

Data Path : W:\HPCHEM1\MSVOA_D\Data\VD092115\
 Data File : VD046820.D
 Acq On : 21 Sep 2015 21:08
 Operator : FY/SY
 Sample : G3727-03
 Misc : 6.18gm/5mL/MSVOA_D/SOIL
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 MMB6B

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 : W:\HPCHEM1\MSVOA_D\METHOD\82D091815S.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.237	69	72	73	rBV	47998	51991	4.13%	0.281%
2	1.266	73	75	84	rVB	552418	1004790	79.75%	5.428%
3	2.154	160	165	170	rVB2	23308	41734	3.31%	0.225%
4	2.410	187	191	195	rBV	40053	71985	5.71%	0.389%
5	2.538	202	204	210	rVB2	8858	19010	1.51%	0.103%
6	2.942	237	245	248	rBV	7799	18741	1.49%	0.101%
7	3.514	294	303	312	rBV3	48216	217725	17.28%	1.176%
8	3.879	335	340	347	rVB	26818	72974	5.79%	0.394%
9	4.155	361	368	375	rVB5	5176	21386	1.70%	0.116%
10	4.303	375	383	390	rBV	71213	188167	14.93%	1.017%
11	4.549	403	408	413	rBV4	8201	26501	2.10%	0.143%
12	5.229	471	477	483	rBV	79727	220788	17.52%	1.193%
13	5.968	547	552	558	rVB	13879	37400	2.97%	0.202%
14	6.195	569	575	578	rBV	176802	450700	35.77%	2.435%
15	6.284	578	584	590	rVB2	344258	1052929	83.57%	5.688%
16	6.511	601	607	612	rBV2	38551	111422	8.84%	0.602%
17	6.570	612	613	619	rVB5	10425	24205	1.92%	0.131%
18	6.698	619	626	635	rBV	131008	306757	24.35%	1.657%
19	6.856	635	642	647	rVV3	34522	102286	8.12%	0.553%
20	6.944	647	651	655	rVV2	24564	62698	4.98%	0.339%
21	7.033	655	660	671	rVB2	40837	120806	9.59%	0.653%
22	7.210	671	678	692	rBV	120170	294804	23.40%	1.593%
23	7.417	692	699	710	rBV	429873	1065793	84.59%	5.758%
24	8.068	754	765	772	rBV	152712	473781	37.60%	2.560%
25	8.176	772	776	783	rVB4	7938	21634	1.72%	0.117%
26	8.344	788	793	799	rVB3	19795	48954	3.89%	0.264%
27	8.482	799	807	813	rBV2	14490	43503	3.45%	0.235%
28	8.620	818	821	825	rVV2	7488	19098	1.52%	0.103%
29	8.699	825	829	837	rVB3	19794	54671	4.34%	0.295%
30	8.866	841	846	851	rBV2	10962	25009	1.98%	0.135%
31	8.955	851	855	859	rVV3	7689	20010	1.59%	0.108%
32	9.064	861	866	876	rVB2	58673	169878	13.48%	0.918%
33	9.202	876	880	884	rBV5	8048	24865	1.97%	0.134%
34	9.280	884	888	897	rVB	26621	69924	5.55%	0.378%

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35	9.409	897	901	911	rVB2	26865	74472	5.91%	0.402%
36	9.596	914	920	929	rBV	512467	1259976	100.00%	6.807%
37	9.704	929	931	936	rVB2	15209	34218	2.72%	0.185%
38	9.823	936	943	944	rBV4	12217	28454	2.26%	0.154%
39	9.862	944	947	950	rVV5	12183	39020	3.10%	0.211%
40	9.911	950	952	954	rVV3	14540	29538	2.34%	0.160%
41	9.970	954	958	963	rVB	121384	266694	21.17%	1.441%
42	10.148	966	976	985	rVB4	27202	107109	8.50%	0.579%
43	10.315	987	993	998	rBV2	26785	73618	5.84%	0.398%
44	10.651	1022	1027	1032	rBV	21926	47924	3.80%	0.259%
45	10.789	1037	1041	1044	rVV2	13756	33510	2.66%	0.181%
46	10.848	1044	1047	1052	rVV6	18738	58881	4.67%	0.318%
47	10.936	1052	1056	1060	rVB2	49307	99522	7.90%	0.538%
48	11.015	1060	1064	1068	rBV2	45062	97263	7.72%	0.525%
49	11.173	1078	1080	1083	rVV3	9962	19230	1.53%	0.104%
50	11.341	1088	1097	1105	rVB4	33875	132185	10.49%	0.714%
51	11.479	1105	1111	1115	rBV	43823	96112	7.63%	0.519%
52	11.557	1115	1119	1127	rVV	600504	1097952	87.14%	5.932%
53	11.666	1127	1130	1133	rVV2	13118	28722	2.28%	0.155%
54	11.715	1133	1135	1143	rVB3	15639	43683	3.47%	0.236%
55	11.863	1143	1150	1154	rBV	407286	901325	71.54%	4.869%
56	11.932	1154	1157	1164	rVB	140315	303108	24.06%	1.638%
57	12.050	1164	1169	1171	rBV5	8149	18204	1.44%	0.098%
58	12.238	1182	1188	1189	rBV4	22777	54155	4.30%	0.293%
59	12.277	1189	1192	1196	rVB	131045	214665	17.04%	1.160%
60	12.445	1206	1209	1211	rBV	41208	66509	5.28%	0.359%
61	12.474	1211	1212	1216	rVV	22509	39407	3.13%	0.213%
62	12.563	1217	1221	1225	rVV3	47563	99158	7.87%	0.536%
63	12.661	1228	1231	1234	rVV	72136	115780	9.19%	0.626%
64	12.730	1234	1238	1242	rVV6	15582	40849	3.24%	0.221%
65	12.839	1242	1249	1253	rVV	448883	783591	62.19%	4.233%
66	12.928	1256	1258	1262	rVV2	24172	45388	3.60%	0.245%
67	13.006	1262	1266	1269	rVV4	14951	31463	2.50%	0.170%
68	13.066	1269	1272	1275	rVV	60998	90920	7.22%	0.491%
69	13.144	1275	1280	1283	rVV	308674	540867	42.93%	2.922%
70	13.184	1283	1284	1287	rVV	143605	194131	15.41%	1.049%
71	13.243	1287	1290	1297	rVB2	204960	457578	36.32%	2.472%

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72	13.411	1303	1307	1316	rVB	156473	257023	20.40%	1.389%
73	13.519	1316	1318	1320	rBV2	15482	20164	1.60%	0.109%
74	13.578	1320	1324	1330	rVV	581102	878320	69.71%	4.745%
75	13.716	1335	1338	1342	rBV3	23189	48780	3.87%	0.264%
76	13.805	1342	1347	1350	rBV	59159	103453	8.21%	0.559%
77	13.864	1350	1353	1354	rVV	25941	39421	3.13%	0.213%
78	13.903	1354	1357	1359	rVV	578100	929496	73.77%	5.022%
79	13.943	1359	1361	1366	rVB	226412	334927	26.58%	1.809%
80	14.022	1366	1369	1374	rVB3	14834	30648	2.43%	0.166%
81	14.130	1374	1380	1383	rBV3	65882	170678	13.55%	0.922%
82	14.179	1383	1385	1390	rVV3	96139	169340	13.44%	0.915%
83	14.288	1392	1396	1399	rVV2	29549	57246	4.54%	0.309%
84	14.337	1399	1401	1405	rVB	28184	38856	3.08%	0.210%
85	14.406	1405	1408	1409	rBV	52681	73529	5.84%	0.397%
86	14.436	1409	1411	1414	rVB	67726	97255	7.72%	0.525%
87	14.495	1414	1417	1420	rVB	81784	116786	9.27%	0.631%
88	14.603	1424	1428	1432	rVB	115561	200959	15.95%	1.086%
89	14.741	1439	1442	1444	rBV	31891	52126	4.14%	0.282%
90	14.820	1447	1450	1452	rBV2	20345	27933	2.22%	0.151%
91	14.860	1452	1454	1457	rVB	63477	86300	6.85%	0.466%
92	14.919	1457	1460	1461	rBV	15941	25110	1.99%	0.136%
93	15.096	1475	1478	1481	rVV	42501	69853	5.54%	0.377%
94	15.155	1481	1484	1486	rVB3	15408	25050	1.99%	0.135%
95	15.214	1486	1490	1494	rBV2	108692	234734	18.63%	1.268%
96	15.382	1504	1507	1510	rVB	15041	25431	2.02%	0.137%
97	15.500	1512	1519	1521	rBV2	21391	51223	4.07%	0.277%
98	15.619	1527	1531	1537	rVB	24754	62218	4.94%	0.336%
99	15.816	1547	1551	1556	rVB	18093	35883	2.85%	0.194%
100	16.930	1660	1664	1672	rVB3	7889	25071	1.99%	0.135%

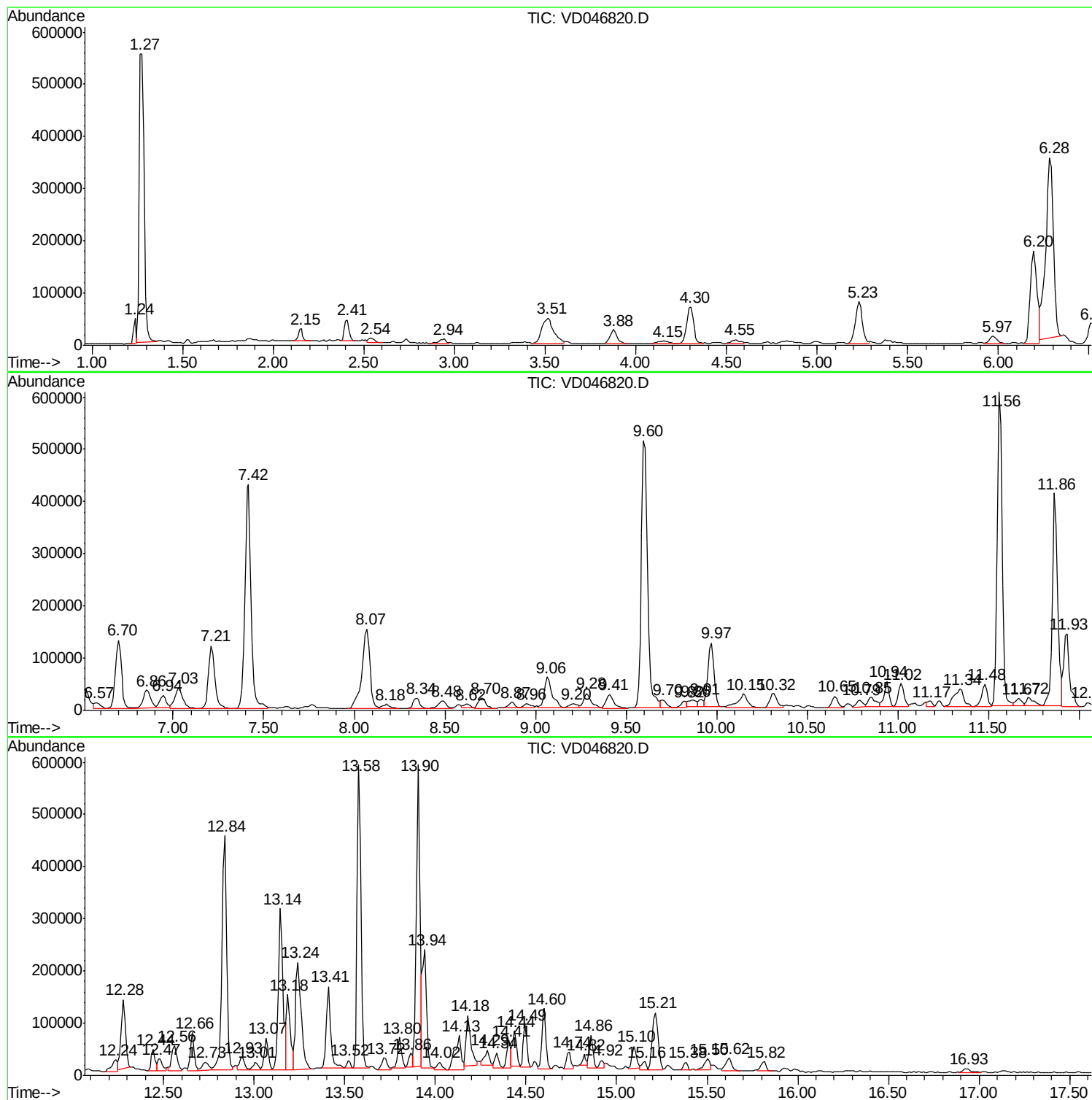
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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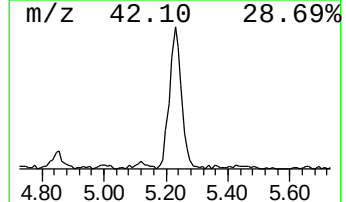
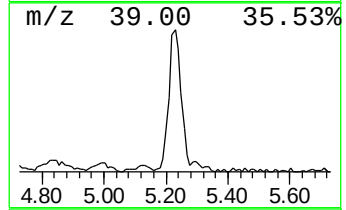
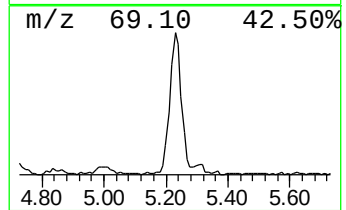
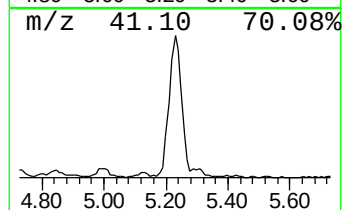
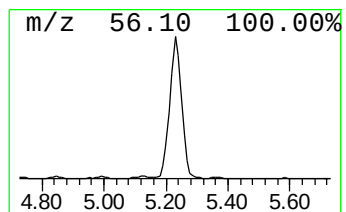
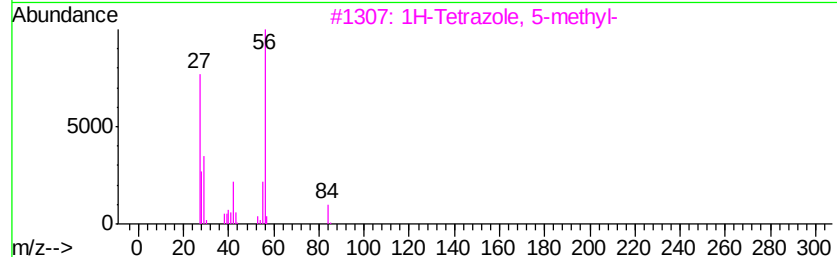
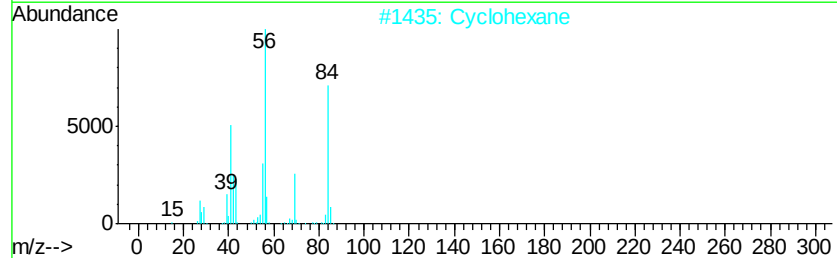
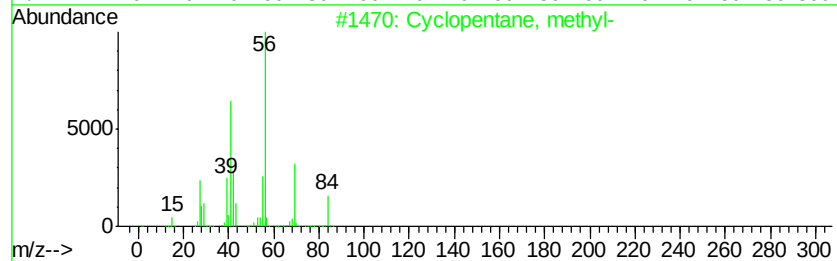
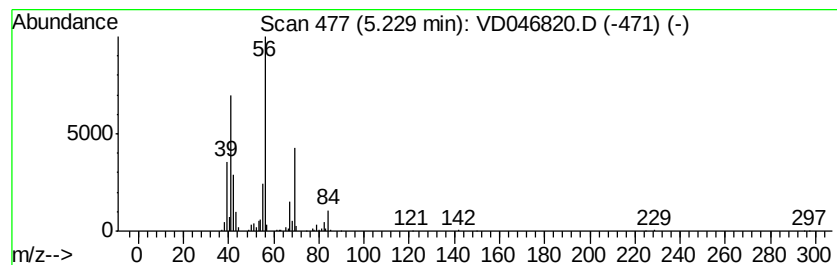
Quant Method : W:\HPCHEM1\MSVOA_D\METHOD\82D091815S.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclopentane, methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.23	10.48 ug/l	220788	Pentafluorobenzene	6.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	83
2		Cyclohexane	84	C6H12	000110-82-7	72
3		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	52
4		Cyclobutane, 1,2-diethyl-, cis-	112	C8H16	061141-50-2	50
5		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	50



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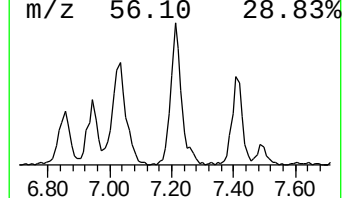
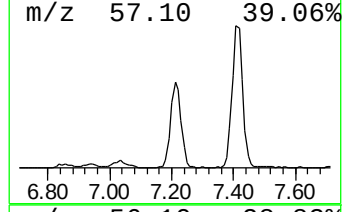
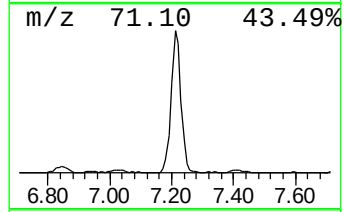
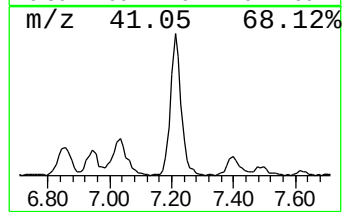
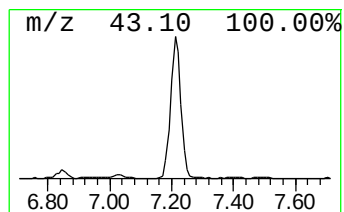
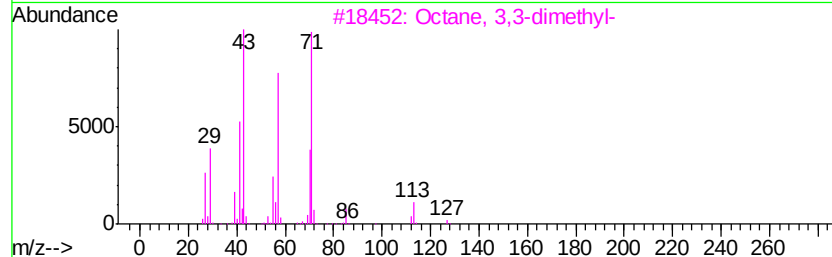
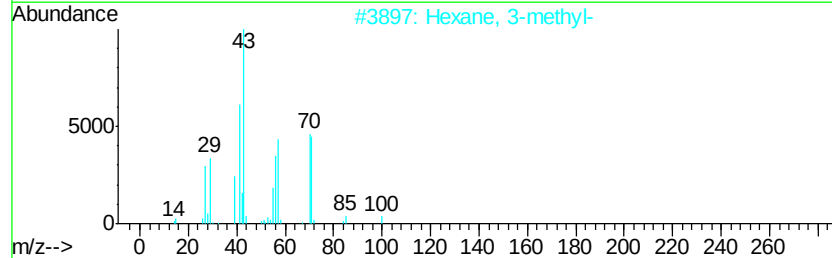
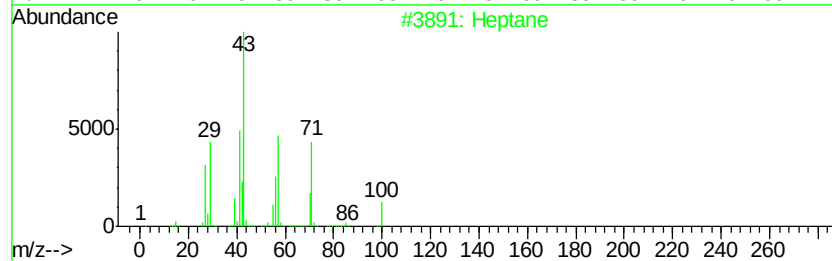
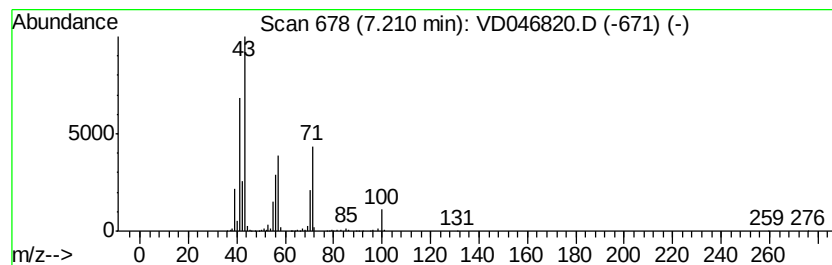
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Heptane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.21	13.83 ug/l	294804	1,4-Difluorobenzene	7.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane	100	C7H16	000142-82-5	91
2		Hexane, 3-methyl-	100	C7H16	000589-34-4	42
3		Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	37
4		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	28
5		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	28



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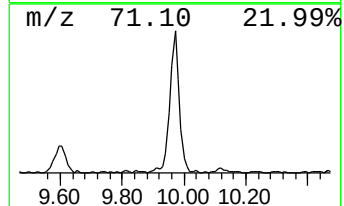
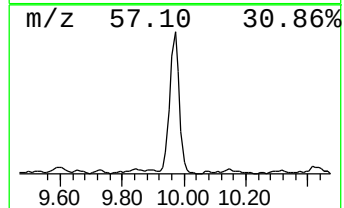
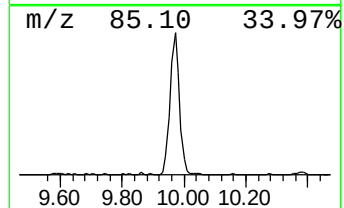
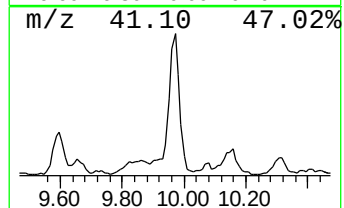
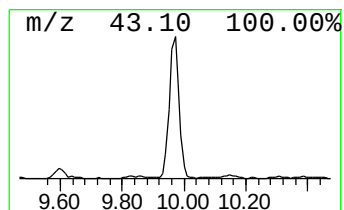
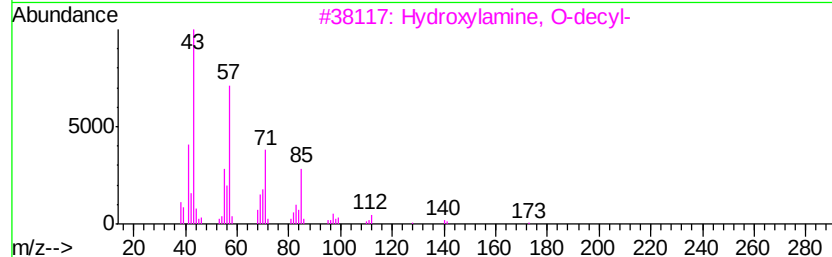
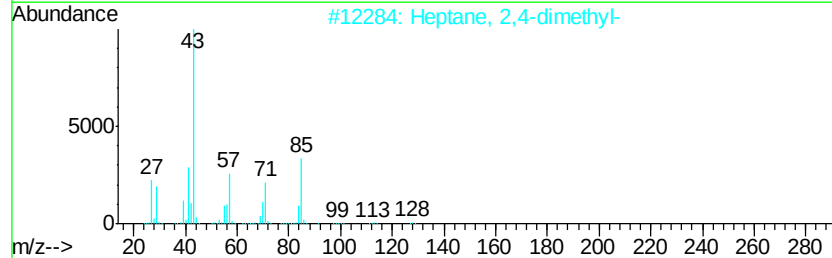
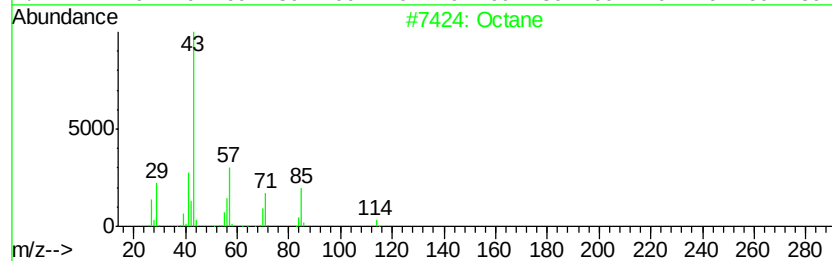
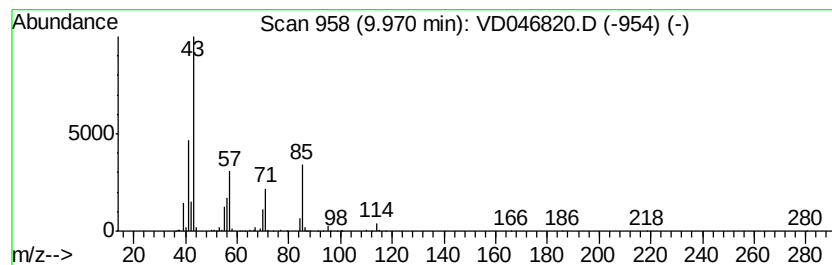
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TIC Integration Parameters: LSCINT.P

Peak Number 4 Octane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.97	12.15 ug/l	266694	Chlorobenzene-d5	11.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane		114	C8H18	000111-65-9	91
2	Heptane, 2,4-dimethyl-		128	C9H20	002213-23-2	72
3	Hydroxylamine, O-decyl-		173	C10H23NO	029812-79-1	64
4	Ether, hexyl pentyl		172	C11H24O	032357-83-8	56
5	Undecane, 2,4-dimethyl-		184	C13H28	017312-80-0	53



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Acq On : 21 Sep 2015 21:08
Operator : FY/SY
Sample : G3727-03
Misc : 6.18gm/5mL/MSVOA_D/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
MMB6B

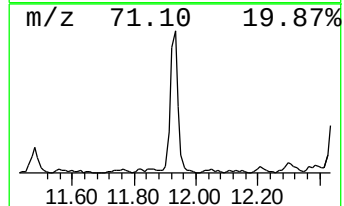
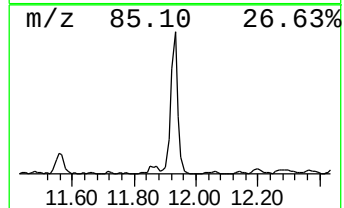
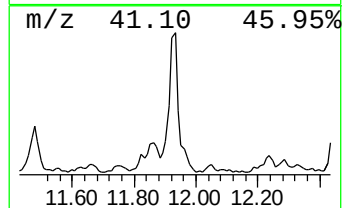
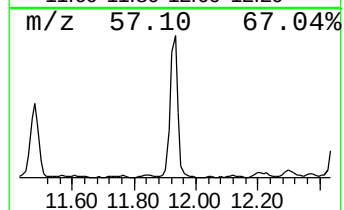
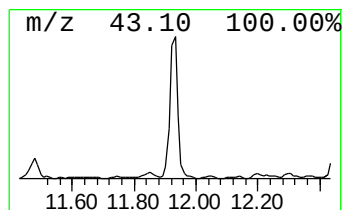
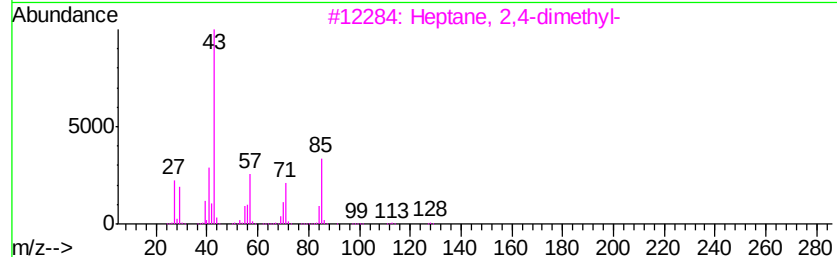
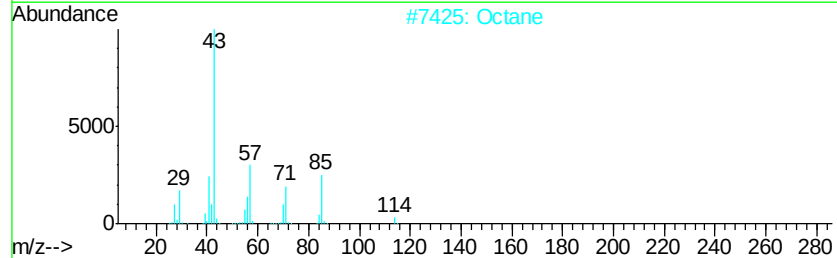
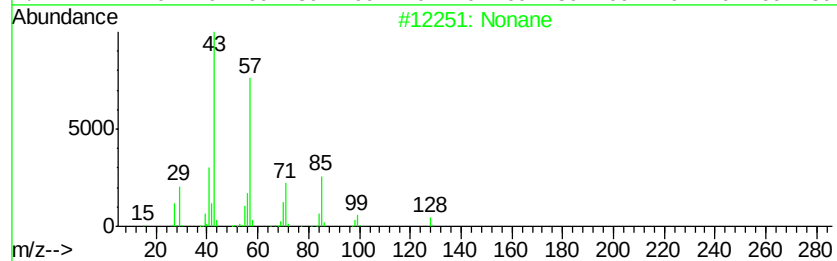
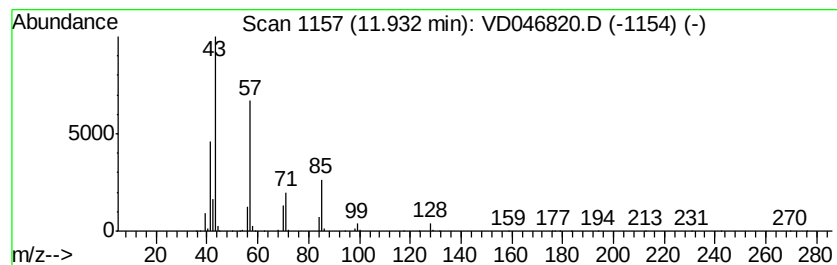
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Quant Title : SW846 8260

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

Peak Number 5 Nonane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	13.80 ug/l	303108	Chlorobenzene-d5	11.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane		128	C9H20	000111-84-2	91
2	Octane		114	C8H18	000111-65-9	72
3	Heptane, 2,4-dimethyl-		128	C9H20	002213-23-2	72
4	Undecane, 2,4-dimethyl-		184	C13H28	017312-80-0	59
5	Hexane, 2,4-dimethyl-		114	C8H18	000589-43-5	53



Data Path : W:\HPCHEM1\MSVOA_D\Data\VD092115\
Data File : VD046820.D
Acq On : 21 Sep 2015 21:08
Operator : FY/SY
Sample : G3727-03
Misc : 6.18gm/5mL/MSVOA_D/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
MMB6B

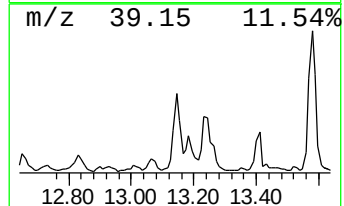
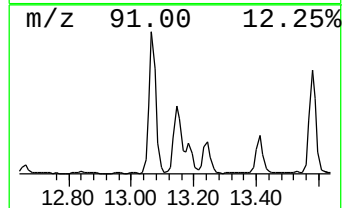
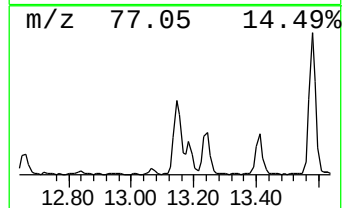
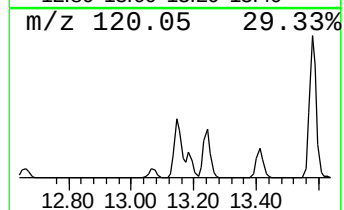
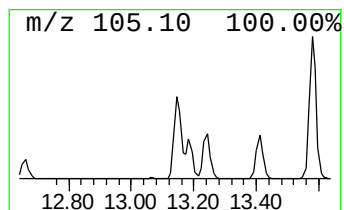
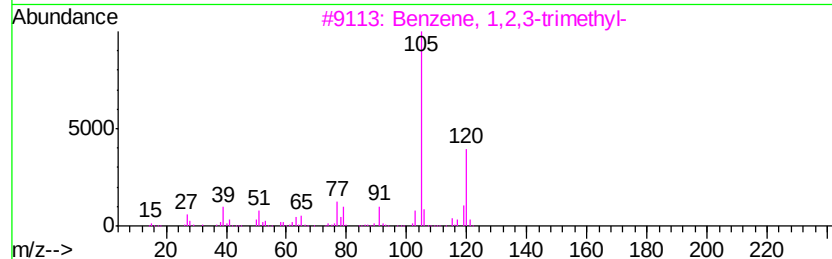
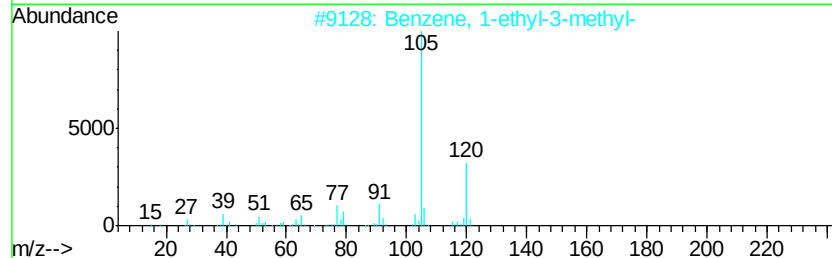
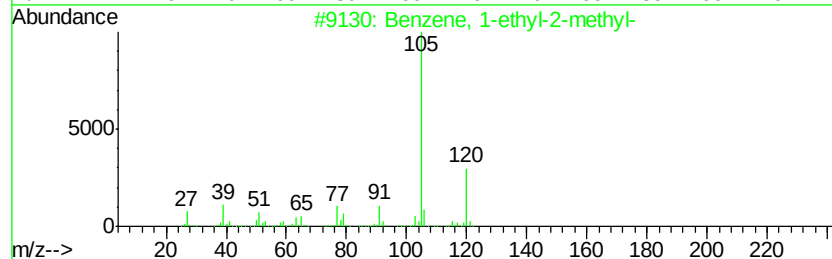
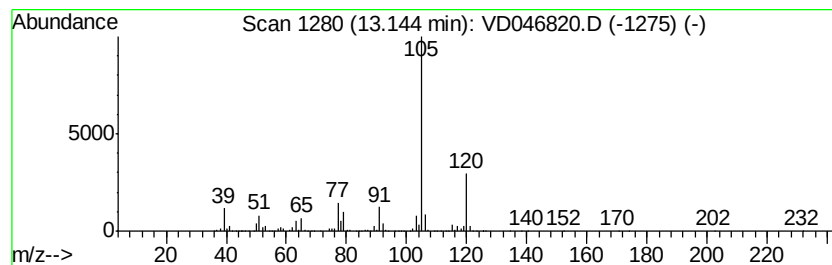
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.14	29.09 ug/l	540867	1,4-Dichlorobenzene-d4	13.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : W:\HPCHEM1\MSVOA_D\Data\VD092115\
Data File : VD046820.D
Acq On : 21 Sep 2015 21:08
Operator : FY/SY
Sample : G3727-03
Misc : 6.18gm/5mL/MSVOA_D/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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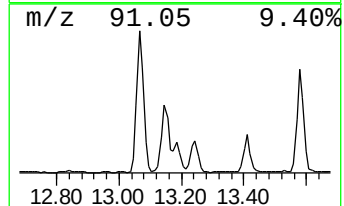
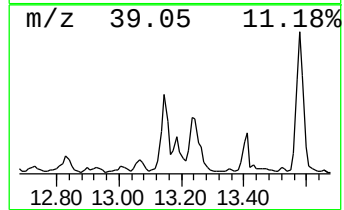
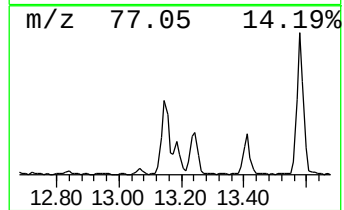
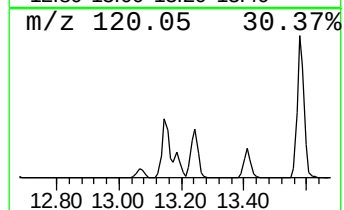
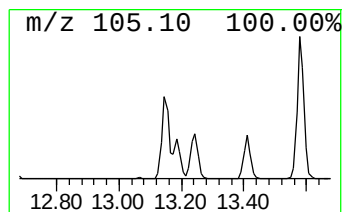
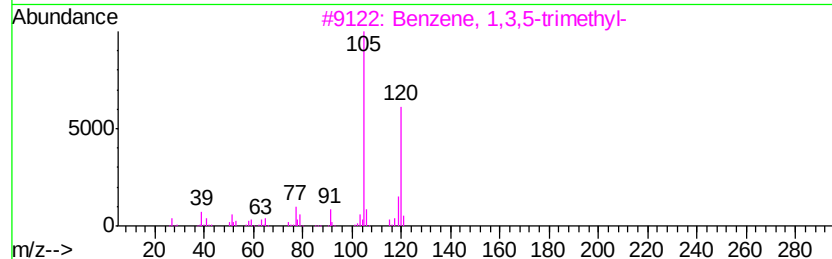
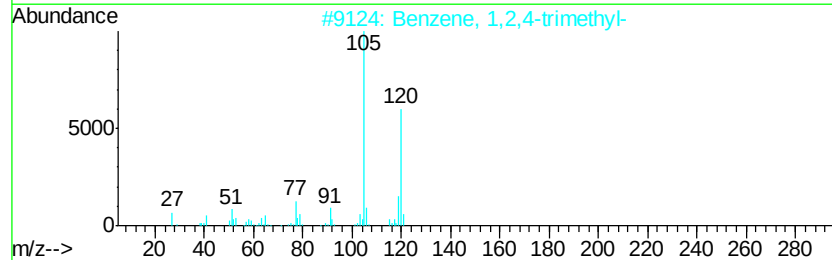
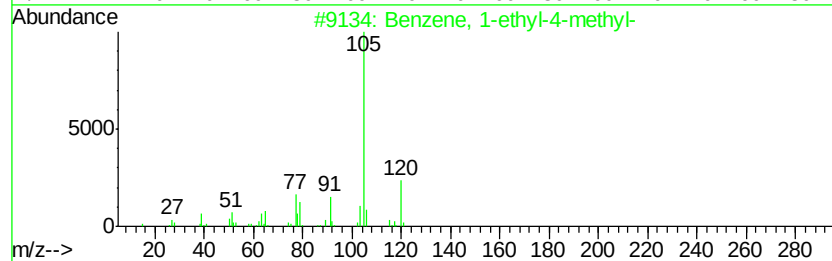
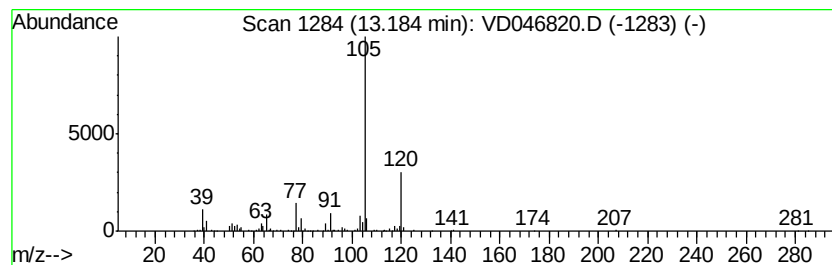
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-4-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.18	10.44 ug/l	194131	1,4-Dichlorobenzene-d4	13.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
3		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	87
4		Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	86
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	78



Data Path : W:\HPCHEM1\MSVOA_D\Data\VD092115\
Data File : VD046820.D
Acq On : 21 Sep 2015 21:08
Operator : FY/SY
Sample : G3727-03
Misc : 6.18gm/5mL/MSVOA_D/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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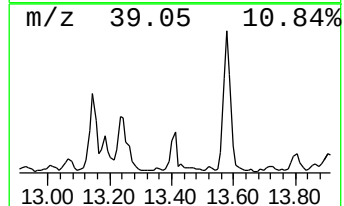
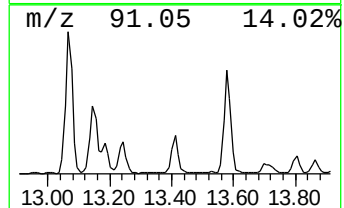
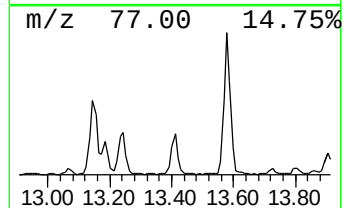
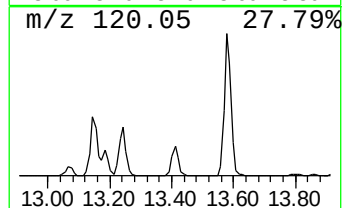
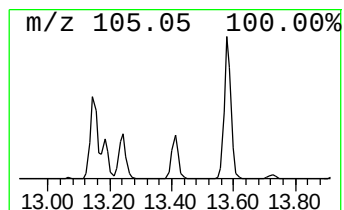
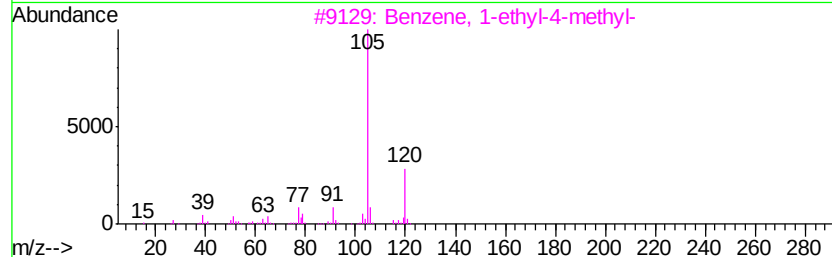
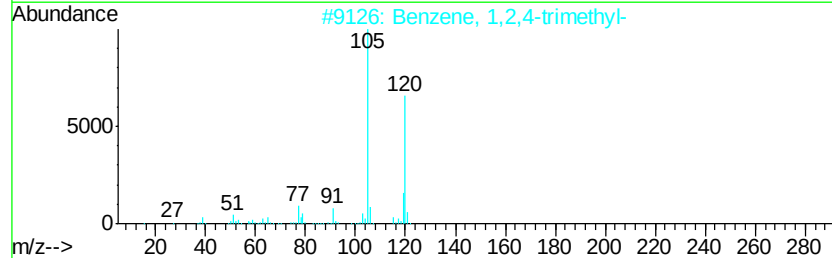
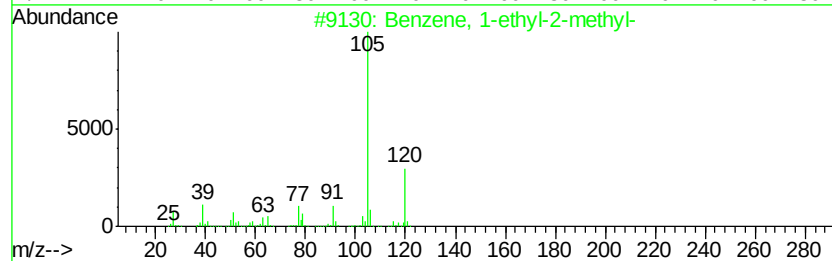
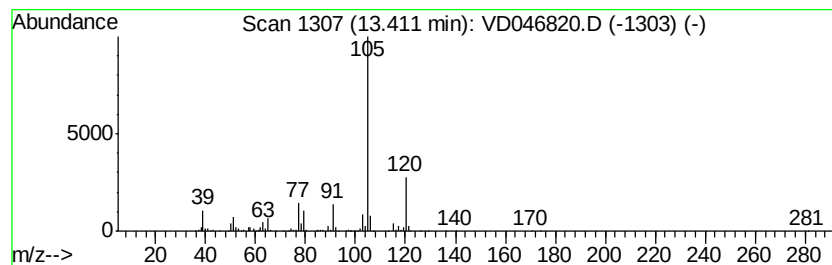
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,4-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.41	13.83 ug/l	257023	1,4-Dichlorobenzene-d4	13.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91



Data Path : W:\HPCHEM1\MSVOA_D\Data\VD092115\
Data File : VD046820.D
Acq On : 21 Sep 2015 21:08
Operator : FY/SY
Sample : G3727-03
Misc : 6.18gm/5mL/MSVOA_D/SOIL
ALS Vial : 24 Sample Multiplier: 1

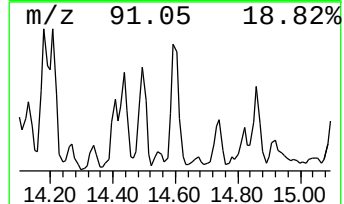
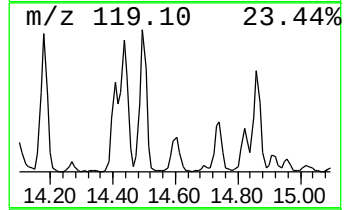
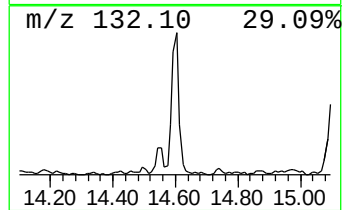
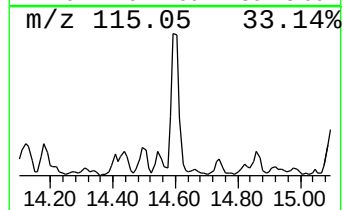
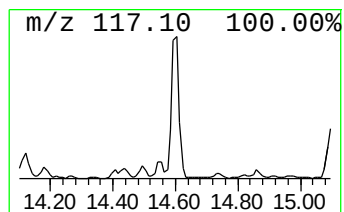
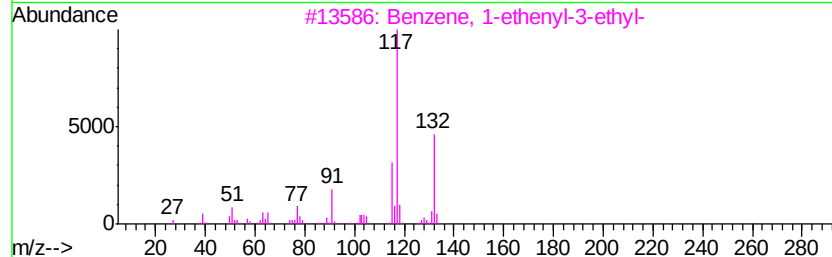
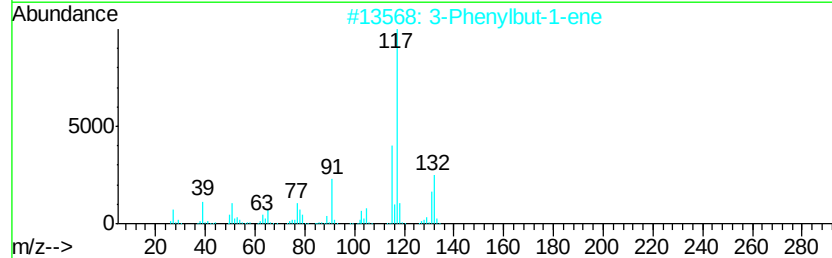
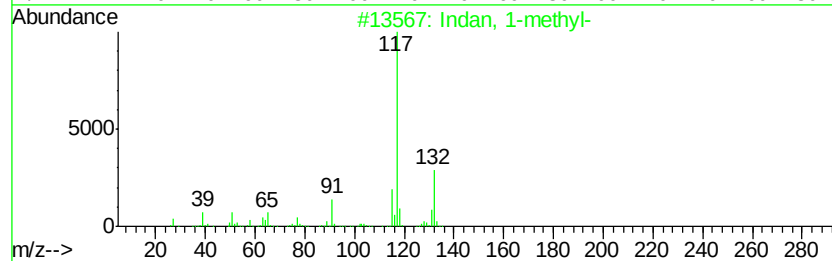
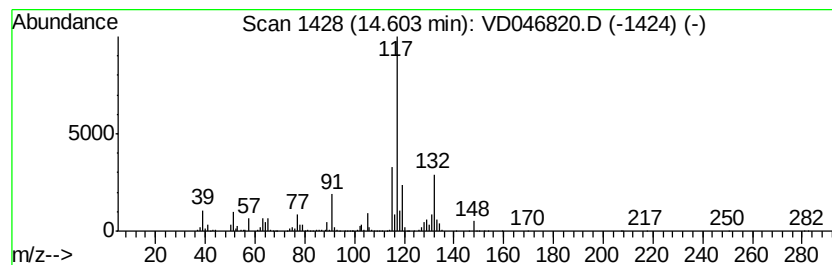
Instrument :
MSVOA_D
ClientSampleId :
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Quant Method : W:\HPCHEM1\MSVOA_D\METHOD\82D091815S.M
Quant Title : SW846 8260

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

Peak Number 9 Indan, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	10.81 ug/l	200959	1,4-Dichlorobenzene-d4	13.90
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indan, 1-methyl-	132 C10H12	000767-58-8	81
2	3-Phenylbut-1-ene	132 C10H12	000934-10-1	76
3	Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4	74
4	Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0	74
5	Benzene, (2-methyl-2-propenyl)-	132 C10H12	003290-53-7	72



Data Path : W:\HPCHEM1\MSVOA_D\Data\VD092115\
Data File : VD046820.D
Acq On : 21 Sep 2015 21:08
Operator : FY/SY
Sample : G3727-03
Misc : 6.18gm/5mL/MSVOA_D/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
MMB6B

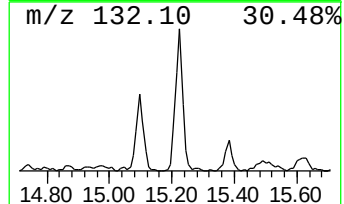
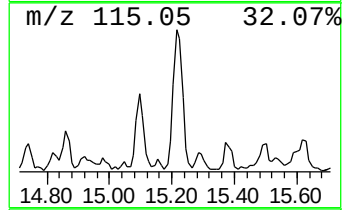
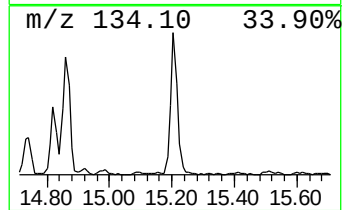
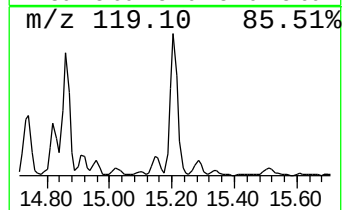
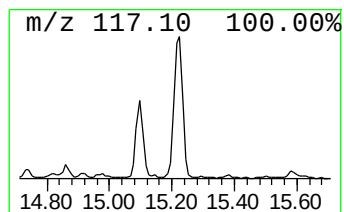
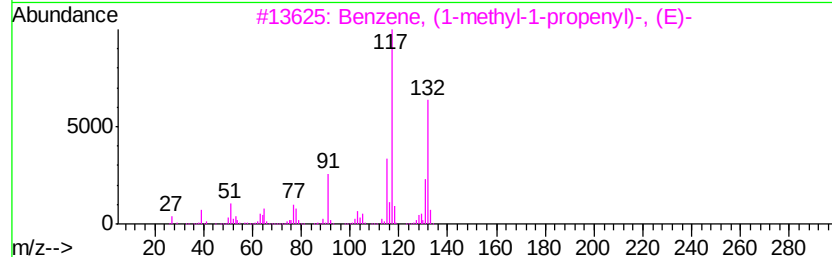
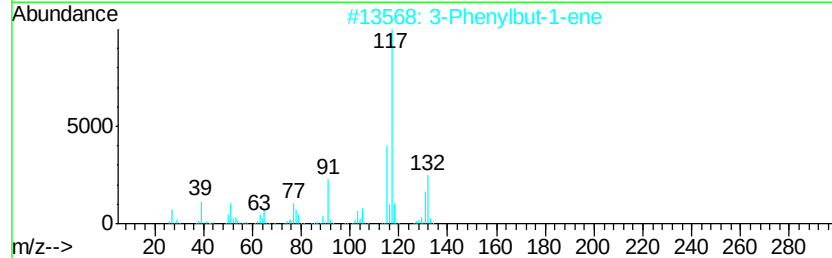
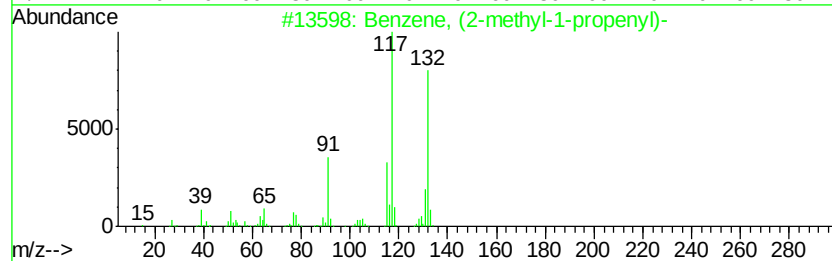
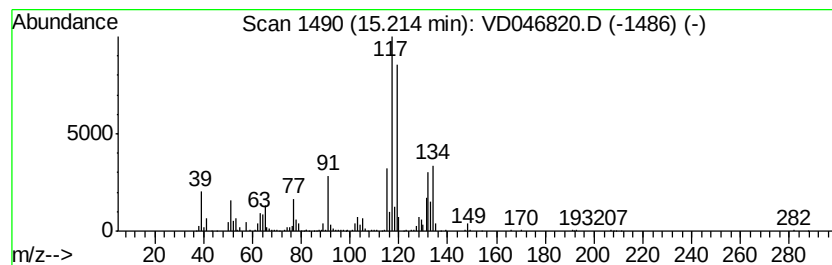
Quant Method : W:\HPCHEM1\MSVOA_D\METHOD\82D091815S.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, (2-methyl-1-propen... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.21	12.63 ug/l	234734	1,4-Dichlorobenzene-d4	13.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	55
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	55
3		Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	49
4		1-Phenyl-1-butene	132	C10H12	000824-90-8	49
5		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	49



Data Path : W:\HPCHEM1\MSVOA_D\Data\VD092115\
Data File : VD046820.D
Acq On : 21 Sep 2015 21:08
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Sample : G3727-03
Misc : 6.18gm/5mL/MSVOA_D/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
MMB6B

Quant Method : W:\HPCHEM1\MSVOA_D\METHOD\82D091815S.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
Cyclopentane, met...	5.23	10.5	ug/l	220788	1	6.29	1052930	50.0
Heptane	7.21	13.8	ug/l	294804	2	7.42	1065790	50.0
Octane	9.97	12.2	ug/l	266694	3	11.56	1097950	50.0
Nonane	11.93	13.8	ug/l	303108	3	11.56	1097950	50.0
Benzene, 1-ethyl-...	13.14	29.1	ug/l	540867	4	13.90	929496	50.0
Benzene, 1-ethyl-...	13.18	10.4	ug/l	194131	4	13.90	929496	50.0
Benzene, 1,2,4-tr...	13.41	13.8	ug/l	257023	4	13.90	929496	50.0
Indan, 1-methyl-	14.60	10.8	ug/l	200959	4	13.90	929496	50.0
Benzene, (2-methy...	15.21	12.6	ug/l	234734	4	13.90	929496	50.0