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	81



Data Path : W:\HPCHEM1\MSVOA H\DATA\XH092115\
Data File : VH057094.D
Acq On : 21 Sep 2015 16:08
Operator : SY/FY
Sample : G3747-02
Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_H
ClientSampled :
TRENCH-1-2

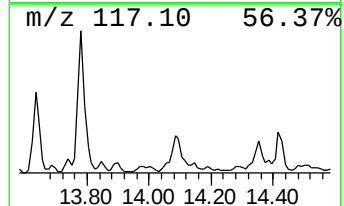
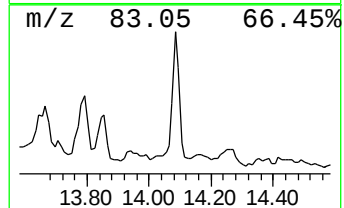
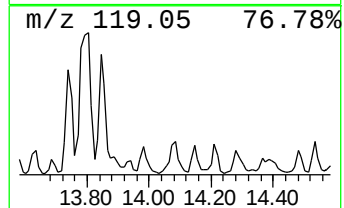
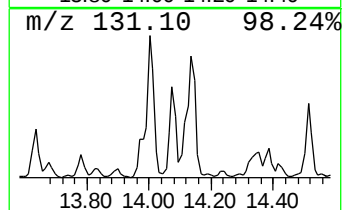
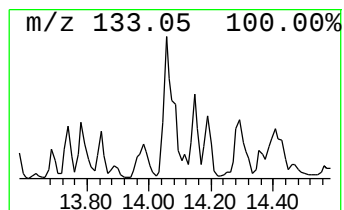
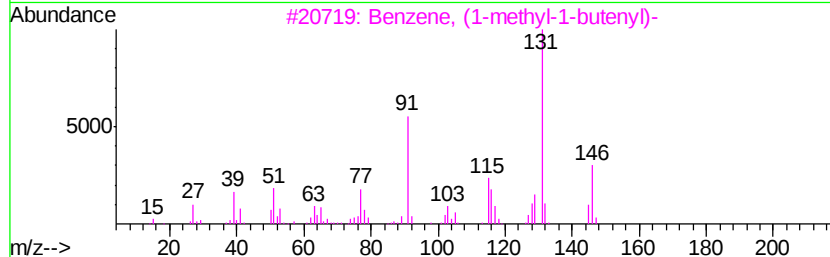
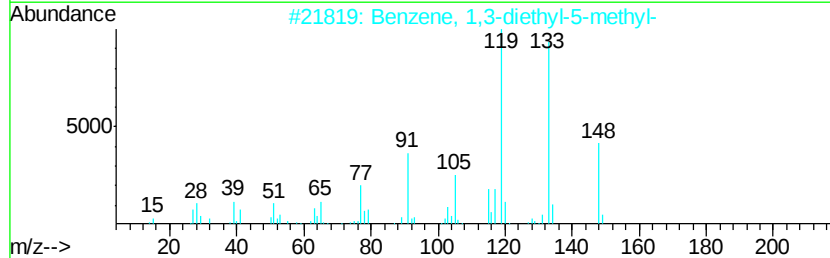
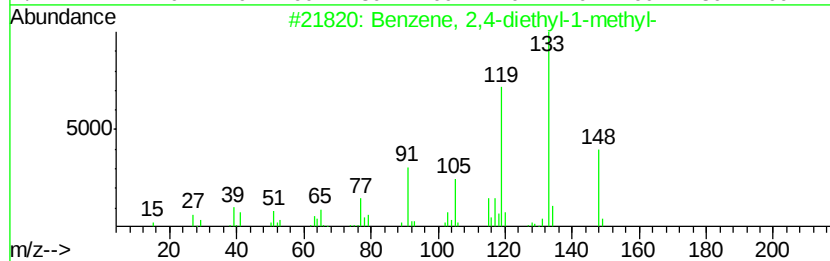
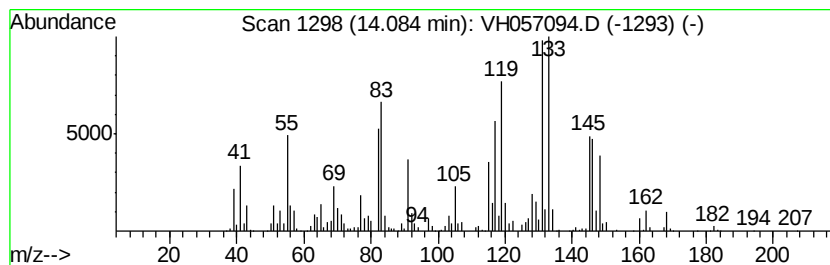
Quant Method : W:\HPCHEM1\MSVOA_H\METHODS\82H091715W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2,4-diethyl-1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.08	112.41 ug/l	9843990	1,4-Dichlorobenzene-d4	12.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2,4-diethyl-1-methyl-	148	C11H16	001758-85-6	80
2		Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	42
3		Benzene, (1-methyl-1-butenvl)-	146	C11H14	053172-84-2	35
4		Benzene, 1,4-diethyl-2-methyl-	148	C11H16	013632-94-5	30
5		Ethanone, 1-(3,4-dimethylphenyl)-	148	C10H12O	003637-01-2	25



Data Path : W:\HPCHEM1\MSVOA_H\DATA\VH092115\
Data File : VH057094.D
Acq On : 21 Sep 2015 16:08
Operator : SY/FY
Sample : G3747-02
Misc : 6.55/10mL/100uL/5mL/MSVOA_H/MEOH
ALS Vial : 9 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
Benzene, 1-bromo-...	11.35	54.2	ug/l	4745180	4	12.48	4378550	50.0
Indene	12.93	58.3	ug/l	5101570	4	12.48	4378550	50.0
Benzene, 1-methyl...	13.03	64.6	ug/l	5654780	4	12.48	4378550	50.0
Naphthalene, deca...	13.12	53.8	ug/l	4708820	4	12.48	4378550	50.0
Cyclohexane, pentyl-	13.22	71.2	ug/l	6236890	4	12.48	4378550	50.0
Benzene, 1,2,4,5-...	13.46	85.9	ug/l	7524920	4	12.48	4378550	50.0
1-Phenyl-1-butene	13.78	133.0	ug/l	11650600	4	12.48	4378550	50.0
Benzene, (3-methy...	14.00	65.2	ug/l	5705420	4	12.48	4378550	50.0
Benzene, 2,4-diet...	14.08	112.4	ug/l	9843990	4	12.48	4378550	50.0