

Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
 Data File : BG018203.D
 Acq On : 1 Aug 2015 4:38
 Operator : TP/UM
 Sample : G3065-12
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C02A6

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % 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:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG073115.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.139	413	417	425	rVB2	12148	23702	0.36%	0.079%
2	5.509	472	480	485	rBV4	12438	23239	0.36%	0.078%
3	5.767	516	524	530	rBV8	7144	17001	0.26%	0.057%
4	6.155	587	590	596	rVB5	13640	23160	0.35%	0.078%
5	6.296	608	614	621	rVB3	12303	22177	0.34%	0.074%
6	6.373	621	627	634	rBV8	9819	23389	0.36%	0.078%
7	6.478	634	645	650	rBV6	14712	38973	0.60%	0.131%
8	6.619	665	669	673	rBV4	19841	37838	0.58%	0.127%
9	6.672	673	678	683	rVB2	61247	94498	1.45%	0.317%
10	6.825	696	704	708	rBV	59659	103773	1.59%	0.348%
11	6.966	720	728	732	rBV2	50323	83656	1.28%	0.281%
12	7.013	732	736	742	rVB	70791	113375	1.74%	0.380%
13	7.113	748	753	760	rBV3	38658	74642	1.14%	0.250%
14	7.195	762	767	774	rVB7	8214	20024	0.31%	0.067%
15	7.289	779	783	790	rBV8	9425	21983	0.34%	0.074%
16	7.383	790	799	800	rBV5	11494	18606	0.28%	0.062%
17	7.442	805	809	811	rVV	60619	97257	1.49%	0.326%
18	7.477	811	815	822	rVB	214099	371179	5.68%	1.245%
19	7.554	822	828	831	rBV7	9428	18834	0.29%	0.063%
20	7.595	831	835	838	rBV2	13166	19004	0.29%	0.064%
21	7.671	844	848	851	rVB3	14330	17490	0.27%	0.059%
22	7.753	856	862	871	rVB6	21040	50651	0.78%	0.170%
23	7.918	886	890	896	rBV6	14033	25003	0.38%	0.084%
24	8.024	903	908	913	rVB5	34543	66098	1.01%	0.222%
25	8.118	920	924	927	rBV3	22489	37548	0.57%	0.126%
26	8.200	934	938	944	rBV2	49604	75016	1.15%	0.252%
27	8.294	950	954	959	rVB2	41026	63767	0.98%	0.214%
28	8.482	981	986	991	rVB2	53340	86036	1.32%	0.288%
29	8.582	991	1003	1007	rVV4	44829	121832	1.86%	0.409%
30	8.635	1007	1012	1017	rVV	80575	159195	2.44%	0.534%
31	8.711	1021	1025	1034	rVB	84602	147852	2.26%	0.496%
32	8.823	1034	1044	1051	rBV8	48200	135821	2.08%	0.455%
33	8.911	1056	1059	1063	rBV5	10657	18924	0.29%	0.063%
34	9.122	1089	1095	1099	rBV4	32276	61297	0.94%	0.206%

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35	9.181	1099	1105	1113	rVB5	40354	87789	1.34%	0.294%
36	9.351	1126	1134	1139	rBV4	89899	175573	2.69%	0.589%
37	9.440	1145	1149	1159	rVB7	38631	96689	1.48%	0.324%
38	9.551	1162	1168	1173	rBV7	14474	24865	0.38%	0.083%
39	9.628	1175	1181	1184	rBV6	20905	32302	0.49%	0.108%
40	9.669	1184	1188	1192	rBV5	35163	62003	0.95%	0.208%
41	9.810	1208	1212	1217	rBV5	11902	30390	0.47%	0.102%
42	9.892	1222	1226	1232	rVB2	99866	157284	2.41%	0.527%
43	9.968	1234	1239	1242	rBV4	18505	27254	0.42%	0.091%
44	10.051	1250	1253	1258	rVB7	12978	18759	0.29%	0.063%
45	10.133	1262	1267	1271	rBV4	30945	52896	0.81%	0.177%
46	10.256	1280	1288	1294	rBV	332386	614510	9.41%	2.061%
47	10.315	1294	1298	1305	rVV	66489	114406	1.75%	0.384%
48	10.380	1305	1309	1314	rVV5	28693	46662	0.71%	0.156%
49	10.438	1315	1319	1324	rVB3	46627	72019	1.10%	0.241%
50	10.497	1325	1329	1336	rVB9	12851	23454	0.36%	0.079%
51	10.756	1368	1373	1376	rBV5	14470	23212	0.36%	0.078%
52	10.803	1376	1381	1384	rVV2	19796	28784	0.44%	0.097%
53	10.944	1401	1405	1412	rVB7	39001	82340	1.26%	0.276%
54	11.232	1452	1454	1459	rVB4	15657	21271	0.33%	0.071%
55	11.302	1461	1466	1470	rVV3	75061	134204	2.05%	0.450%
56	11.349	1470	1474	1483	rVB8	53374	128044	1.96%	0.429%
57	11.555	1503	1509	1514	rBV7	31485	61667	0.94%	0.207%
58	11.713	1532	1536	1541	rBV6	42341	95523	1.46%	0.320%
59	11.854	1558	1560	1562	rBV3	16697	16957	0.26%	0.057%
60	11.943	1571	1575	1582	rVB3	83397	155489	2.38%	0.521%
61	12.013	1582	1587	1592	rBV4	66393	125507	1.92%	0.421%
62	12.084	1595	1599	1602	rBV2	31343	47216	0.72%	0.158%
63	12.160	1608	1612	1616	rVB2	42564	66270	1.01%	0.222%
64	12.430	1654	1658	1663	rBV8	35596	58753	0.90%	0.197%
65	12.489	1664	1668	1672	rVV7	67370	109743	1.68%	0.368%
66	12.571	1676	1682	1688	rVB5	119533	220716	3.38%	0.740%
67	12.630	1689	1692	1698	rVV7	53670	102282	1.57%	0.343%
68	12.689	1699	1702	1705	rVV4	57223	73636	1.13%	0.247%
69	12.736	1705	1710	1716	rVV5	122673	219283	3.36%	0.735%
70	12.918	1736	1741	1746	rBV	257176	422613	6.47%	1.417%
71	13.076	1764	1768	1772	rBV6	63202	98508	1.51%	0.330%

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72	13.470	1831	1835	1839	rBV4	157879	245317	3.75%	0.823%
73	13.570	1847	1852	1856	rBV2	917712	1227963	18.79%	4.118%
74	13.617	1856	1860	1864	rBV	345862	526526	8.06%	1.766%
75	13.846	1896	1899	1902	rBV	109881	125664	1.92%	0.421%
76	13.887	1902	1906	1912	rVV4	311032	491254	7.52%	1.647%
77	13.946	1913	1916	1918	rVV3	158135	201212	3.08%	0.675%
78	13.981	1918	1922	1930	rVB	953026	1421420	21.76%	4.766%
79	14.058	1931	1935	1941	rBV5	177436	324286	4.96%	1.087%
80	14.116	1941	1945	1948	rBV2	431919	705196	10.79%	2.365%
81	14.158	1949	1952	1957	rVB2	539301	777032	11.89%	2.606%
82	14.451	2000	2002	2005	rBV	113049	111267	1.70%	0.373%
83	14.487	2005	2008	2012	rBV	258635	375992	5.75%	1.261%
84	14.704	2041	2045	2050	rBV4	381843	546262	8.36%	1.832%
85	14.792	2056	2060	2062	rBV3	277642	446677	6.84%	1.498%
86	14.880	2071	2075	2081	rVB5	331034	608415	9.31%	2.040%
87	14.951	2082	2087	2093	rBV2	1312028	2089684	31.98%	7.007%
88	15.209	2128	2131	2135	rBV2	358282	614382	9.40%	2.060%
89	15.903	2245	2249	2252	rBV	772065	1161479	17.78%	3.895%
90	16.972	2427	2431	2434	rVB	1312611	1726313	26.42%	5.789%
91	19.093	2789	2792	2797	rVB2	5074418	6533550	100.00%	21.908%
92	25.021	3797	3801	3811	rVB2	1477894	3320429	50.82%	11.134%

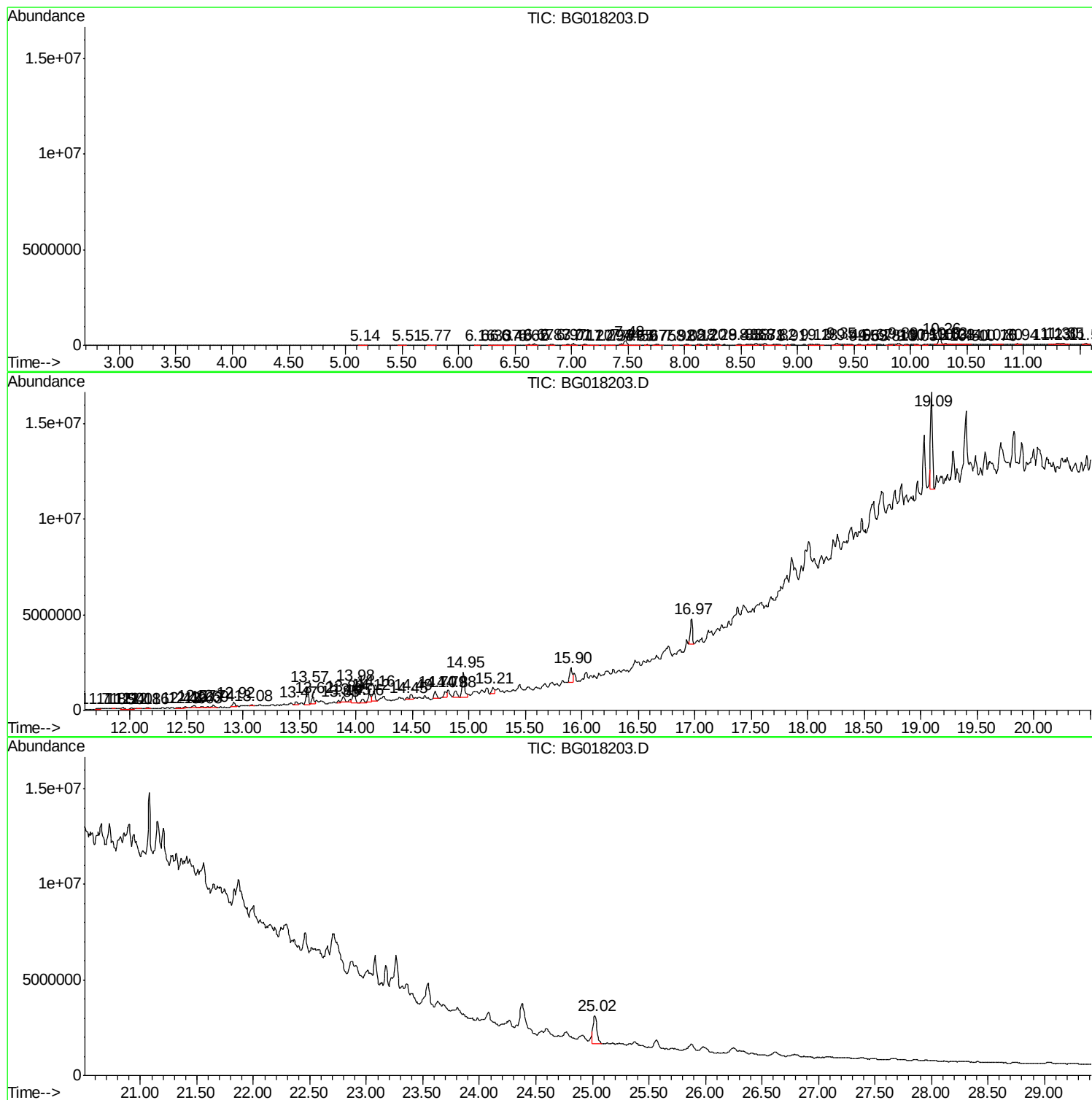
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG073115.M
Quant Title : SVOA CALIBRATION

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



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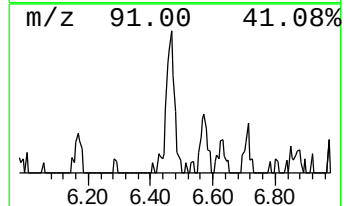
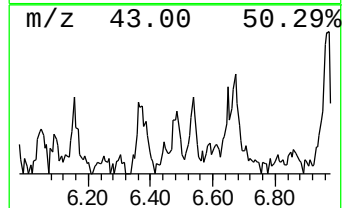
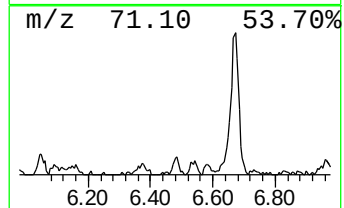
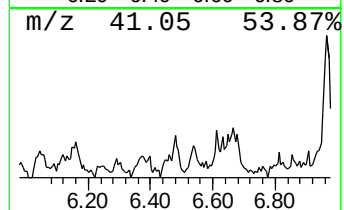
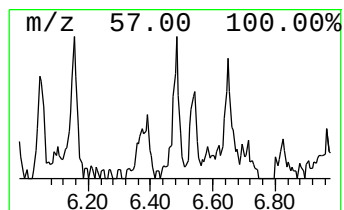
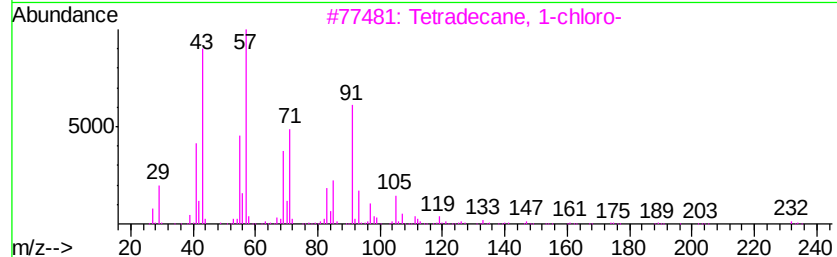
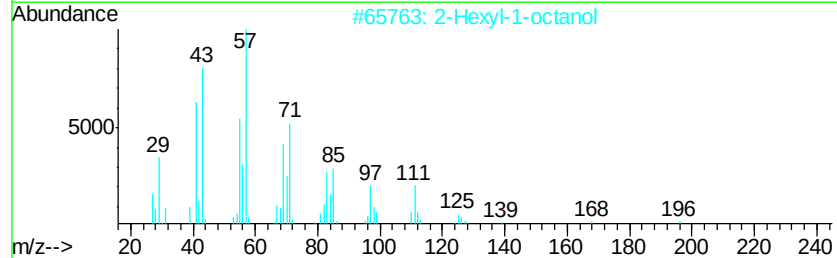
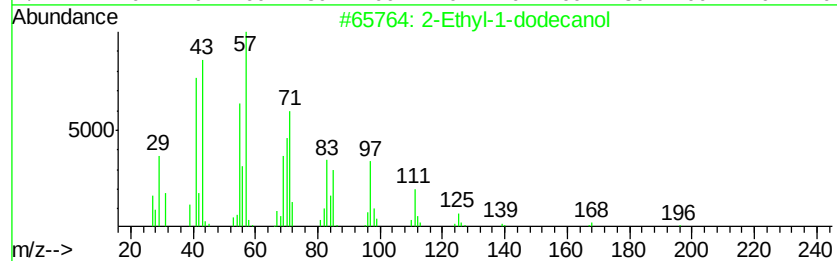
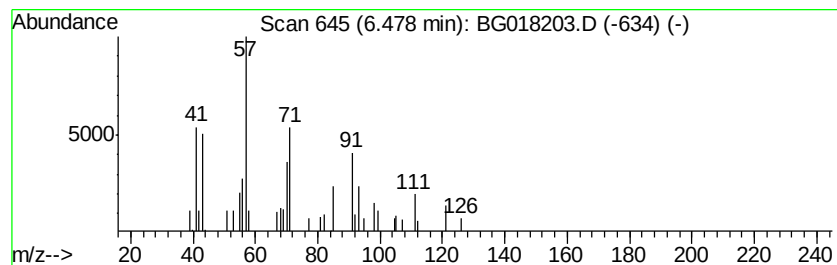
Instrument :
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG073115.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Ethyl-1-dodecanol Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.		
6.48	2.10 ng/ul	38973	1,4-Dichlorobenzene-d4	7.48		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Ethyl-1-dodecanol	214	C14H30O	019780-33-7	59	
2	2-Hexyl-1-octanol	214	C14H30O	019780-79-1	59	
3	Tetradecane, 1-chloro-	232	C14H29Cl	002425-54-9	50	
4	Tetradecane, 1-chloro-	232	C14H29Cl	002425-54-9	47	
5	Tridecane, 6-methyl-	198	C14H30	013287-21-3	43	



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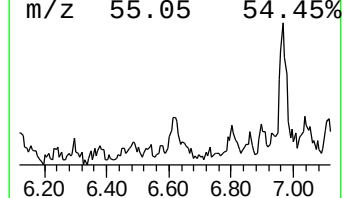
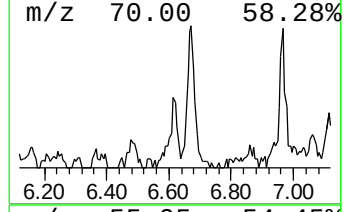
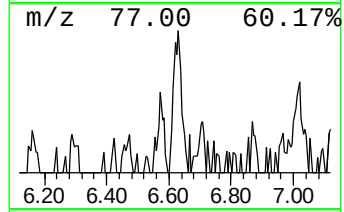
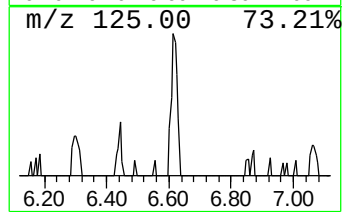
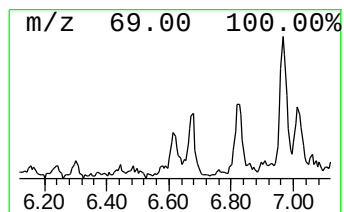
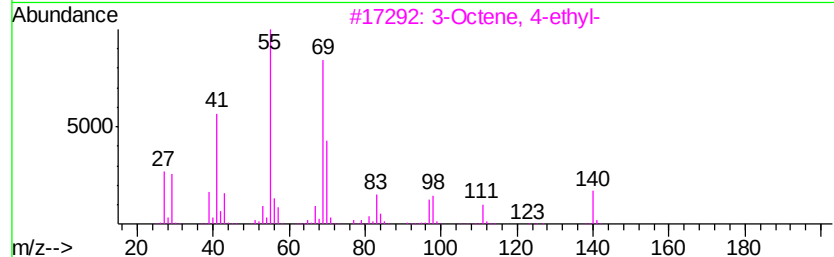
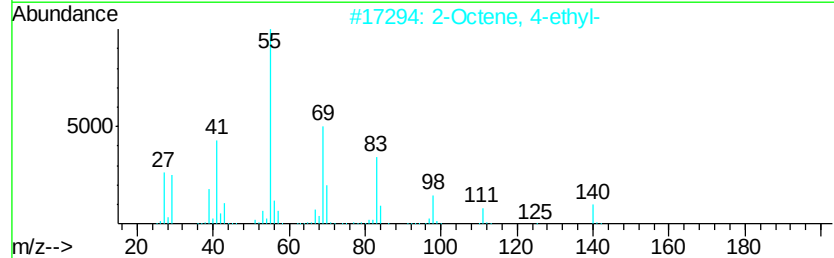
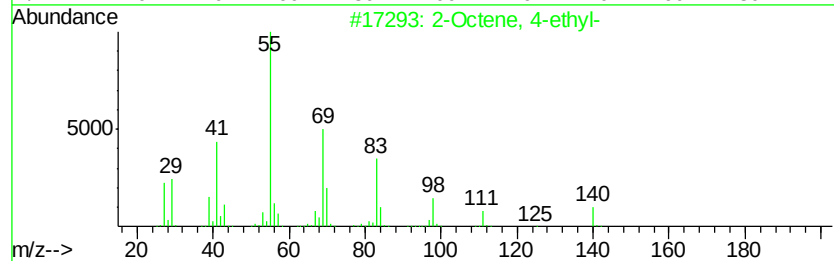
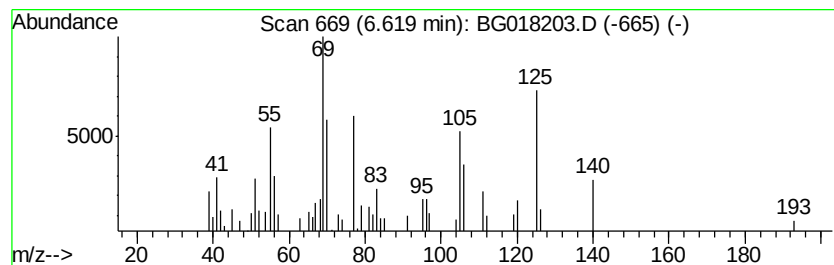
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG073115.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Cyclic6.62 Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	2.04 ng/ul	37838	1,4-Dichlorobenzene-d4	7.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Octene, 4-ethyl-	140	C10H20	053966-52-2	38
2		2-Octene, 4-ethyl-	140	C10H20	053966-52-2	38
3		3-Octene, 4-ethyl-	140	C10H20	053966-51-1	35
4		Cyclohexanone, 3,5-dimethyl-	126	C8H14O	002320-30-1	25
5		cis-3,5-Dimethylcyclohexanone	126	C8H14O	007214-49-5	22



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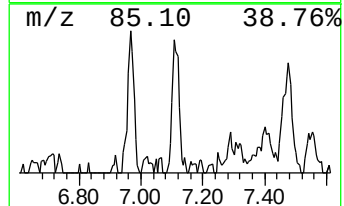
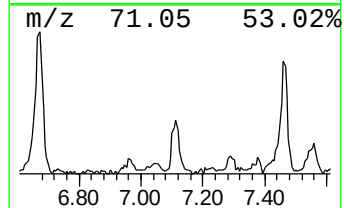
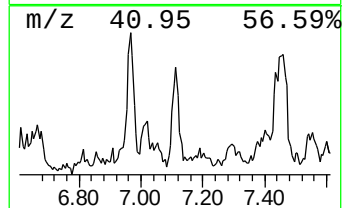
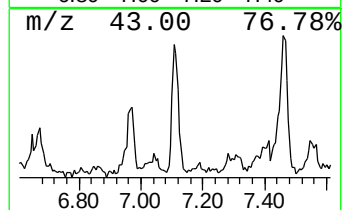
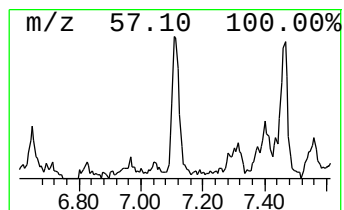
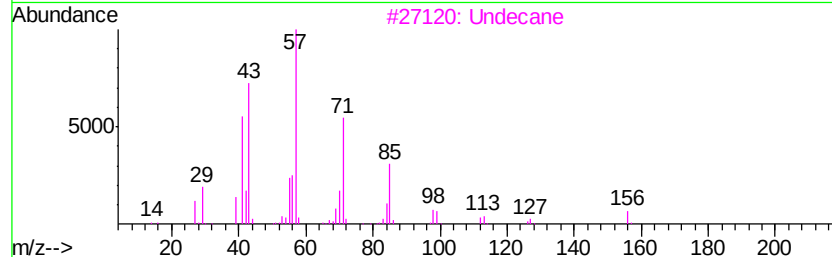
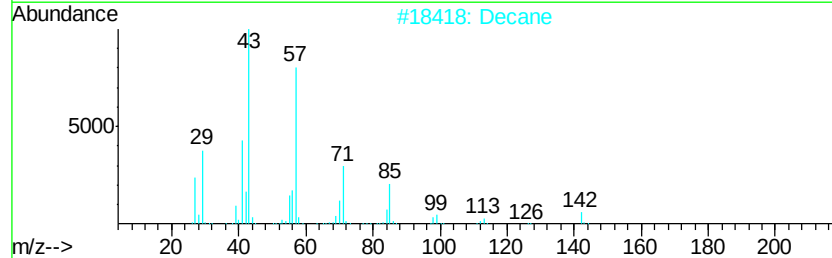
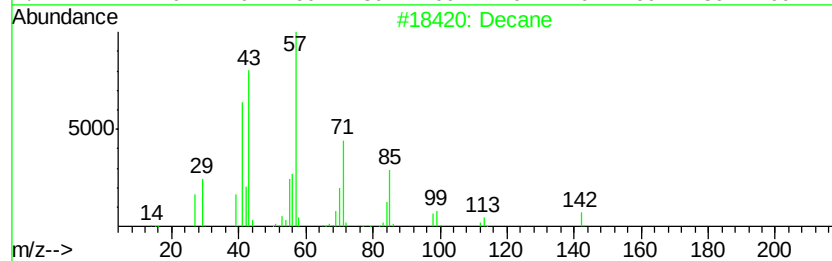
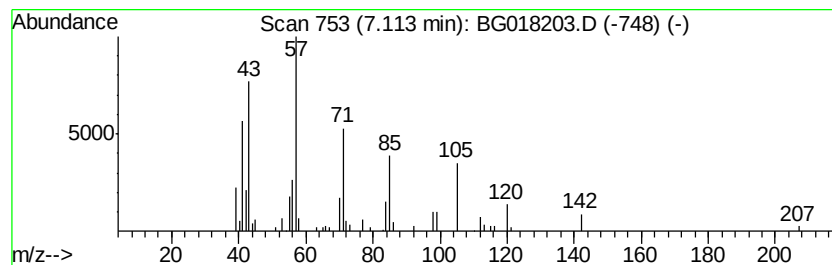
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.11	4.02 ng/ul	74642	1,4-Dichlorobenzene-d4	7.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decane	142	C10H22	000124-18-5	87
2		Decane	142	C10H22	000124-18-5	81
3		Undecane	156	C11H24	001120-21-4	64
4		Decane	142	C10H22	000124-18-5	64
5		Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	53



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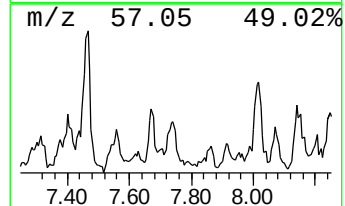
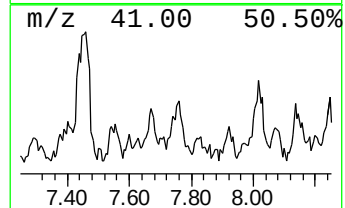
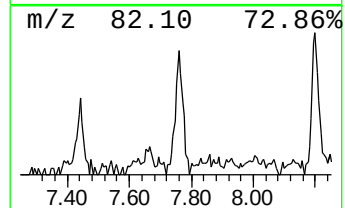
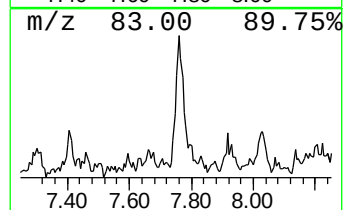
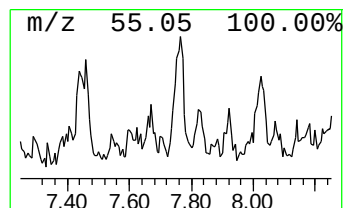
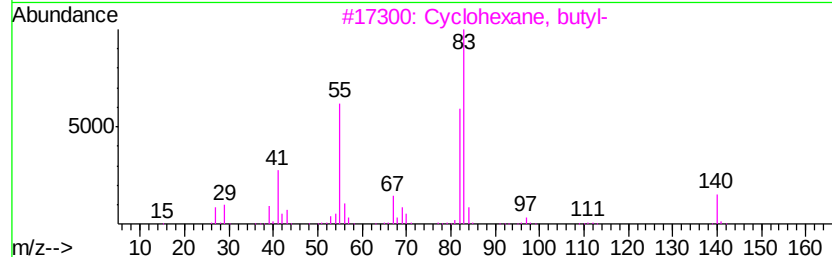
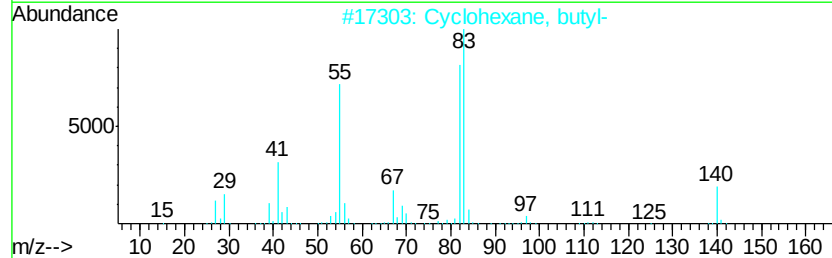
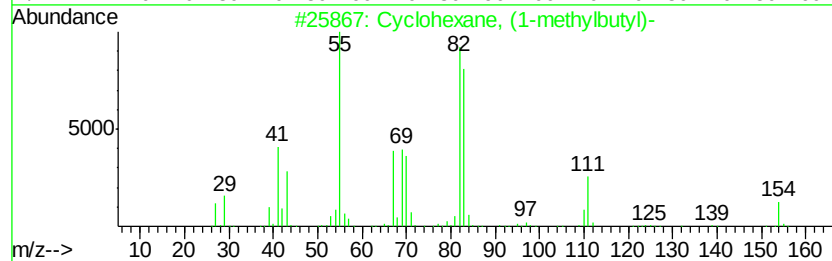
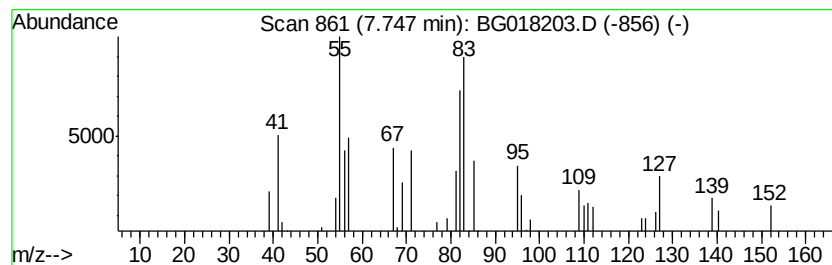
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Cyclic7.75 Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.75	2.73 ng/ul	50651	1,4-Dichlorobenzene-d4	7.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (1-methylbutyl)-	154	C11H22	061208-94-4	64
2			Cyclohexane, butyl-	140	C10H20	001678-93-9	52
3			Cyclohexane, butyl-	140	C10H20	001678-93-9	50
4			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	49
5			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	47



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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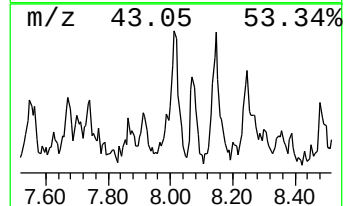
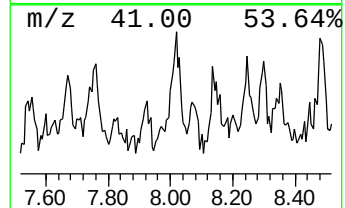
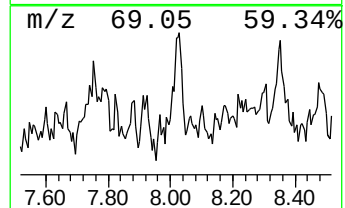
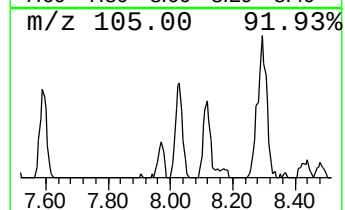
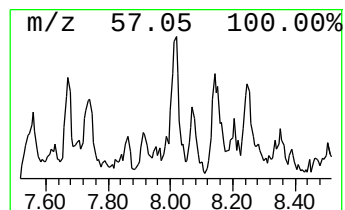
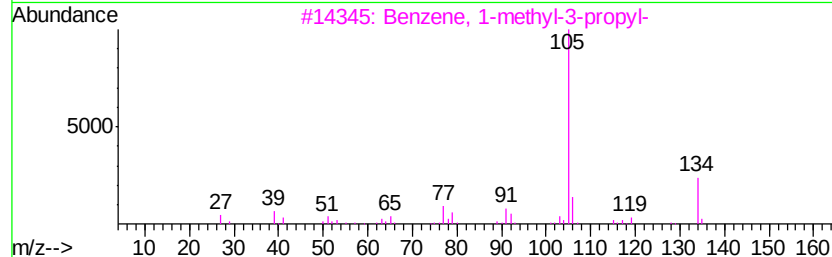
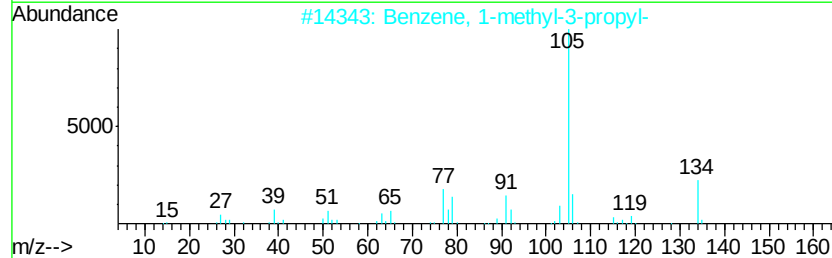
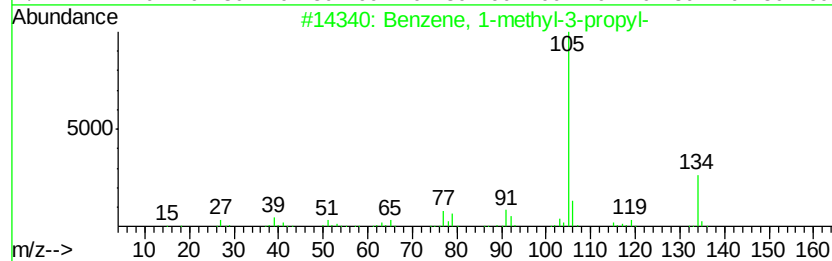
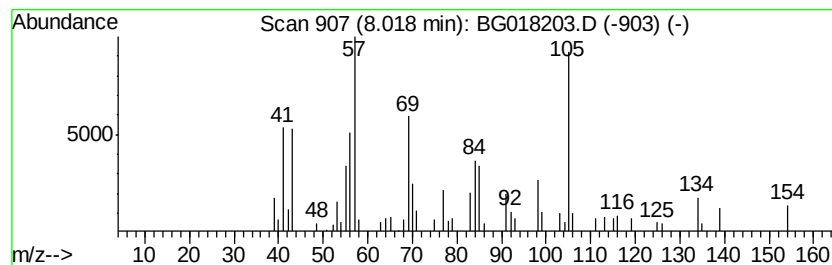
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-01 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.02	3.56 ng/ul	66098	1,4-Dichlorobenzene-d4	7.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	38
2		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	30
3		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	30
4		2-Tolyloxirane	134	C9H10O	002783-26-8	27
5		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	22



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
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ClientSampleId :
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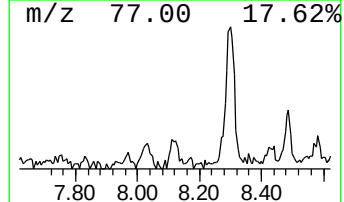
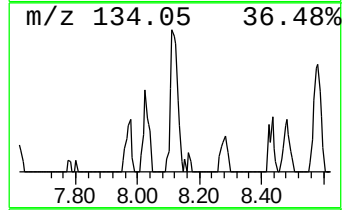
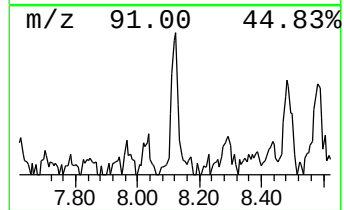
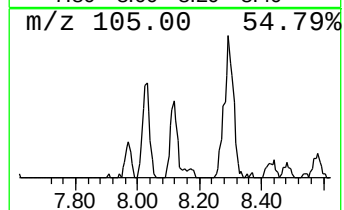
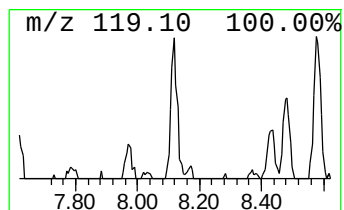
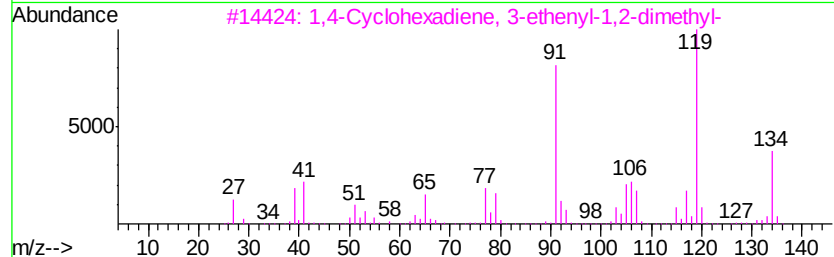
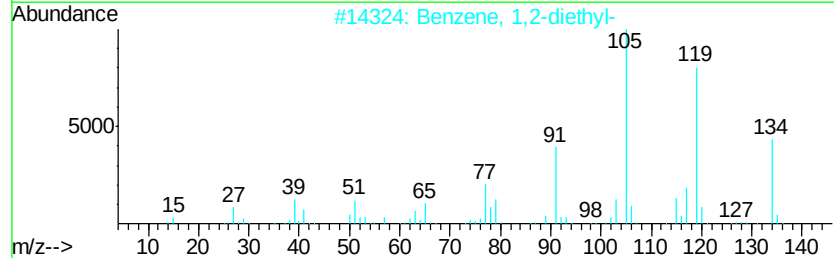
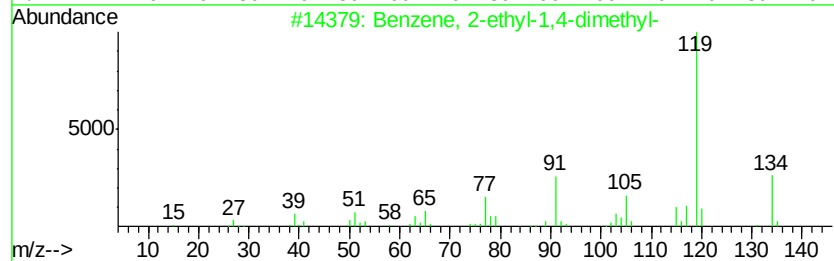
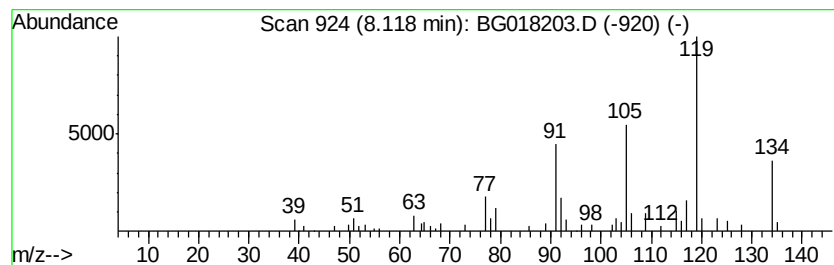
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.12	2.02 ng/ul	37548	1,4-Dichlorobenzene-d4	7.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	76
2		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	76
3		1,4-Cyclohexadiene, 3-ethenyl-1,...	134	C10H14	062338-57-2	72
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	68
5		Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	68



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
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Sample : G3065-12
Misc :
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ClientSampleId :
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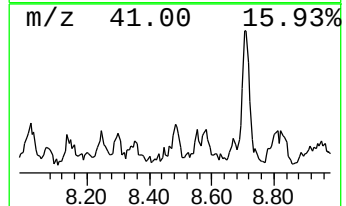
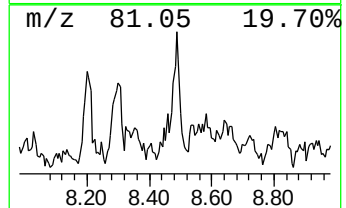
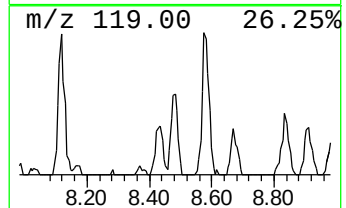
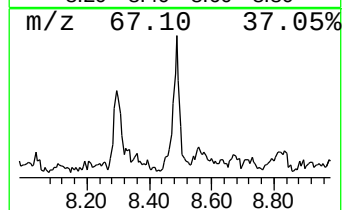
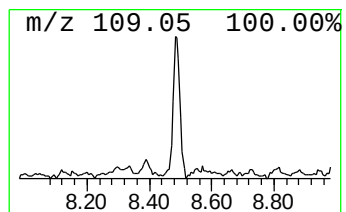
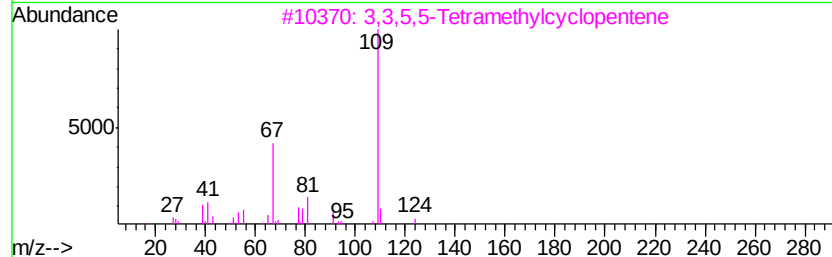
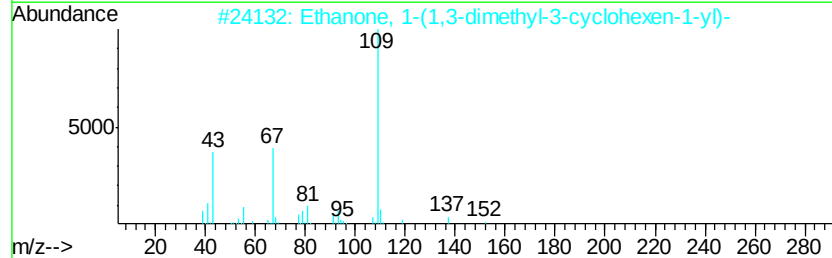
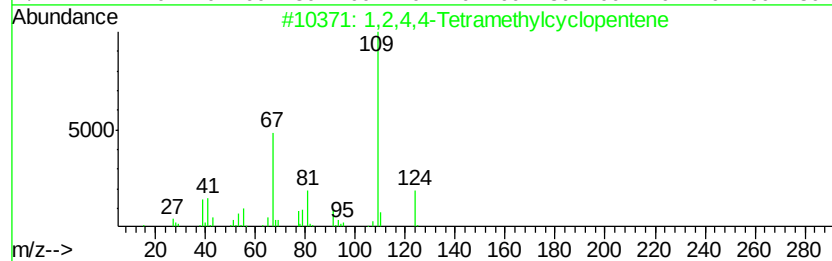
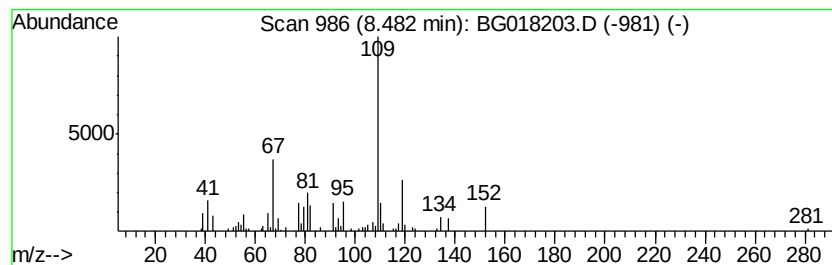
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1,2,4,4-Tetramethylcyclopentene... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.48	4.64 ng/ul	86036	1,4-Dichlorobenzene-d4	7.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2,4,4-Tetramethylcyclopentene	124	C9H16	065378-76-9	64
2			Ethanone, 1-(1,3-dimethyl-3-cycl...	152	C10H16O	051733-68-7	53
3			3,3,5,5-Tetramethylcyclopentene	124	C9H16	038667-10-6	52
4			Ethanone, 1-(1,4-dimethyl-3-cycl...	152	C10H16O	043219-68-7	50
5			1,3-Octadiene, 5,5-dimethyl-	138	C10H18	1000142-73-2	50



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
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Misc :
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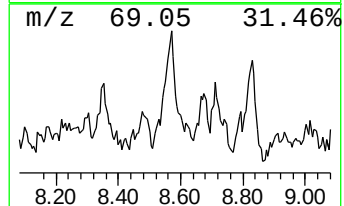
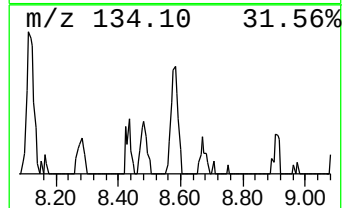
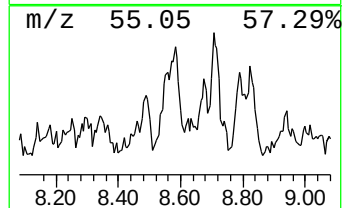
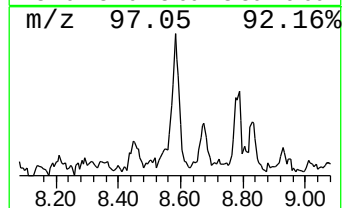
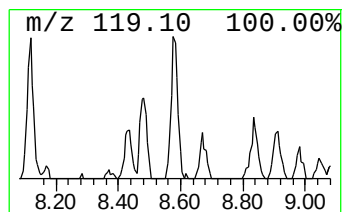
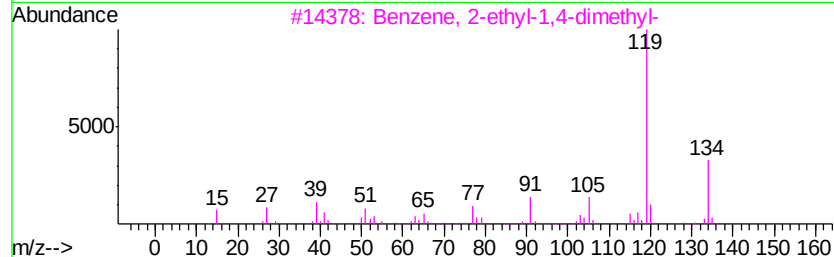
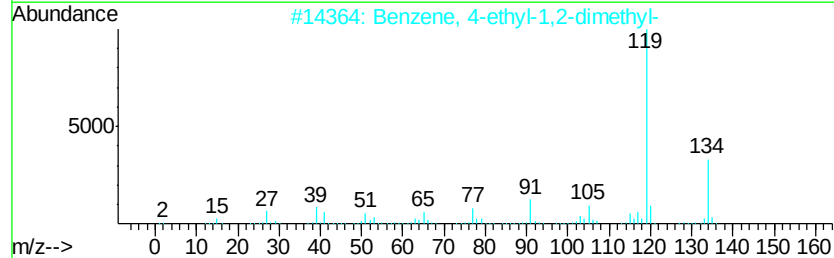
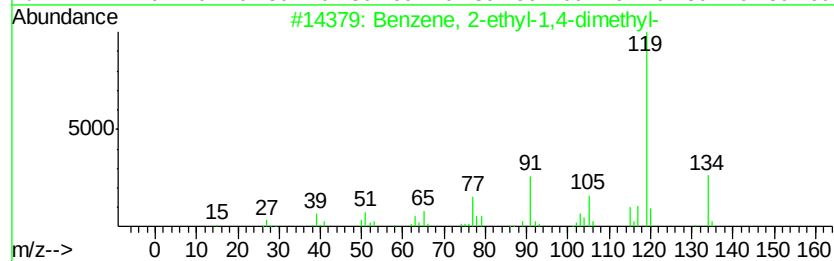
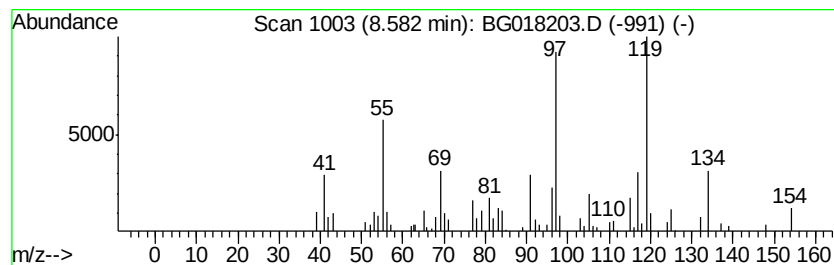
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-02 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.58	6.56 ng/ul	121832	1,4-Dichlorobenzene-d4	7.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	45
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	41
3		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	38
4		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	38
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	38



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
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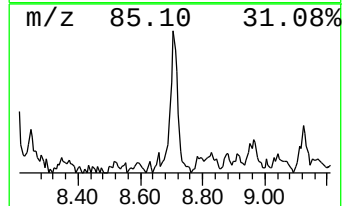
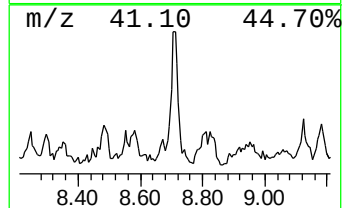
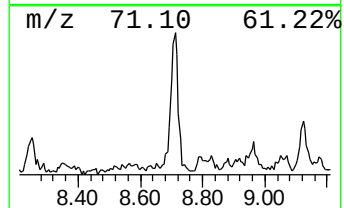
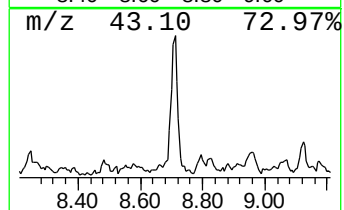
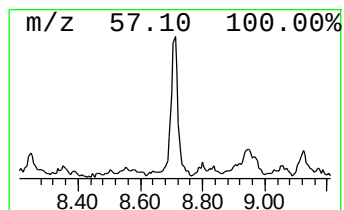
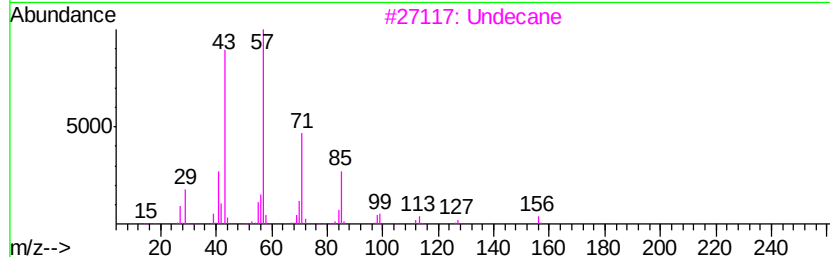
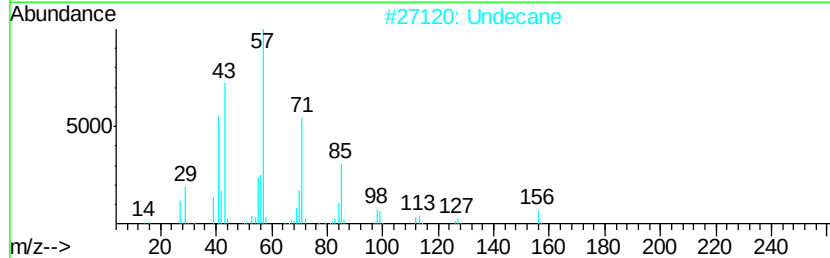
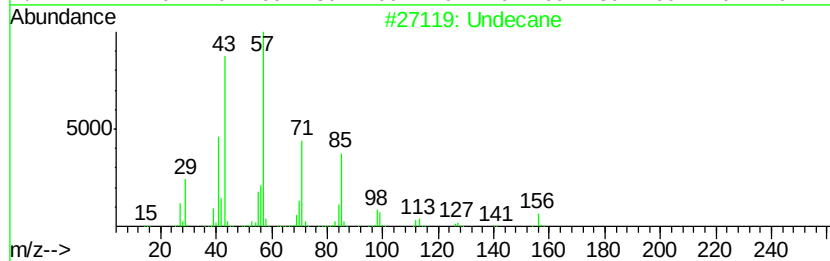
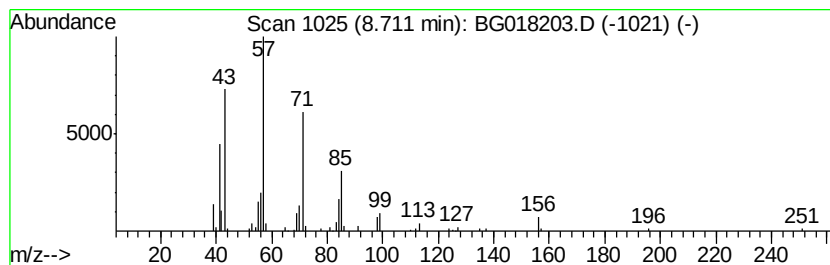
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.71	7.97 ng/ul	147852	1,4-Dichlorobenzene-d4	7.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Undecane	156	C11H24	001120-21-4	93
2		Undecane	156	C11H24	001120-21-4	91
3		Undecane	156	C11H24	001120-21-4	90
4		Undecane	156	C11H24	001120-21-4	90
5		Tetradecane	198	C14H30	000629-59-4	80



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
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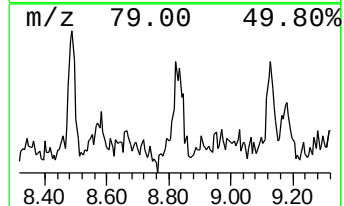
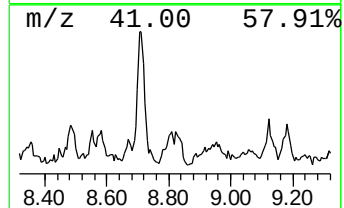
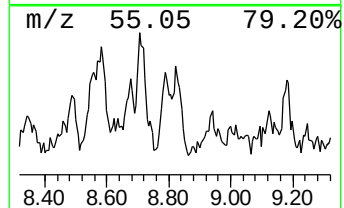
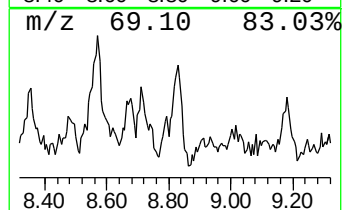
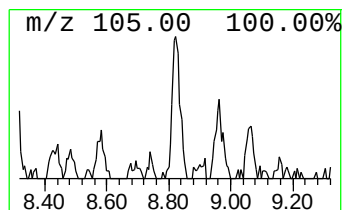
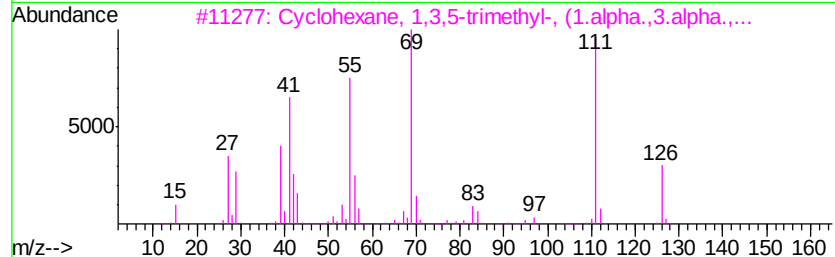
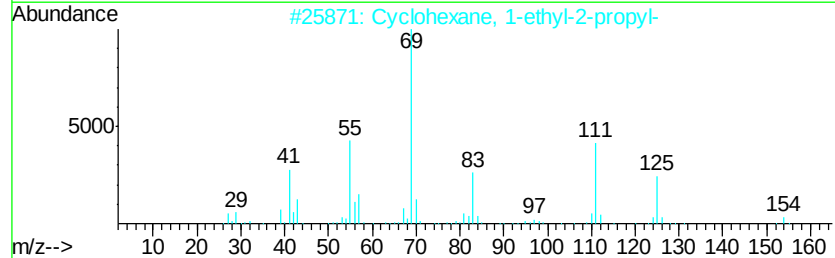
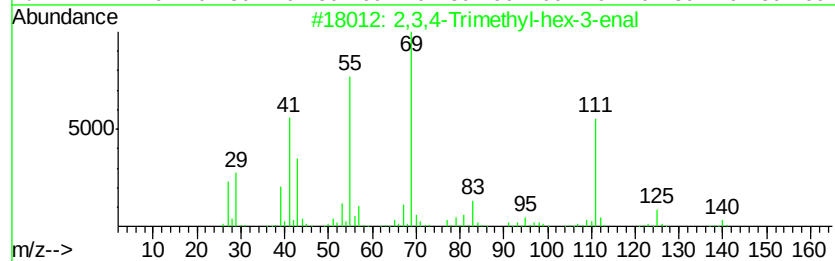
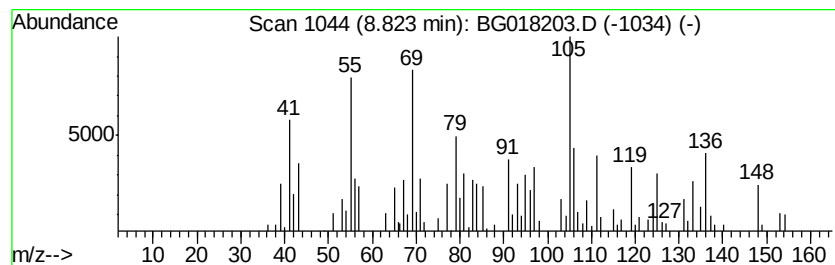
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 unknown-03 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.82	7.32 ng/ul	135821	1,4-Dichlorobenzene-d4	7.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,3,4-Trimethyl-hex-3-enal	140	C9H16O	1000193-72-9	27
2			Cyclohexane, 1-ethyl-2-propyl-	154	C11H22	062238-33-9	25
3			Cyclohexane, 1,3,5-trimethyl-, (...)	126	C9H18	001795-26-2	22
4			4-Decene, 4-methyl-, (E)-	154	C11H22	060366-66-7	22
5			Cyclopentaneethanol, .beta.,2,3-...	156	C10H20O	021721-18-6	22



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Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampled :
C02A6

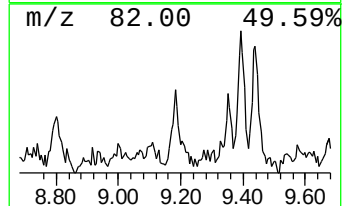
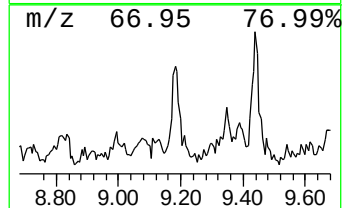
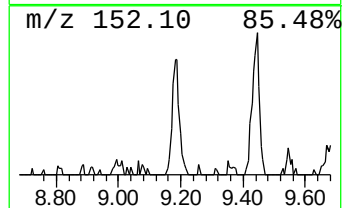
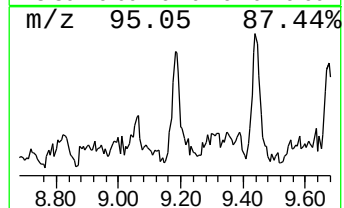
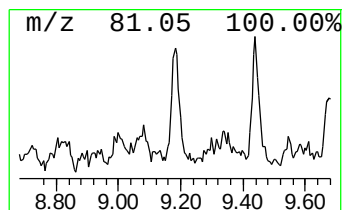
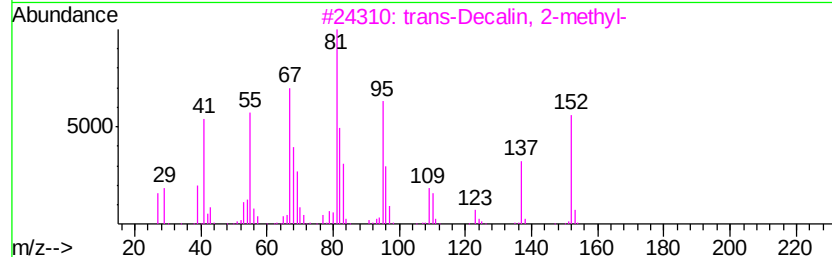
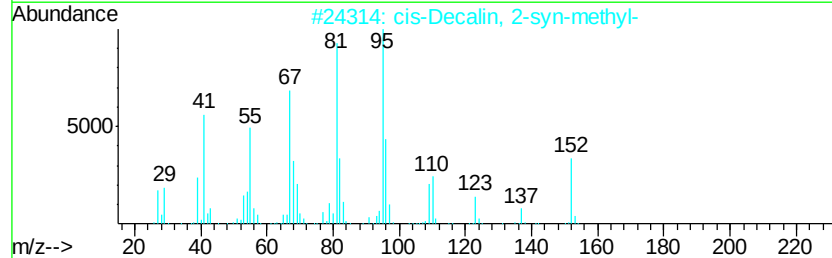
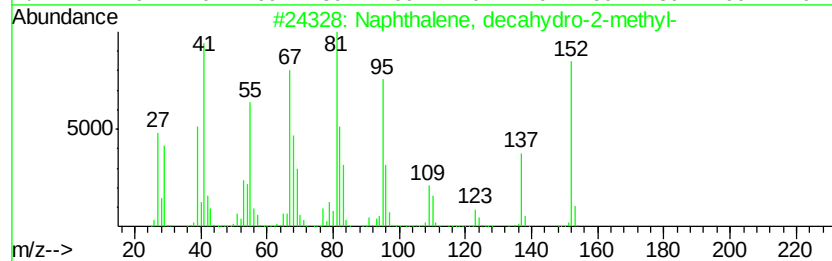
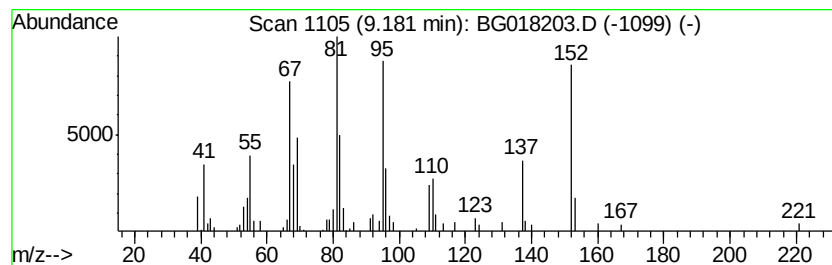
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Naphthalene, decahydro-2-me... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.18	2.86 ng/ul	87789	Naphthalene-d8	10.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	94
2			cis-Decalin, 2-syn-methyl-	152	C11H20	1000155-85-6	93
3			trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	91
4			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	83
5			Pulegone	152	C10H16O	000089-82-7	68



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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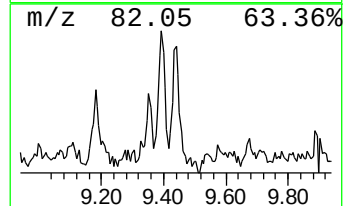
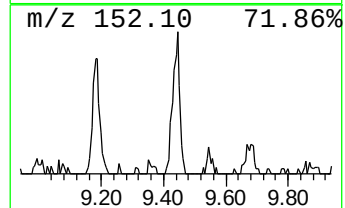
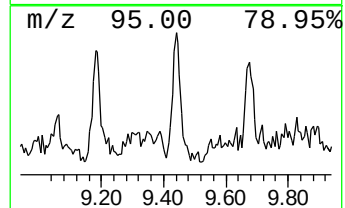
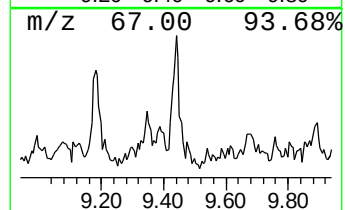
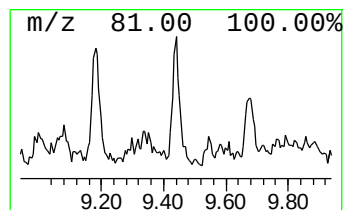
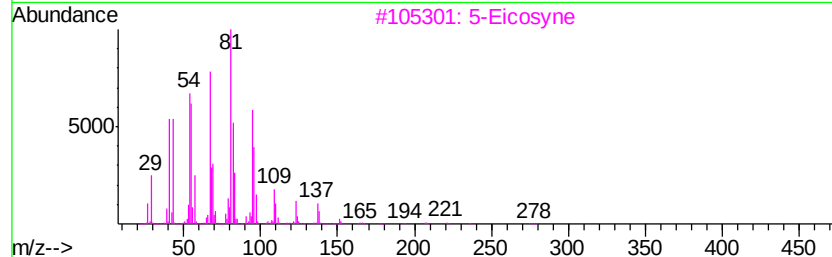
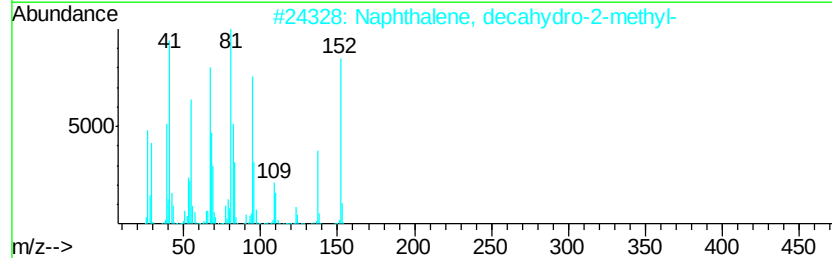
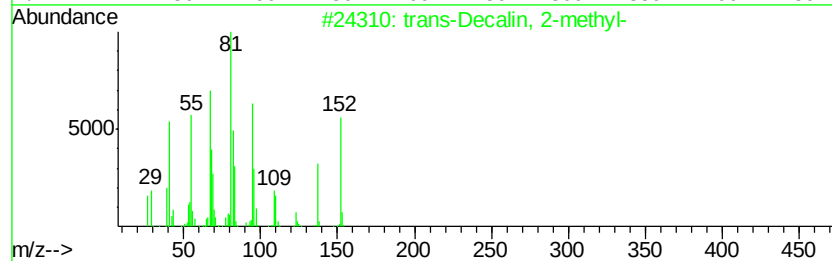
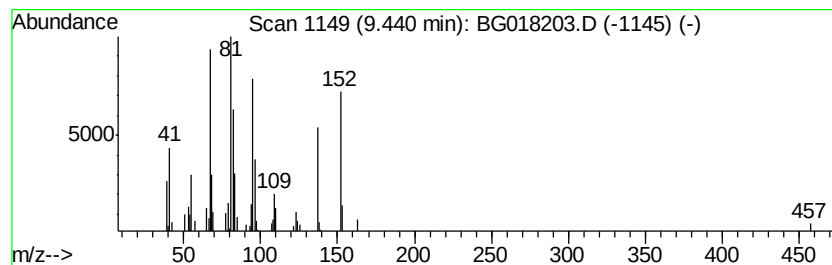
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 trans-Decalin, 2-methyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.44	3.15 ng/ul	96689	Naphthalene-d8	10.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	93
2		Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	90
3		5-Eicosyne	278	C20H38	074685-31-7	74
4		Decalin, anti-1-methyl-, cis-	152	C11H20	1000158-89-0	68
5		Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	64



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
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Instrument :
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ClientSampleId :
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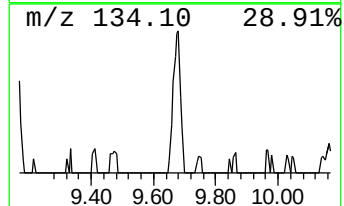
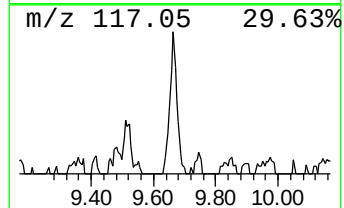
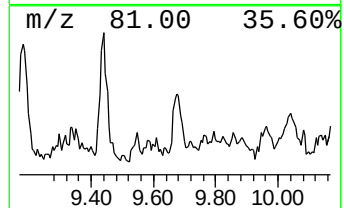
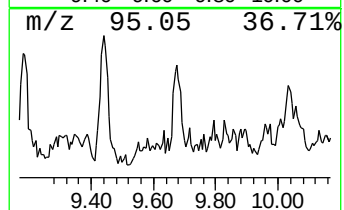
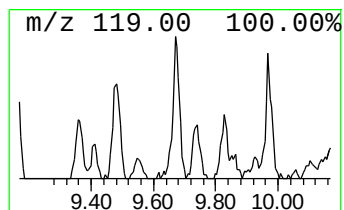
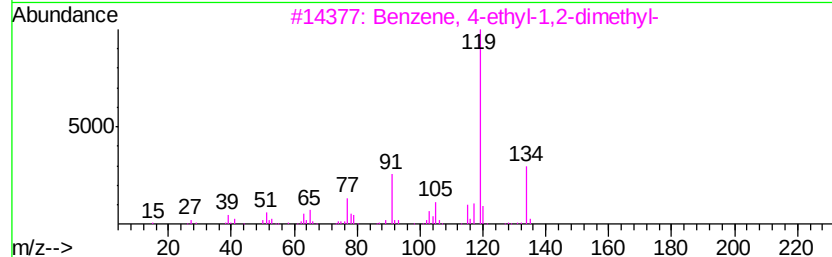
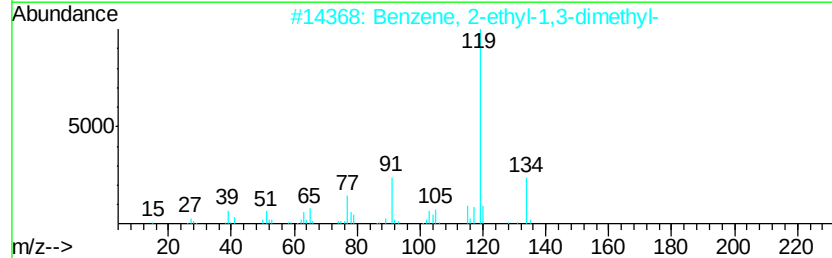
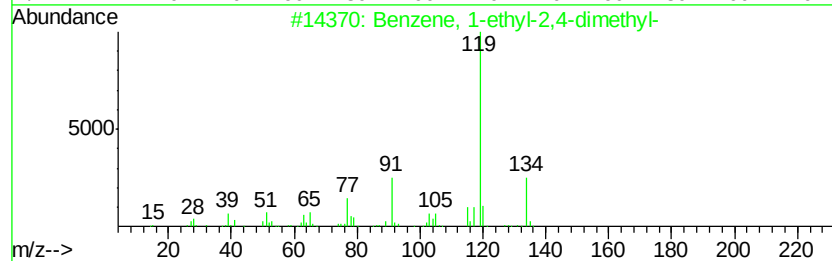
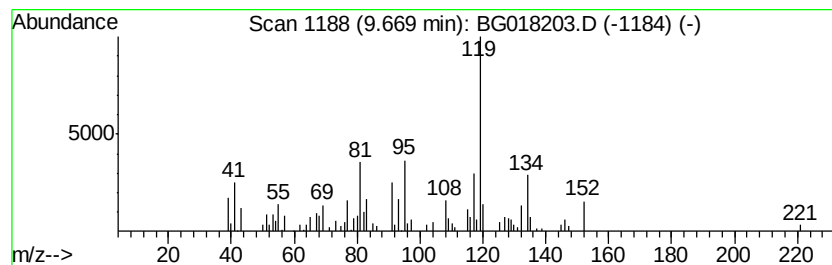
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.67	2.02 ng/ul	62003	Napthalene-d8	10.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	53
2		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	53
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	49
4		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	49
5		Benzene, 1-methyl-4-(1-methyleth...	134	C10H14	000099-87-6	49



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
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Instrument :
BNA_G
ClientSampleId :
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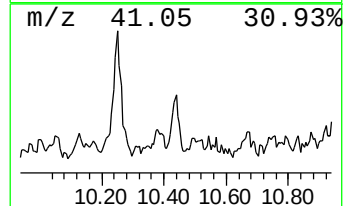
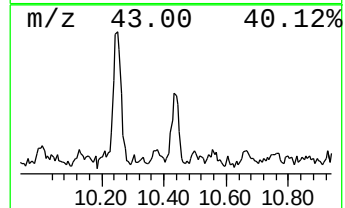
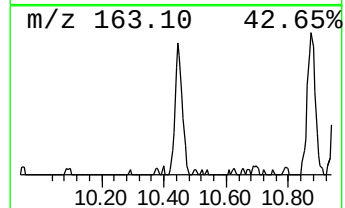
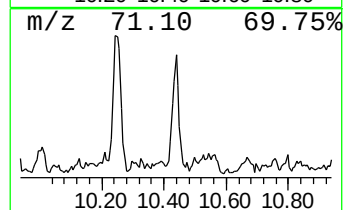
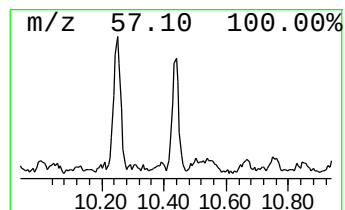
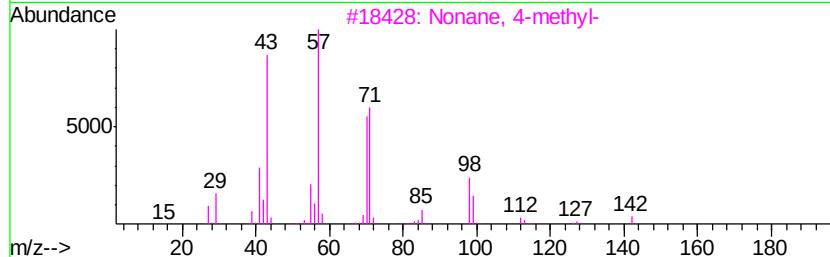
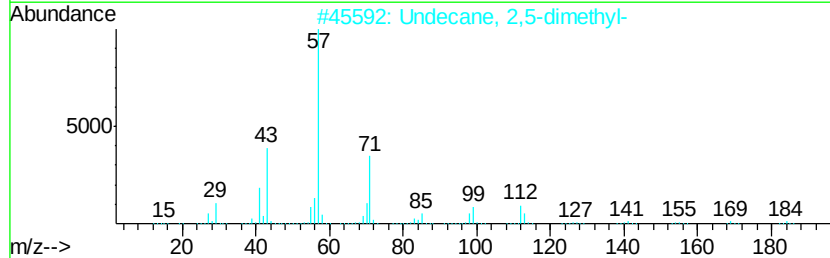
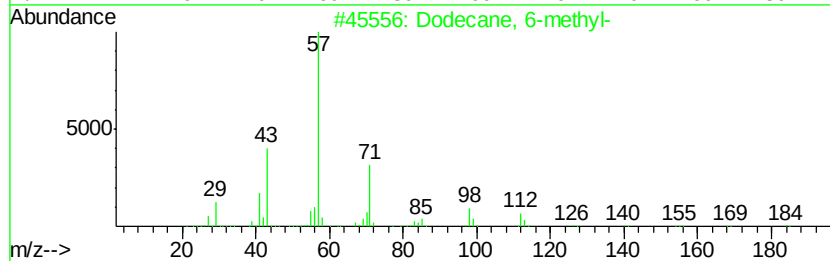
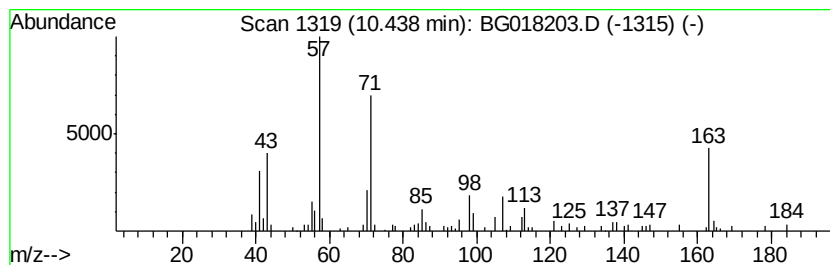
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.44	2.34 ng/ul	72019	Napthalene-d8	10.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane, 6-methyl-	184	C13H28	006044-71-9	52
2		Undecane, 2,5-dimethyl-	184	C13H28	017301-22-3	46
3		Nonane, 4-methyl-	142	C10H22	017301-94-9	38
4		Piperidin-3-one, 2,2,5-methyl-	141	C8H15NO	103634-57-7	35
5		Heptacosane	380	C27H56	000593-49-7	35



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
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Misc :
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ClientSampleId :
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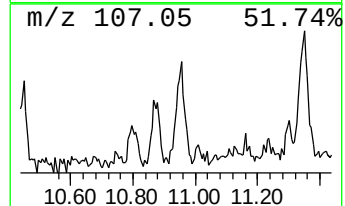
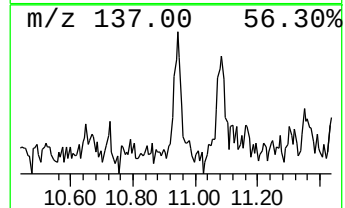
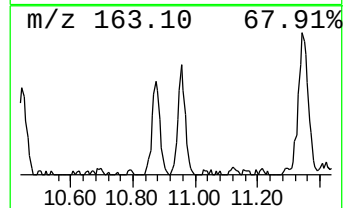
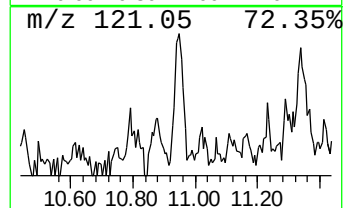
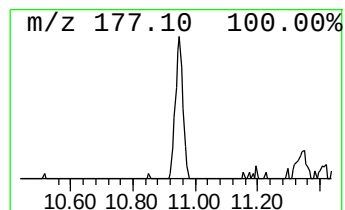
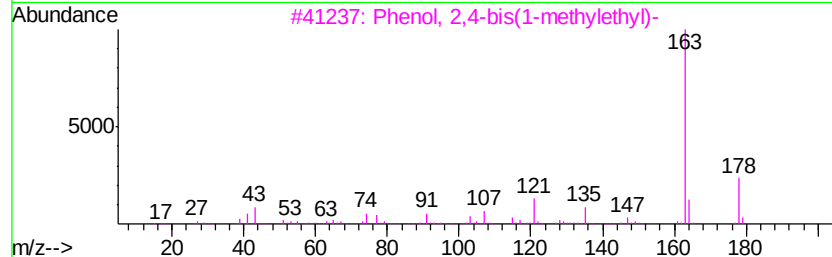
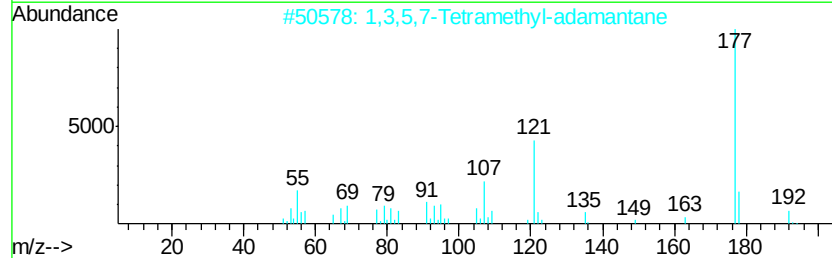
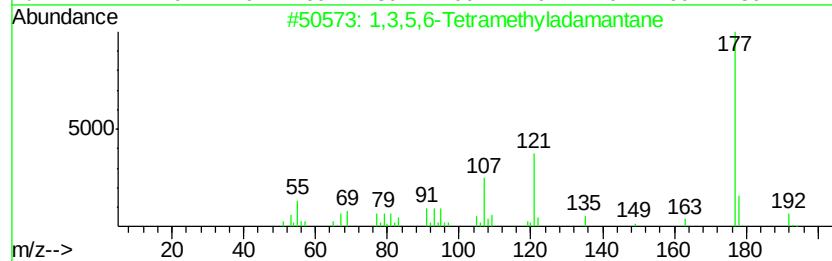
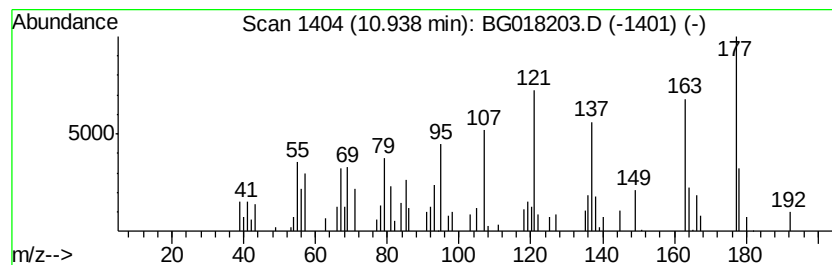
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 unknown-04 Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.94	2.68 ng/ul	82340	Napthalene-d8	10.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3,5,6-Tetramethyladamantane	192	C14H24	034694-70-7	30
2		1,3,5,7-Tetramethyl-adamantane	192	C14H24	001687-36-1	30
3		Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	27
4		Phenol, 2,5-bis(1-methylethyl)-	178	C12H18O	035946-91-9	27
5		1-Ethyl-3-propyladamantane	206	C15H26	019385-94-5	22



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
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Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

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ClientSampleId :
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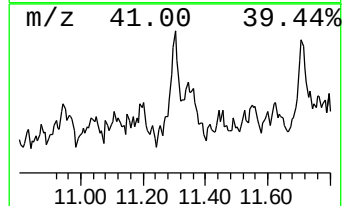
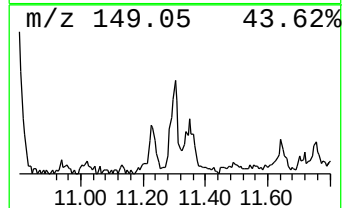
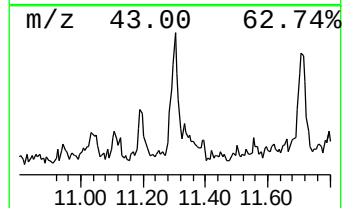
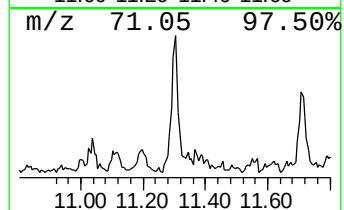
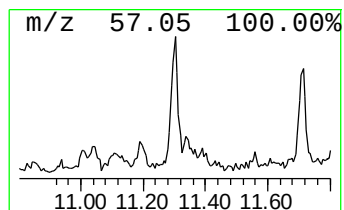
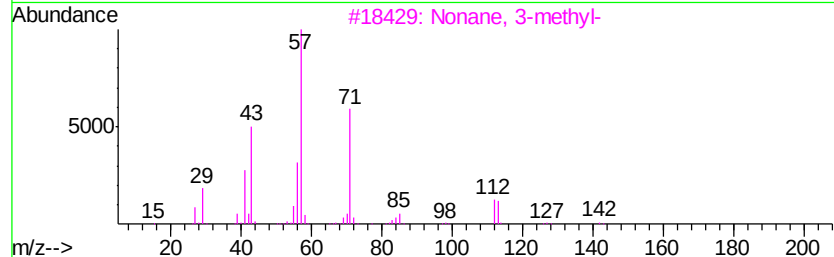
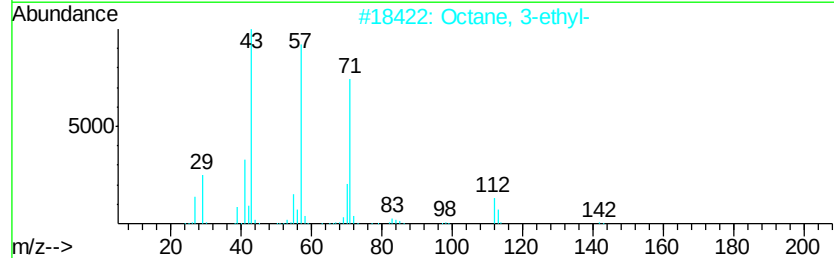
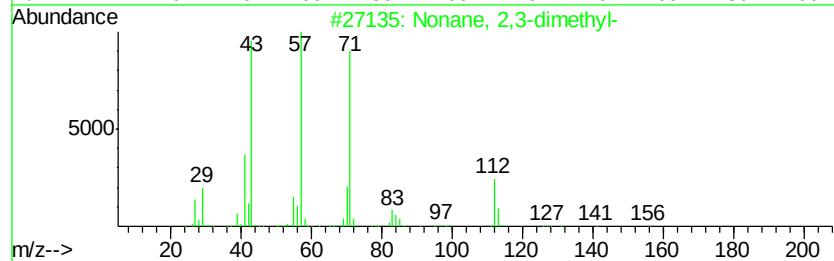
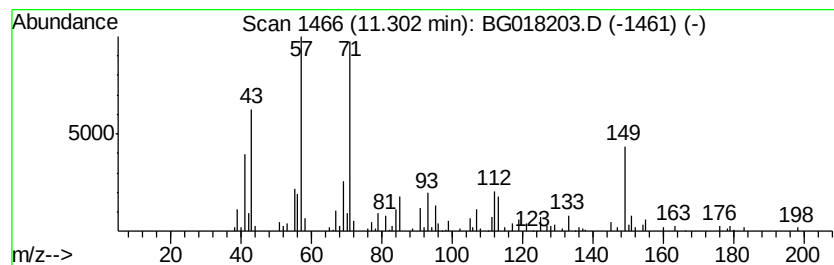
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.30	4.37 ng/ul	134204	Napthalene-d8	10.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 2,3-dimethyl-	156	C11H24	002884-06-2	47
2			Octane, 3-ethyl-	142	C10H22	005881-17-4	43
3			Nonane, 3-methyl-	142	C10H22	005911-04-6	43
4			Octane, 2-bromo-	192	C8H17Br	000557-35-7	38
5			Tridecane, 7-methyl-	198	C14H30	026730-14-3	38



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
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Misc :
ALS Vial : 23 Sample Multiplier: 1

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ClientSampleId :
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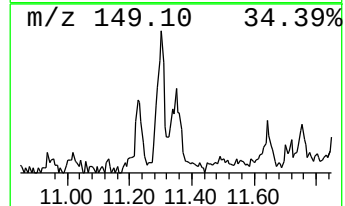
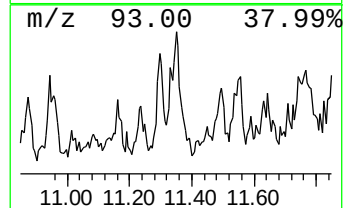
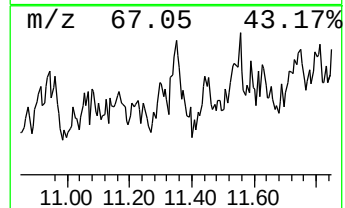
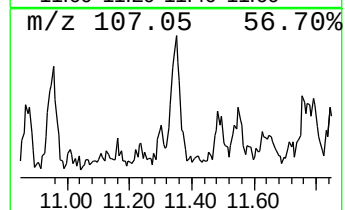
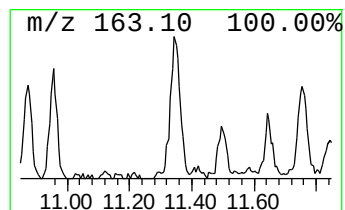
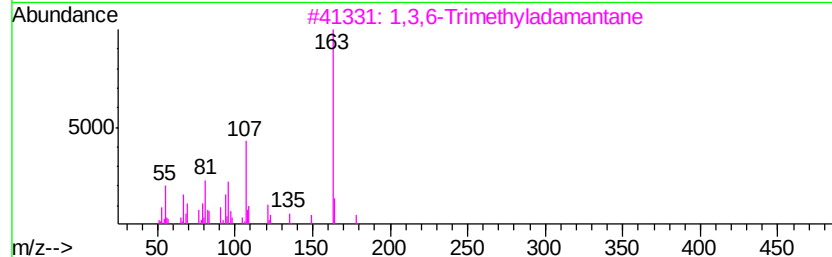
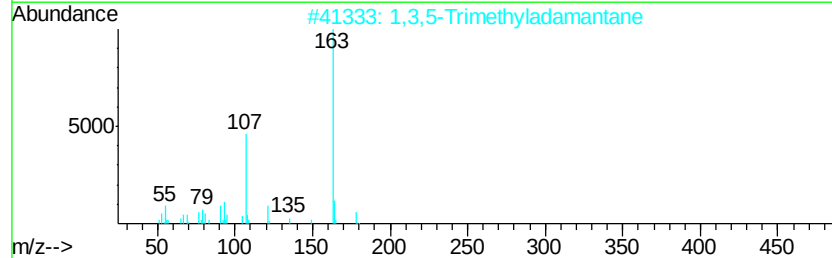
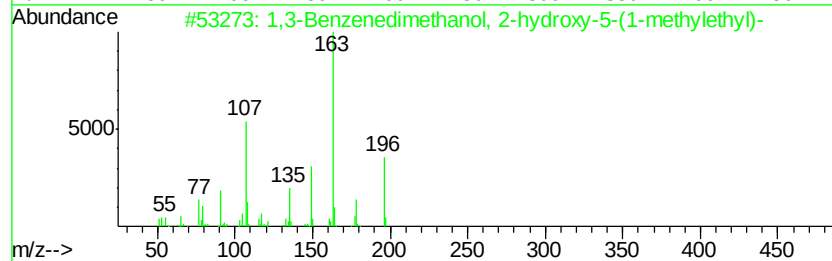
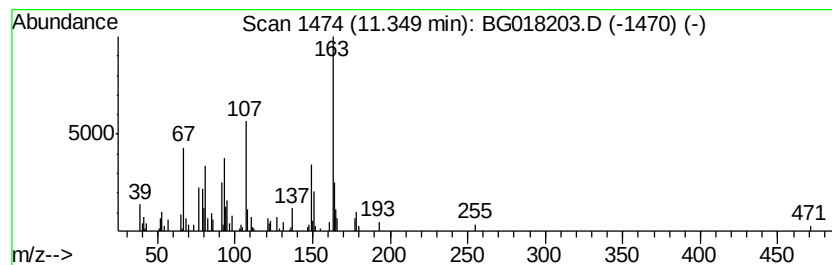
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 unknown-05 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.35	4.17 ng/ul	128044	Naphthalene-d8	10.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Benzenedimethanol, 2-hydroxy...	196	C11H16O3	054845-41-9	47
2		1,3,5-Trimethyladamantane	178	C13H22	000707-35-7	46
3		1,3,6-Trimethyladamantane	178	C13H22	024139-37-5	43
4		Bicyclo[4.3.0]non-3-ene, 3,4,7-t...	164	C12H20	1000155-78-7	38
5		4-[[3-(3,5-Dimethyl-pyrazol-1-yl...	326	C16H18N6O2	1000296-14-6	38



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

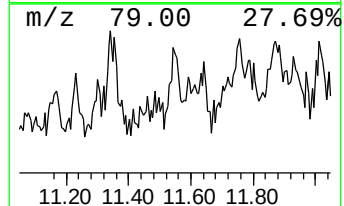
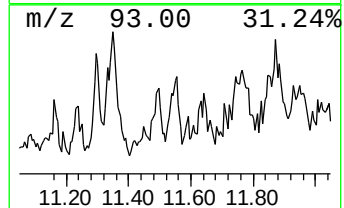
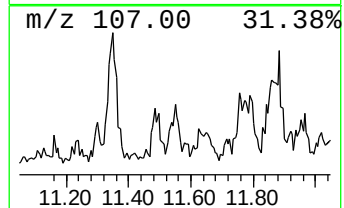
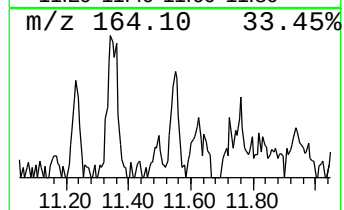
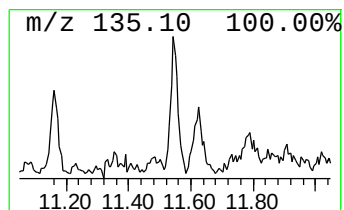
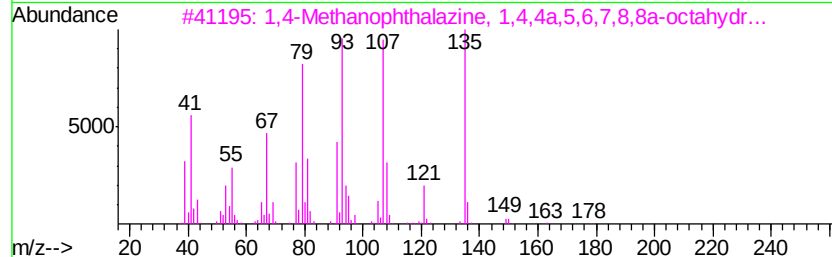
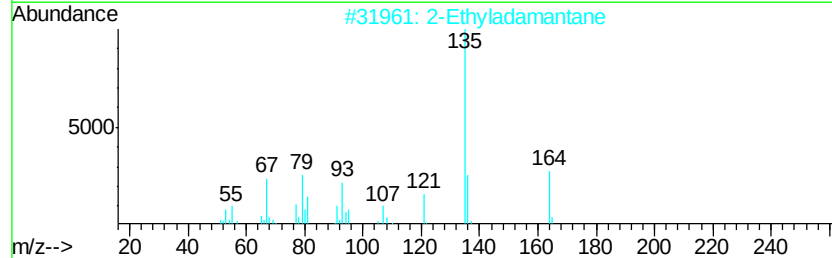
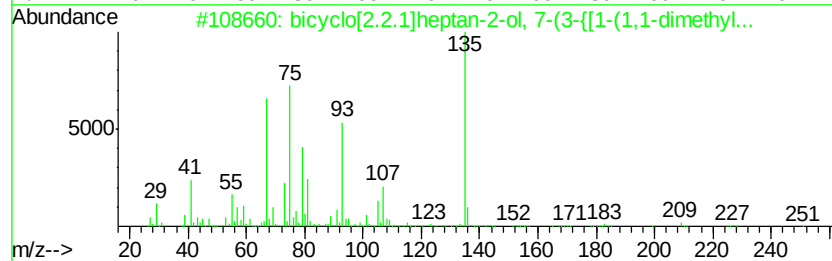
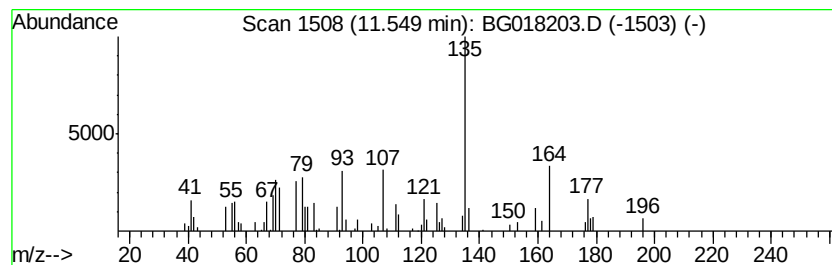
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG073115.M
Quant Title : SVOA CALIBRATION

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

Peak Number 18 unknown-06 Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.55	2.01 ng/ul	61667	Naphthalene-d8	10.26
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	bicyclo[2.2.1]heptan-2-ol, 7-(3-...	284 C16H32O2Si	1000196-96-1	38
2	2-Ethyladamantane	164 C12H20	014451-87-7	38
3	1,4-Methanophthalazine, 1,4,4a,5...	178 C11H18N2	109746-13-6	27
4	Phenol, 4-(1,1-dimethylpropyl)-	164 C11H16O	000080-46-6	27
5	2-Iodoadamantane	262 C10H15I	018971-91-0	27



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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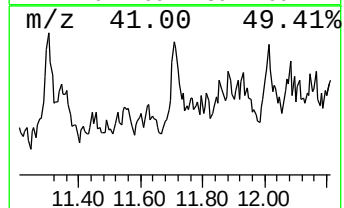
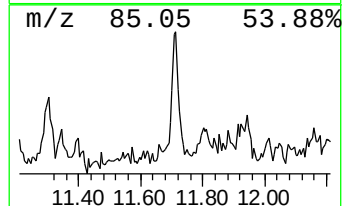
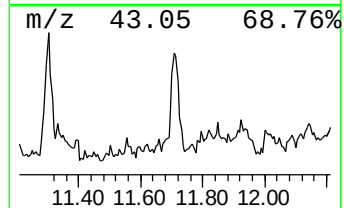
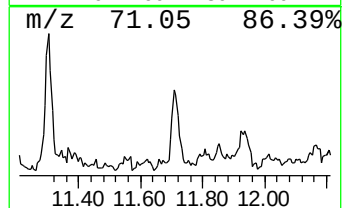
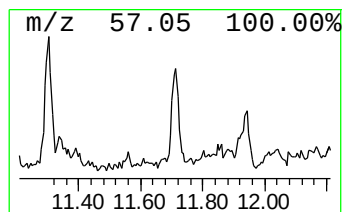
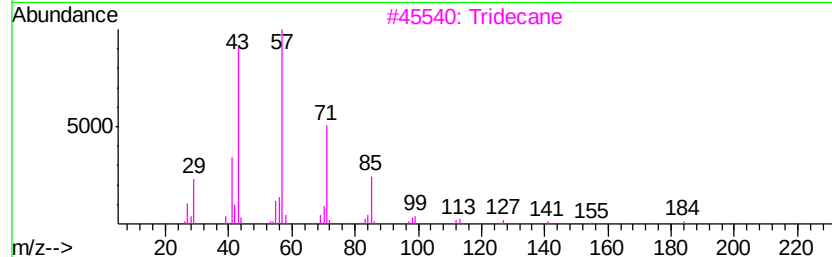
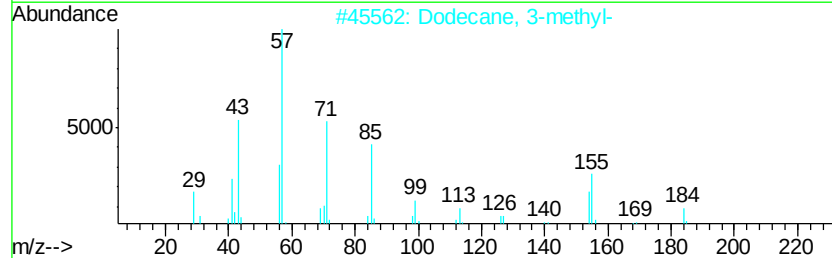
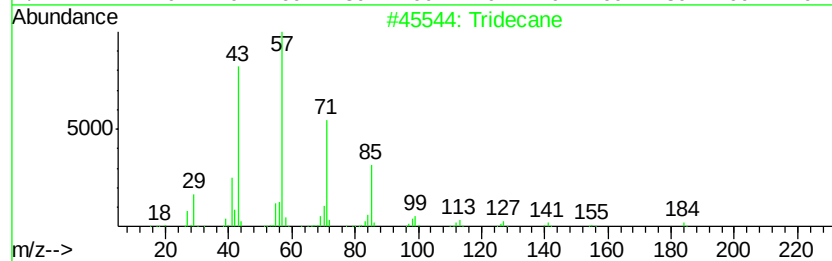
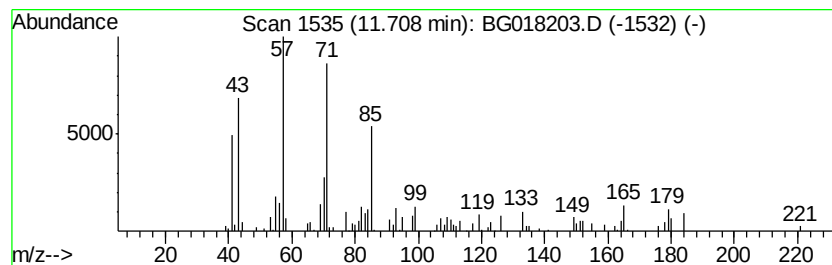
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.71	3.11 ng/ul	95523	Napthalene-d8	10.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	83
2			Dodecane, 3-methyl-	184	C13H28	017312-57-1	81
3			Tridecane	184	C13H28	000629-50-5	68
4			Nonadecane	268	C19H40	000629-92-5	64
5			Tetradecane	198	C14H30	000629-59-4	64



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

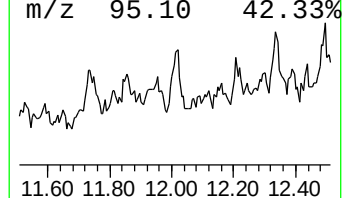
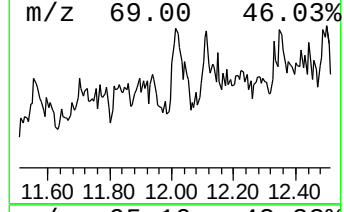
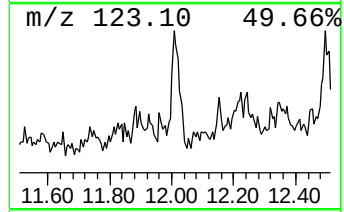
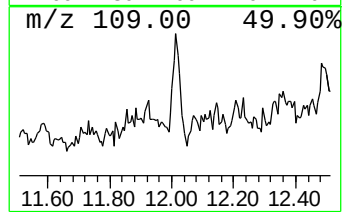
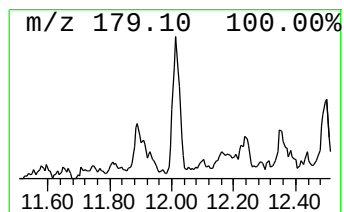
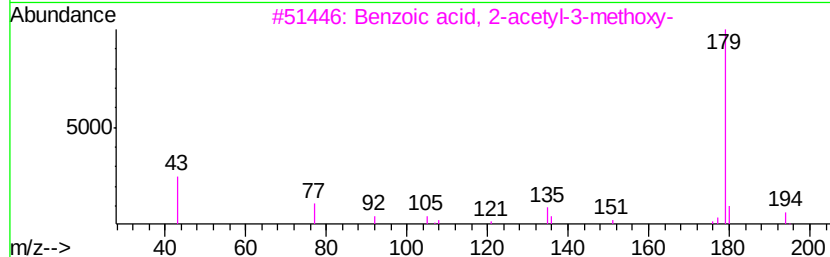
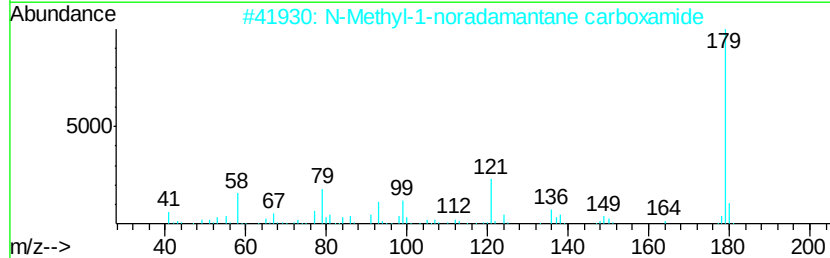
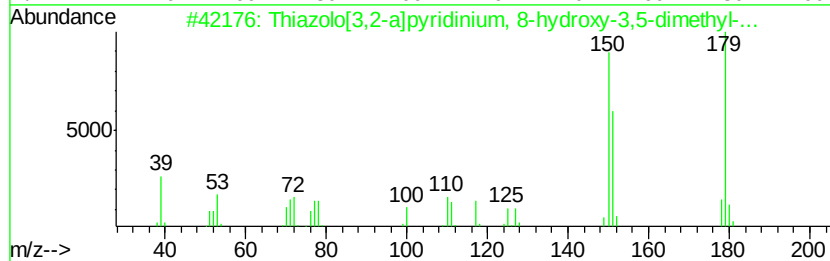
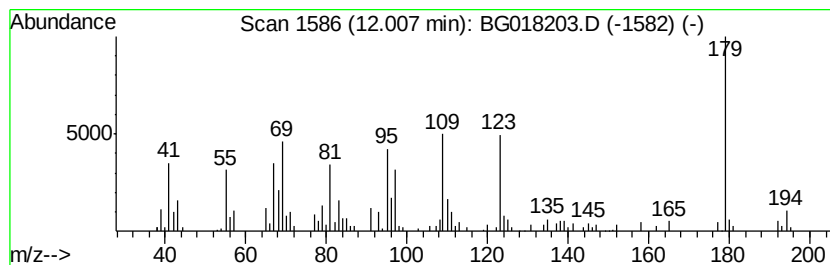
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG073115.M
Quant Title : SVOA CALIBRATION

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

Peak Number 20 unknown-07 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	4.08 ng/ul	125507	Napthalene-d8	10.26
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Thiazolo[3,2-a]pyridinium, 8-hyd...	179 C9H9NOS	030277-00-0 27
2		N-Methyl-1-noradamantane carboxa...	179 C11H17NO	1000211-82-0 22
3		Benzoic acid, 2-acetyl-3-methoxy-	194 C10H10O4	120714-42-3 22
4		Anthracene, 9,10-dihydro-9-methyl-	194 C15H14	017239-99-5 22
5		Benzoic acid, 4-nitroso-, ethyl ...	179 C9H9NO3	007476-79-1 22



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

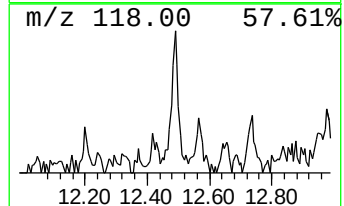
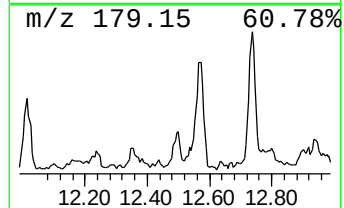
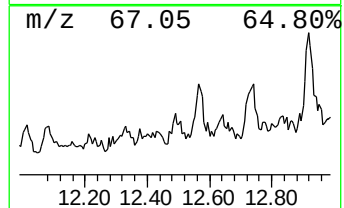
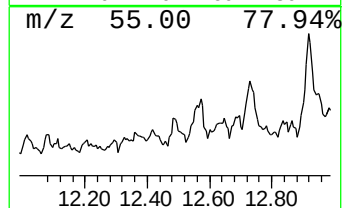
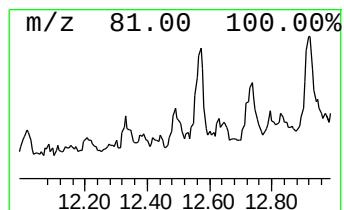
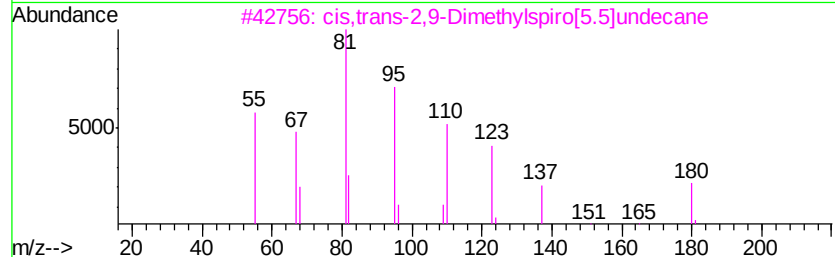
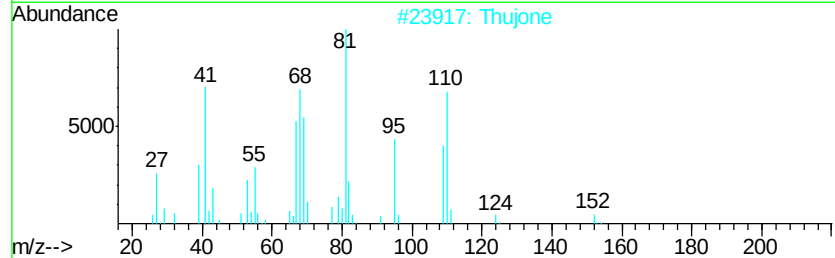
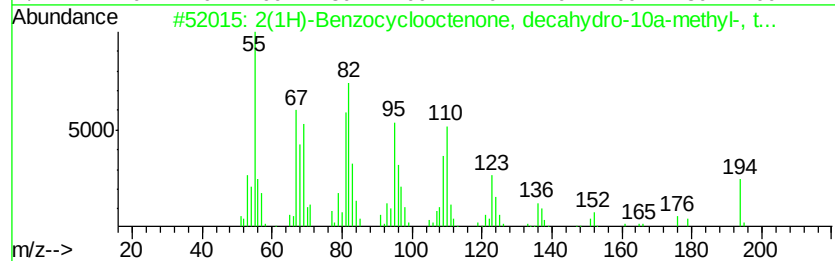
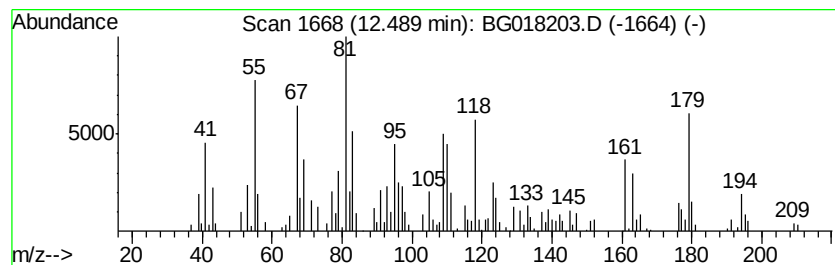
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG073115.M
Quant Title : SVOA CALIBRATION

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

Peak Number 21 2(1H)-Benzocyclooctenone, d... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.49	2.82 ng/ul	109743	Acenaphthene-d10	14.16
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		2(1H)-Benzocyclooctenone, decahy...	194 C13H22O	055103-68-9 52
2		Thujone	152 C10H16O	000546-80-5 41
3		cis,trans-2,9-Dimethylspiro[5.5]...	180 C13H24	095530-74-8 41
4		3-Heptyne, 5-methyl-	110 C8H14	061228-09-9 35
5		Cyclopentane, 1-ethenyl-3-ethyl-...	138 C10H18	055170-98-4 35



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampled :
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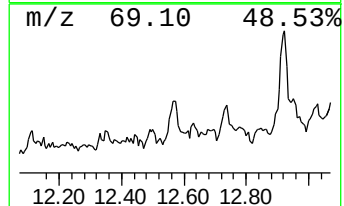
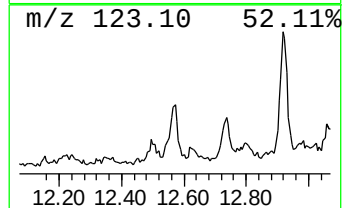
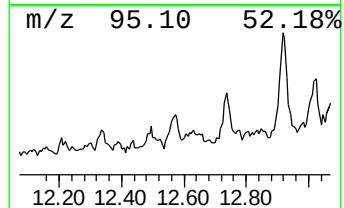
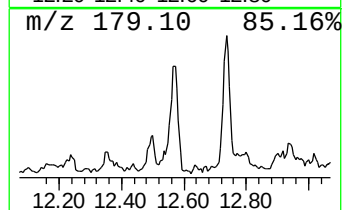
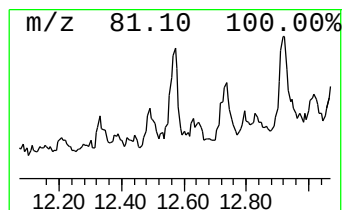
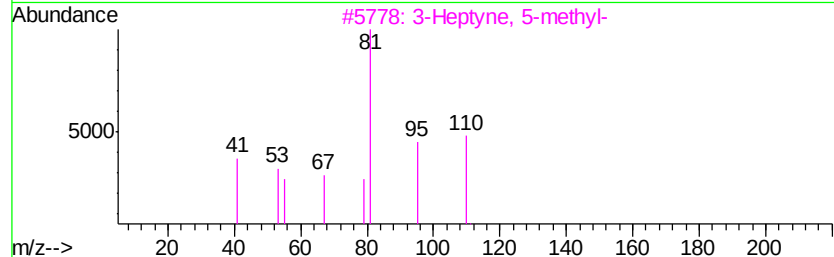
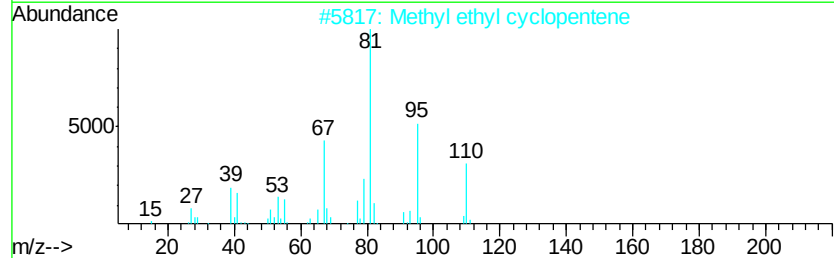
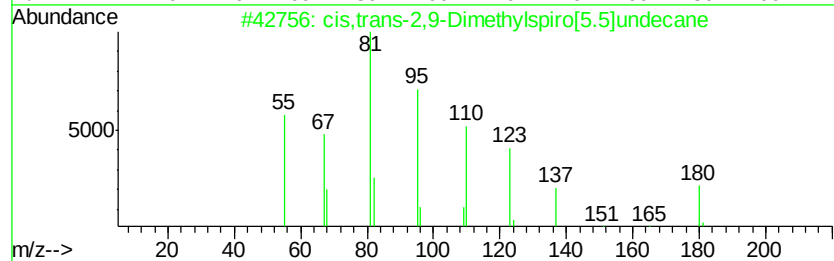
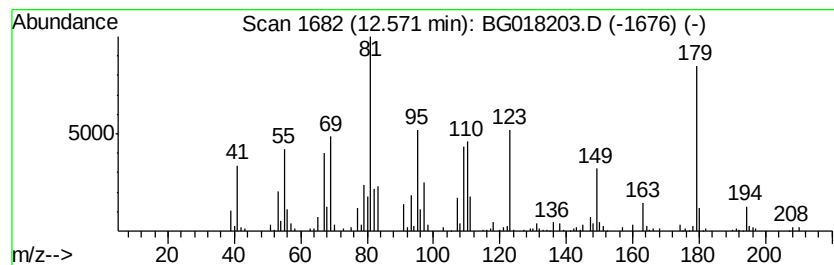
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 cis,trans-2,9-Dimethylspiro... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.57	5.68 ng/ul	220716	Acenaphthene-d10	14.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis,trans-2,9-Dimethylspiro[5.5]undecane	180	C13H24	095530-74-8	55
2			Methyl ethyl cyclopentene	110	C8H14	019780-56-4	43
3			3-Heptyne, 5-methyl-	110	C8H14	061228-09-9	30
4			Cis-3-methyl-endo-tricyclo[5.2.1]undecane	150	C11H18	1000215-29-0	27
5			Cyclohexane, ethylidene-	110	C8H14	001003-64-1	27



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
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Acq On : 1 Aug 2015 4:38
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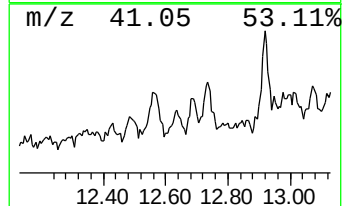
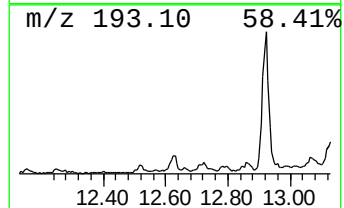
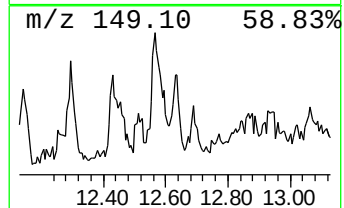
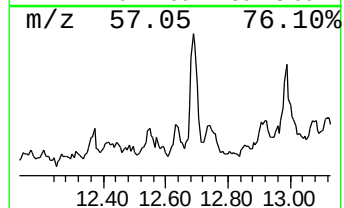
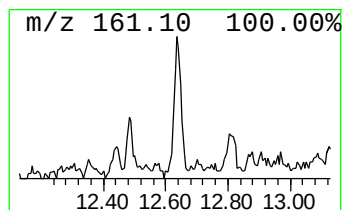
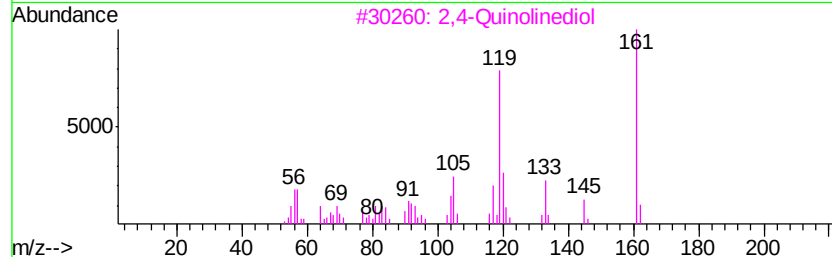
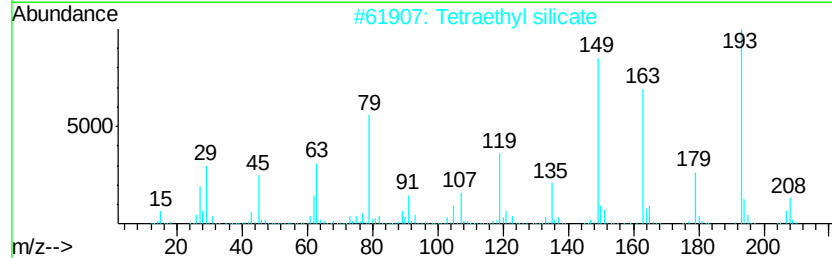
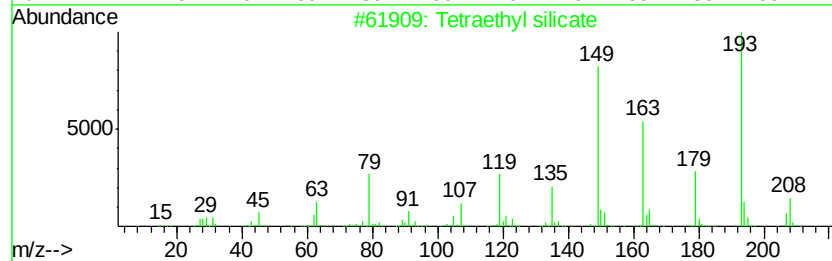
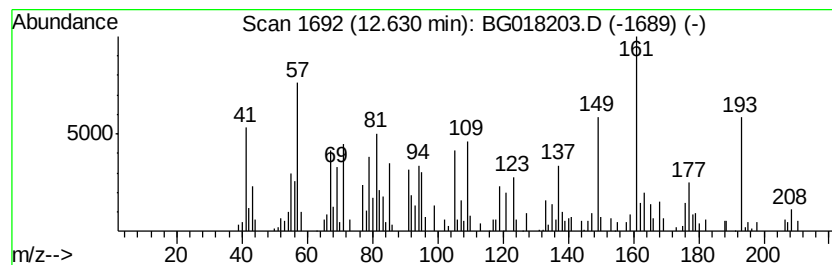
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 unknown-08 Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.63	2.63 ng/ul	102282	Acenaphthene-d10	14.16
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tetraethyl silicate	208 C8H2004Si	000078-10-4 20
2		Tetraethyl silicate	208 C8H2004Si	000078-10-4 20
3		2,4-Quinolinediol	161 C9H7N02	000086-95-3 11
4		2,4-2H-Benzo[c]furanone, 3,3,4-t...	176 C11H12O2	146950-80-3 11
5		Ethanone, 1-(4-butylphenyl)-2-ch...	210 C12H15ClO	1000195-58-2 10



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

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ClientSampleId :
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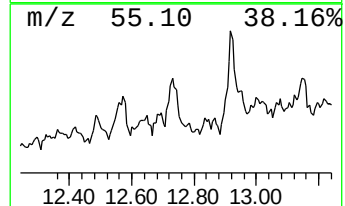
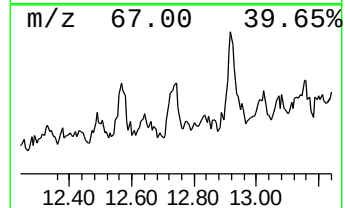
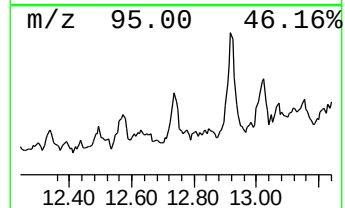
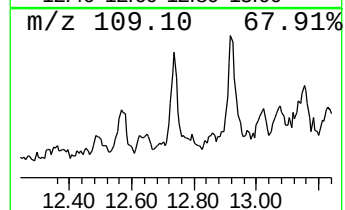
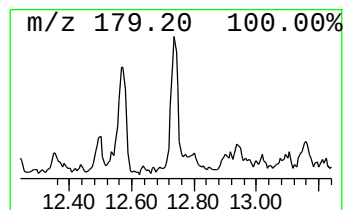
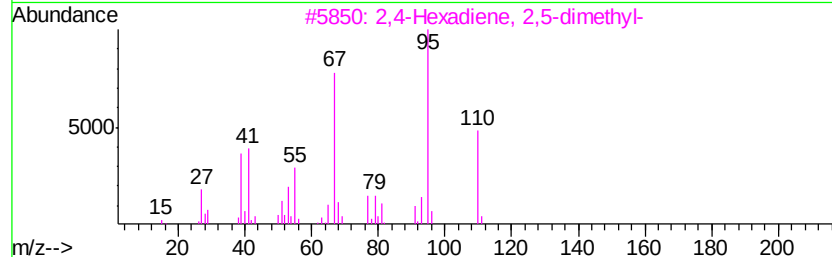
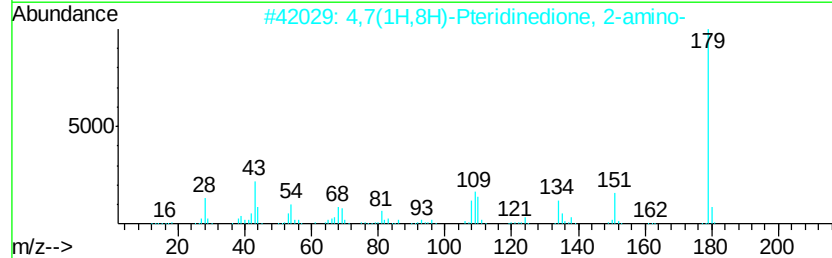
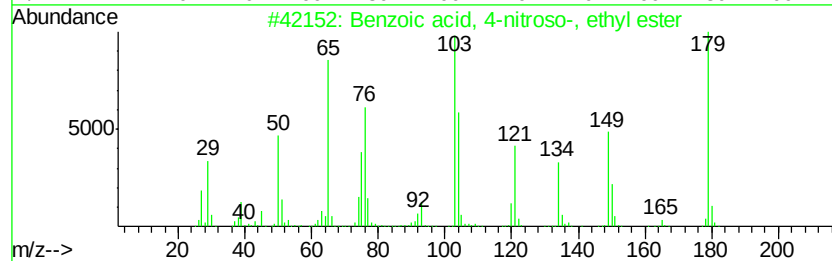
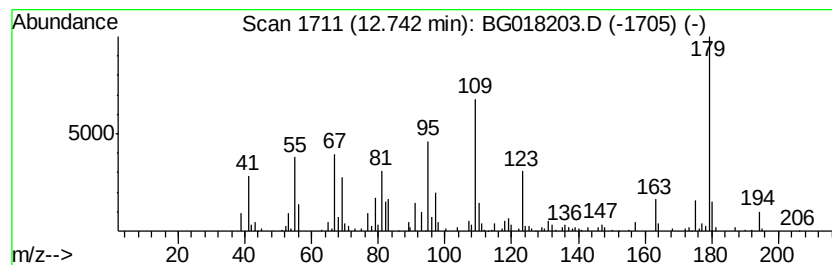
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 unknown-09 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	5.64 ng/ul	219283	Acenaphthene-d10	14.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, 4-nitroso-, ethyl ...	179	C9H9NO3	007476-79-1	35
2			4,7(1H,8H)-Pteridinedione, 2-amino-	179	C6H5N5O2	000529-69-1	27
3			2,4-Hexadiene, 2,5-dimethyl-	110	C8H14	000764-13-6	25
4			Benzoic acid, 2-acetyl-3-methoxy-	194	C10H10O4	120714-42-3	18
5			1,2,3-Trimethyl-cyclopent-2-enec...	138	C9H14O	1000190-18-1	18



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampled :
C02A6

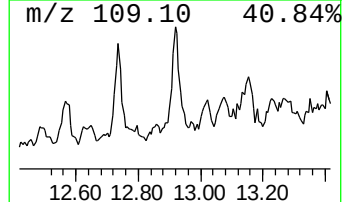
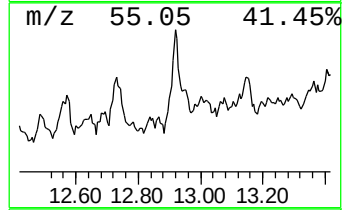
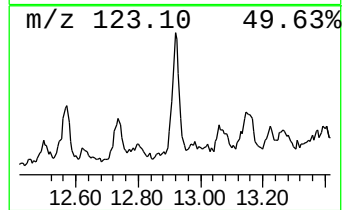
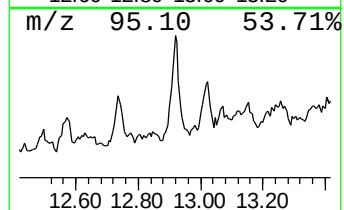
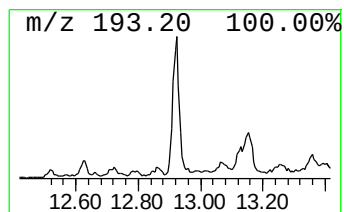
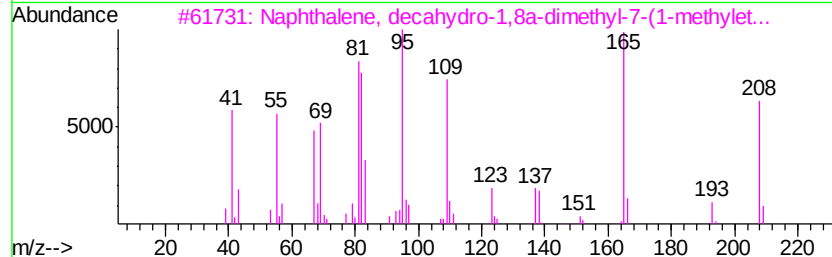
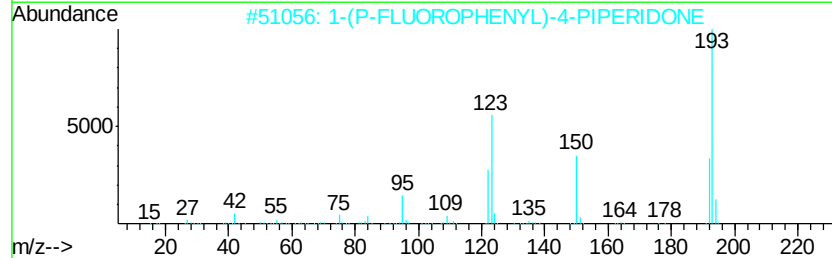
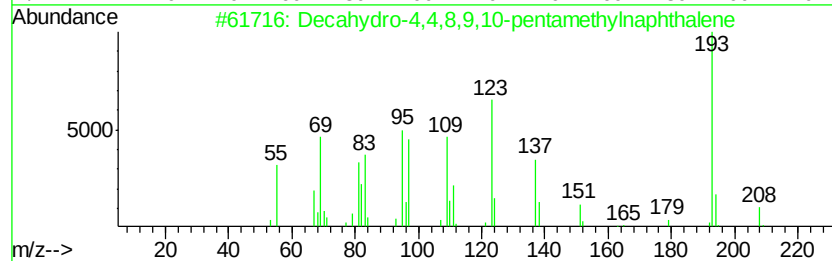
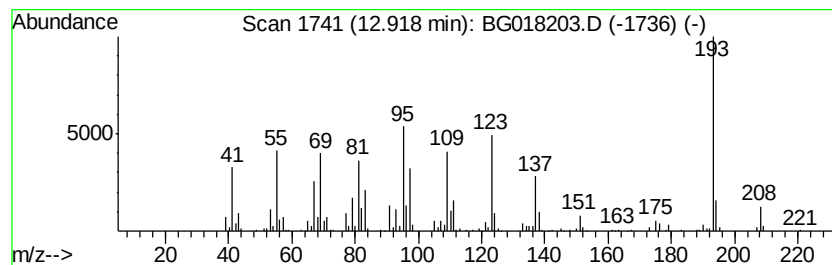
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 Decahydro-4,4,8,9,10-pentam... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.92	10.88 ng/ul	422613	Acenaphthene-d10	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	87
2		1-(P-FLUOROPHENYL)-4-PIPERIDONE	193	C11H12FNO	1000238-56-7	43
3		Naphthalene, decahydro-1,8a-dime...	208	C15H28	015404-63-4	30
4		10.alpha.-Eremophilane	208	C15H28	003242-05-5	30
5		Borinic acid, diethyl-, 3,3,5-tr...	208	C13H25BO	057387-76-5	27



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
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ClientSampleId :
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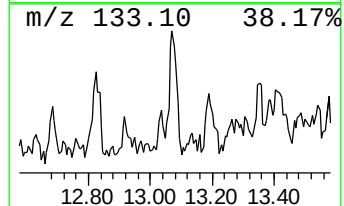
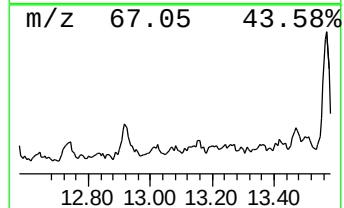
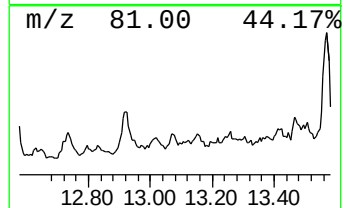
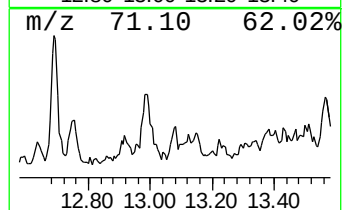
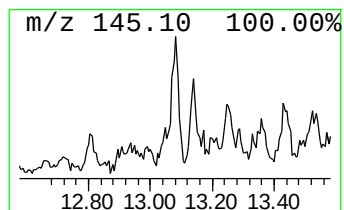
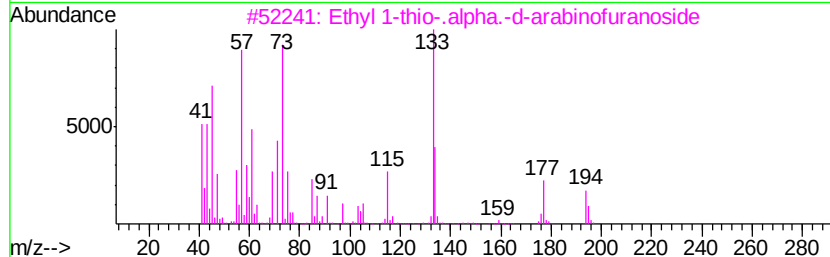
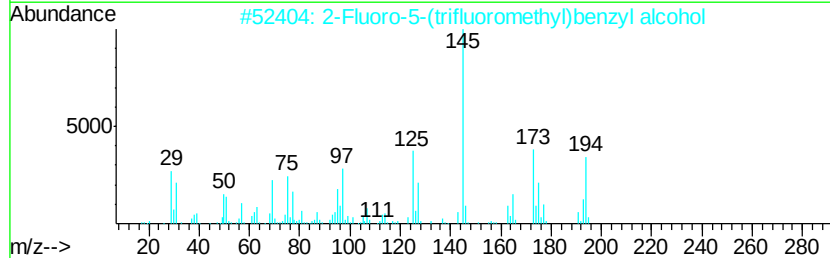
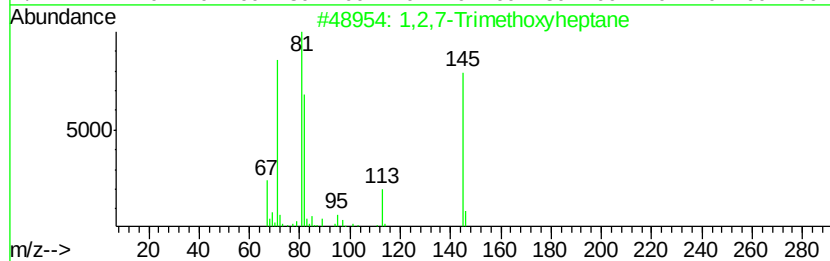
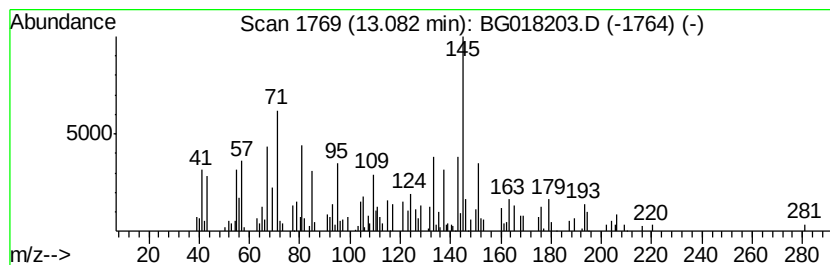
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 unknown-10 Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.08	2.54 ng/ul	98508	Acenaphthene-d10	14.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2,7-Trimethoxyheptane	190	C10H22O3	062635-57-8	38
2			2-Fluoro-5-(trifluoromethyl)benz...	194	C8H6F4O	207974-09-2	15
3			Ethyl 1-thio-.alpha.-d-arabinofu...	194	C7H14O4S	1000129-83-7	12
4			Cyclohexene, 1-pentyl-	152	C11H20	015232-85-6	10
5			4-Heptanone, 3-methyl-	128	C8H16O	015726-15-5	10



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

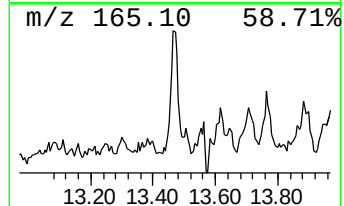
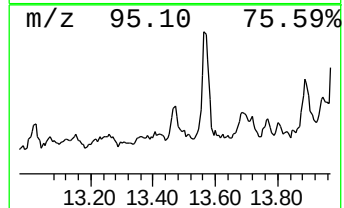
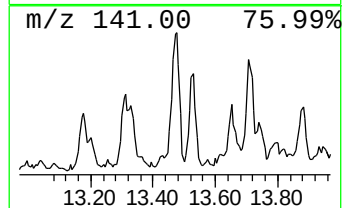
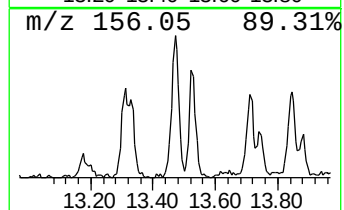
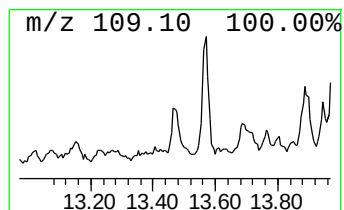
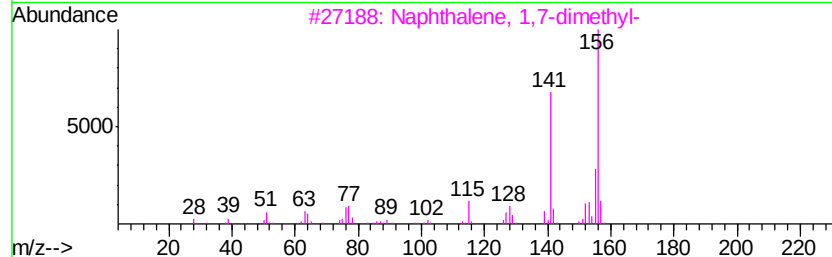
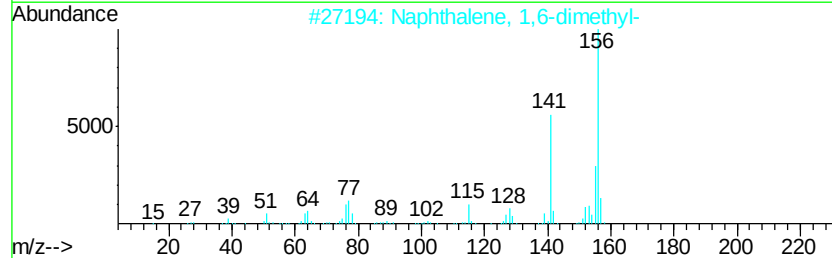
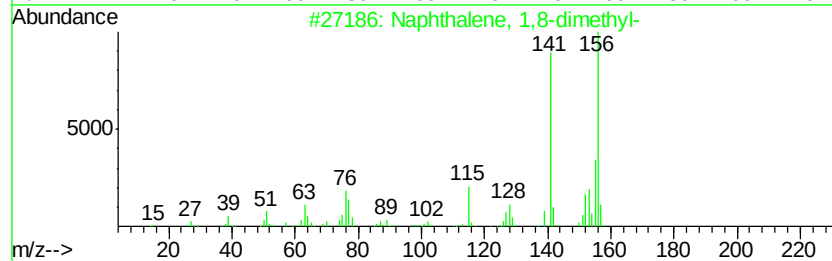
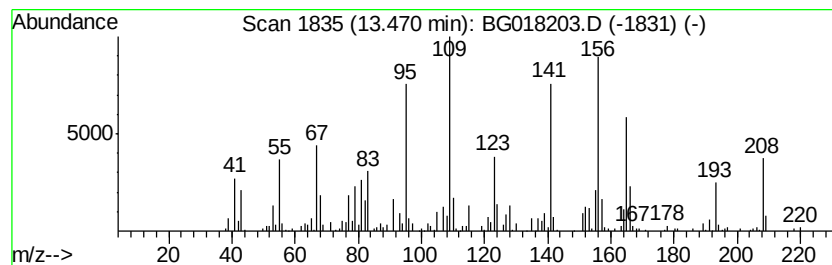
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 Naphthalene, 1,8-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	6.31 ng/ul	245317	Acenaphthene-d10	14.16
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,8-dimethyl-	156 C12H12	000569-41-5 51
2		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 51
3		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 47
4		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 47
5		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 42



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
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Acq On : 1 Aug 2015 4:38
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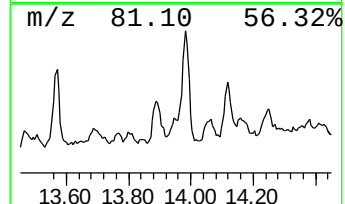
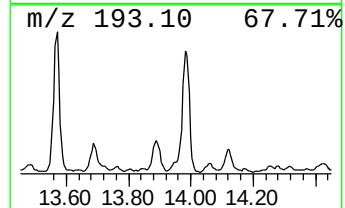
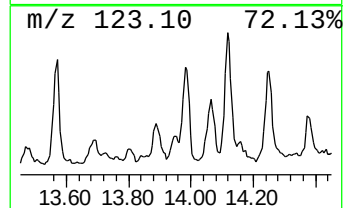
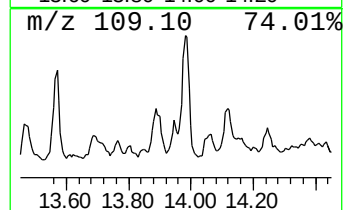
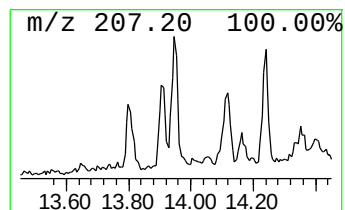
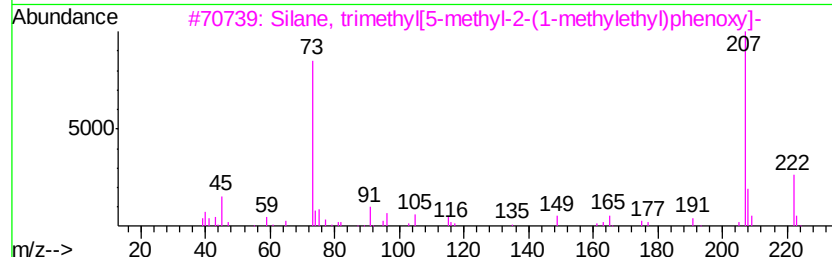
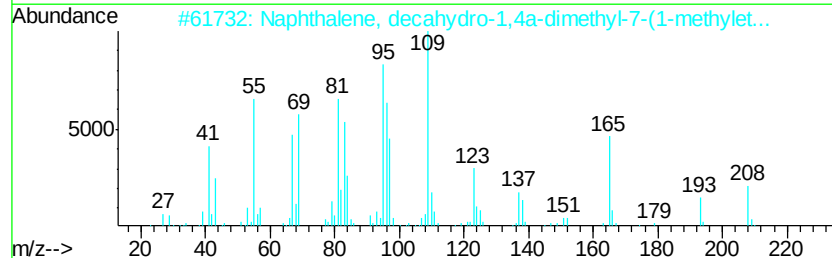
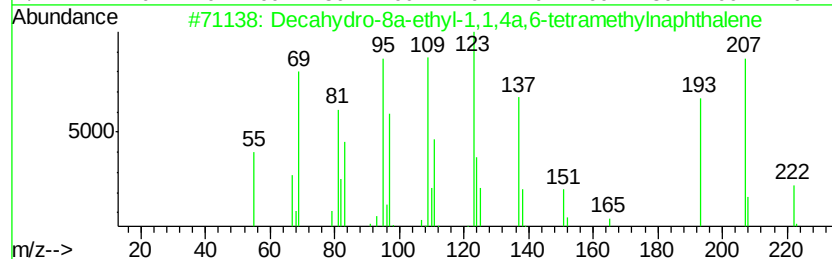
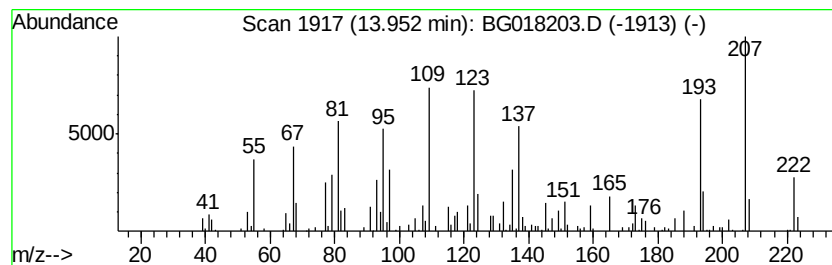
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 unknown-11 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	5.18 ng/ul	201212	Acenaphthene-d10	14.16
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Decahydro-8a-ethyl-1,1,4a,6-tetr...	222 C16H30	1000100-23-6 49
2		Naphthalene, decahydro-1,4a-dime...	208 C15H28	030824-81-8 22
3		Silane, trimethyl[5-methyl-2-(1-...	222 C13H22OSi	055012-80-1 20
4		2,3-Dimethyl-8-oxo-non-2-enal	182 C11H18O2	1000186-82-6 18
5		1,2-Benzenediol, 3,5-bis(1,1-dim...	222 C14H22O2	001020-31-1 18



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
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Operator : TP/UM
Sample : G3065-12
Misc :
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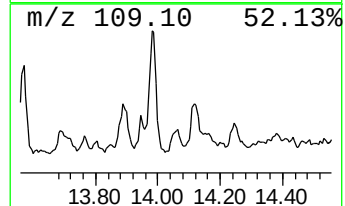
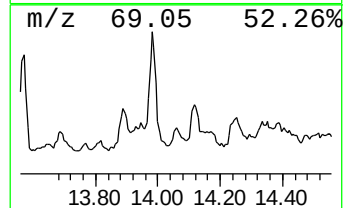
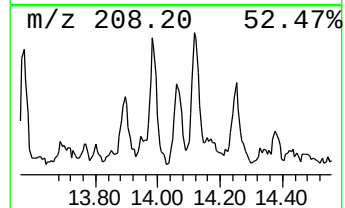
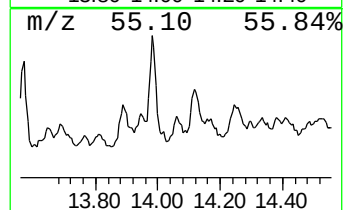
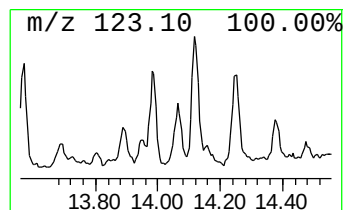
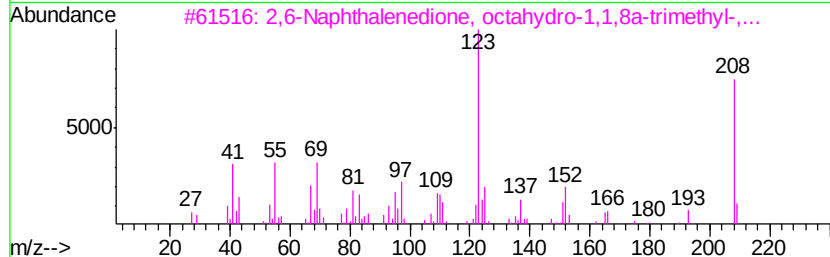
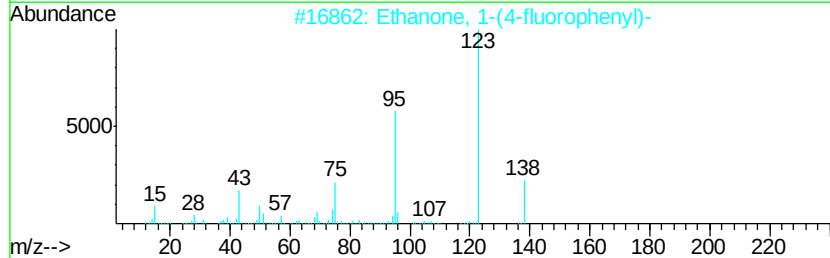
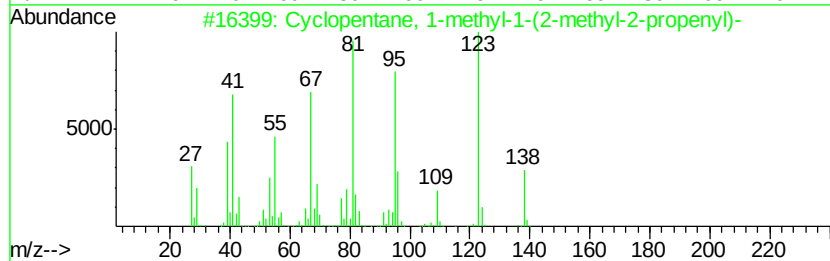
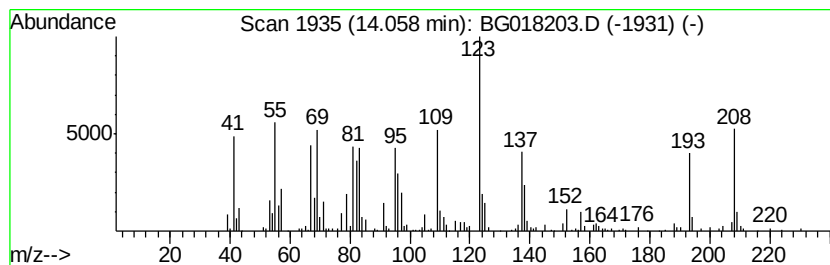
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Quant Title : SVOA CALIBRATION

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

Peak Number 30 unknown-12 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.06	8.35 ng/ul	324286	Acenaphthene-d10	14.16
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclopentane, 1-methyl-1-(2-meth...	138 C10H18	074764-47-9 35
2		Ethanone, 1-(4-fluorophenyl)-	138 C8H7FO	000403-42-9 27
3		2,6-Naphthalenedione, octahydro-...	208 C13H20O2	057289-17-5 25
4		Bicyclo[3.1.1]heptane, 2,6,6-tri...	138 C10H18	004755-33-3 25
5		4-Fluorobenzoic acid, 2-methylpr...	196 C11H13FO2	029240-31-1 22



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
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Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

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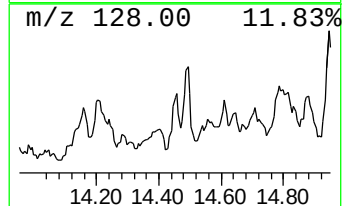
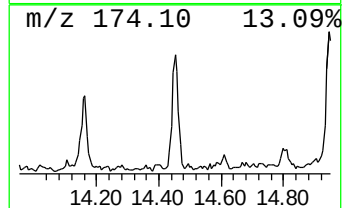
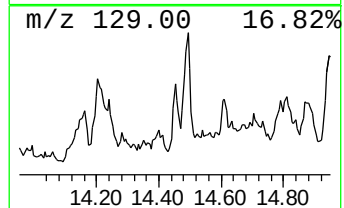
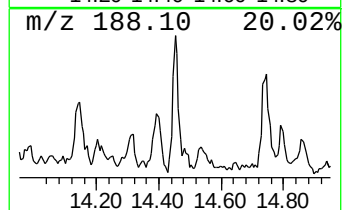
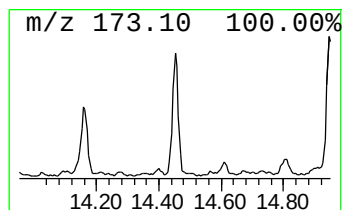
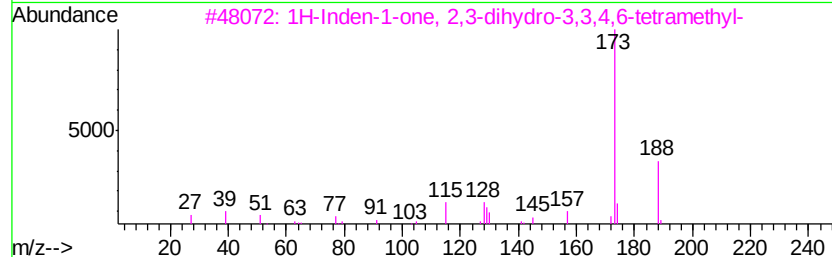
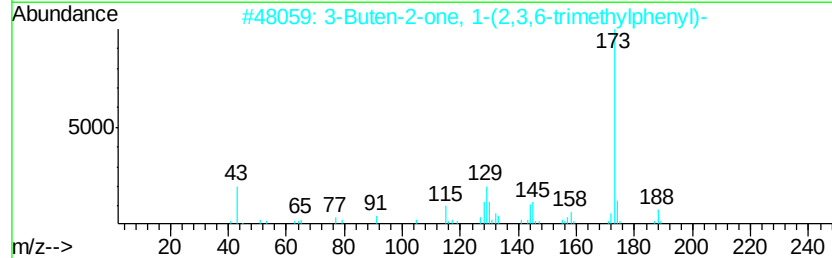
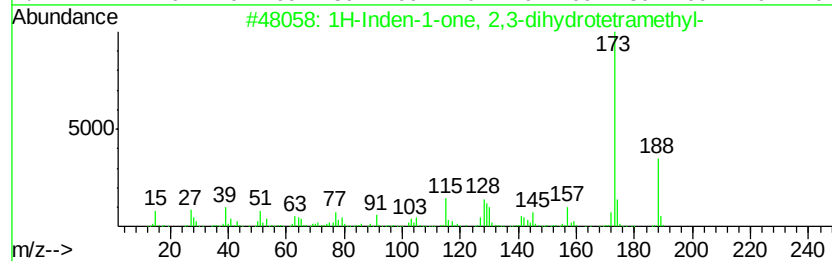
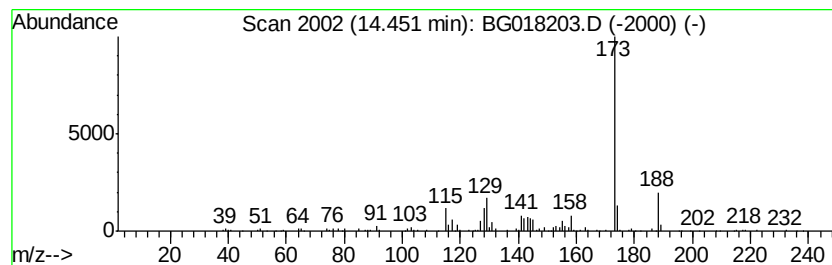
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Quant Title : SVOA CALIBRATION

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

Peak Number 31 1H-Inden-1-one, 2,3-dihydro... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.45	2.86 ng/ul	111267	Acenaphthene-d10	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Inden-1-one, 2,3-dihydrotetra...	188	C13H16O	089907-33-5	87
2		3-Buten-2-one, 1-(2,3,6-trimethy...	188	C13H16O	054789-45-6	87
3		1H-Inden-1-one, 2,3-dihydro-3,3,...	188	C13H16O	055255-42-0	87
4		1H-Inden-1-one, 2,3-dihydro-3,3,...	188	C13H16O	056298-98-7	87
5		1H-Inden-1-one, 2,3-dihydro-3,3,...	188	C13H16O	054789-22-9	86



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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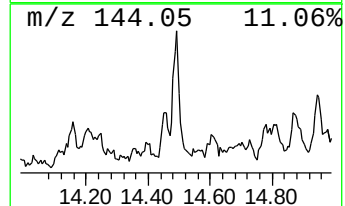
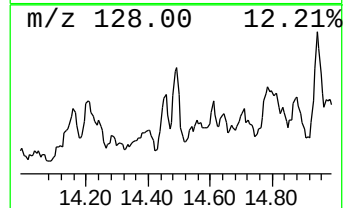
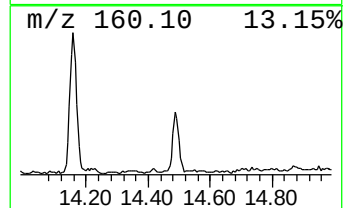
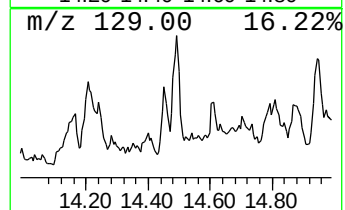
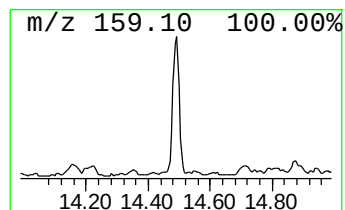
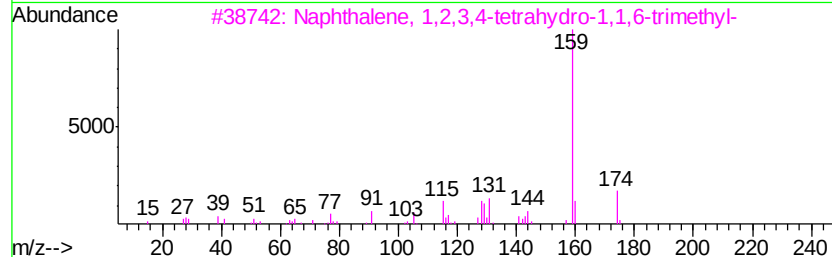
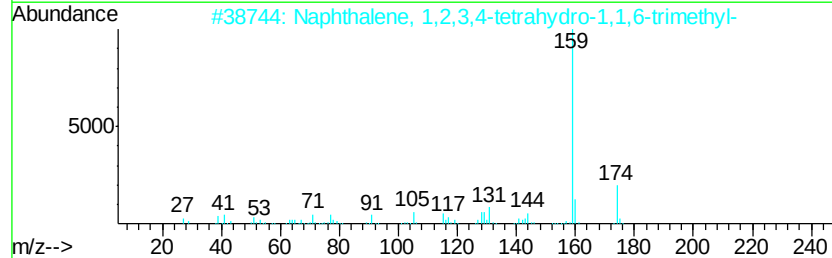
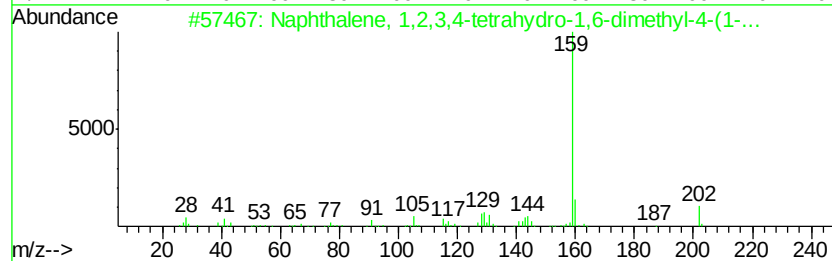
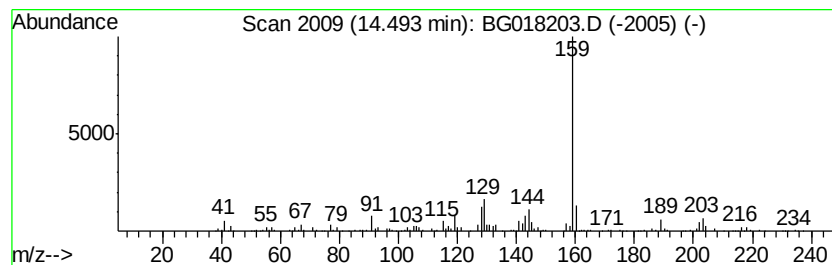
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Quant Title : SVOA CALIBRATION

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

Peak Number 32 Naphthalene, 1,2,3,4-tetra... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.49	9.68 ng/ul	375992	Acenaphthene-d10	14.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	202	C15H22	000483-77-2	87
2			Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	000475-03-6	72
3			Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	000475-03-6	72
4			1H-Indene, 2,3-dihydro-1,1,4,7-t...	174	C13H18	001078-04-2	59
5			4-Methoxy-2-quinolinecarboxylic ...	203	C11H9NO3	015733-83-2	59



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C02A6

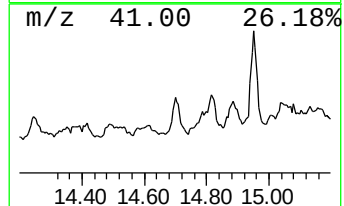
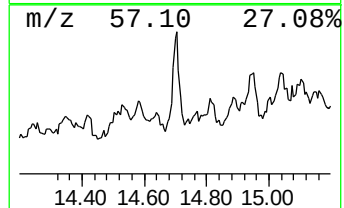
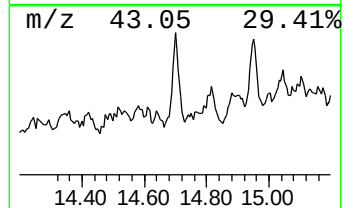
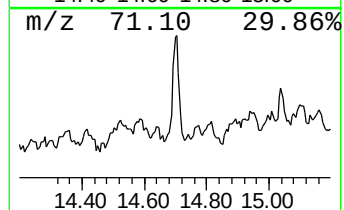
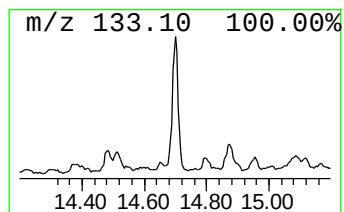
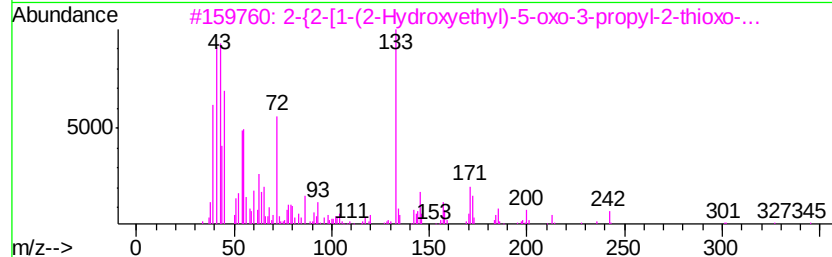
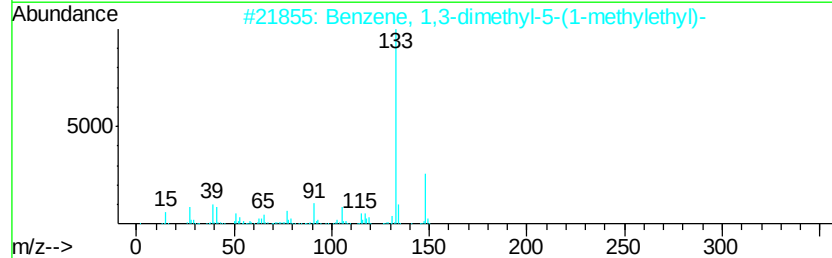
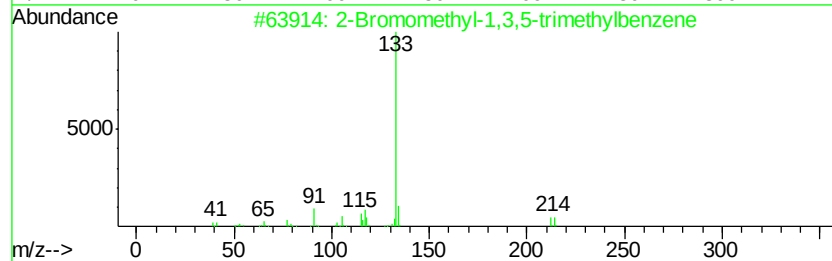
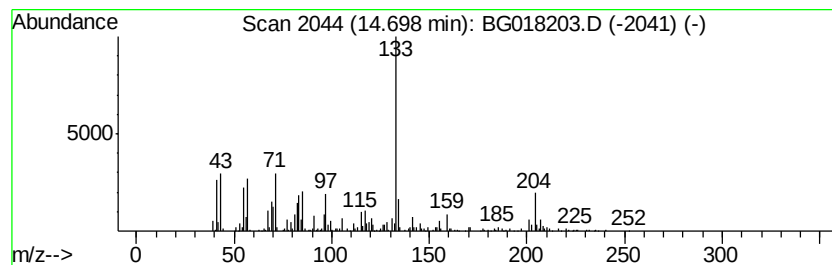
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Quant Title : SVOA CALIBRATION

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

Peak Number 33 unknown-13 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.70	14.06 ng/ul	546262	Acenaphthene-d10	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Bromomethyl-1,3,5-trimethylben...	212	C10H13Br	004761-00-6	46
2		Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	35
3		2-{2-[1-(2-Hydroxyethyl)-5-oxo-3...	417	C20H23N3O5S	1000224-26-4	35
4		1-Propanone, 2-chloro-1-(4-ethyl...	210	C12H15ClO	055012-69-6	35
5		Benzene, 1-(chloromethyl)-2,4,5-...	168	C10H13Cl	010340-77-9	35



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C02A6

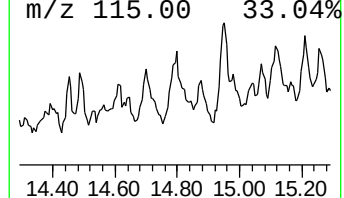
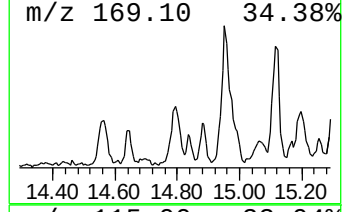
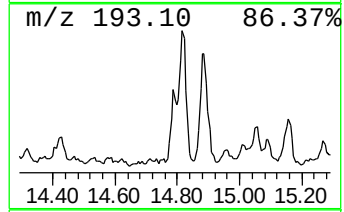
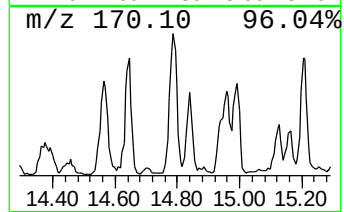
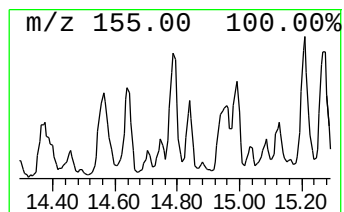
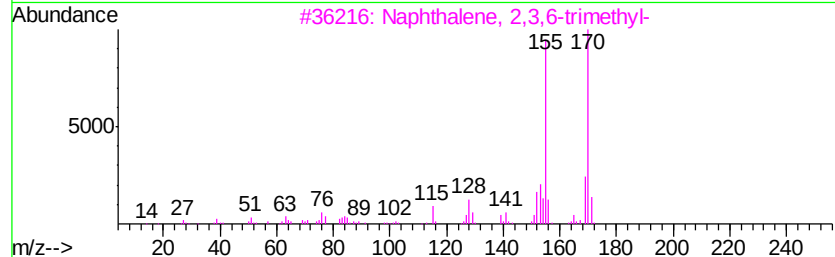
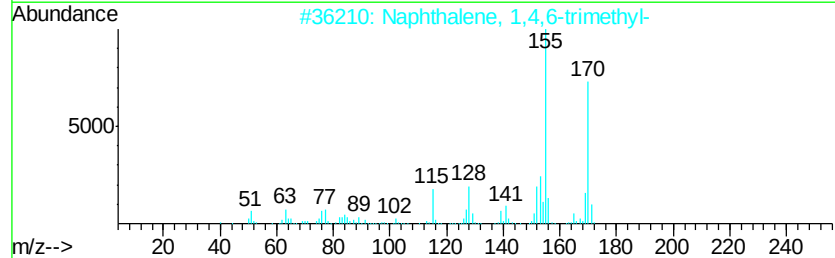
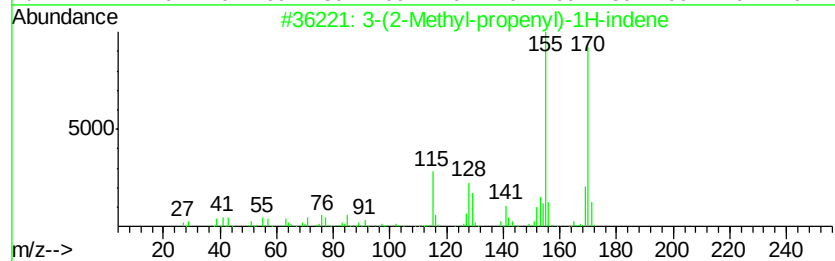
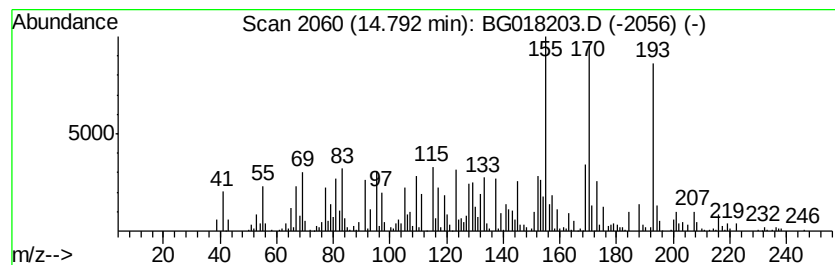
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Quant Title : SVOA CALIBRATION

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

Peak Number 34 3-(2-Methyl-propenyl)-1H-in... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.79	11.50 ng/ul	446677	Acenaphthene-d10	14.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-(2-Methyl-propenyl)-1H-indene	170	C13H14	1000187-78-5	91
2			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	91
3			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	90
4			Azulene, 4,6,8-trimethyl-	170	C13H14	000941-81-1	70
5			1-(2,2-Dimethylcyclopropyl)-2-ph...	170	C13H14	1000294-91-1	70



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C02A6

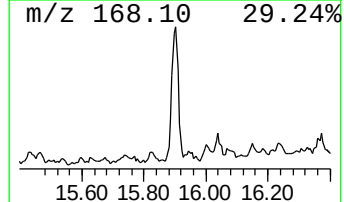
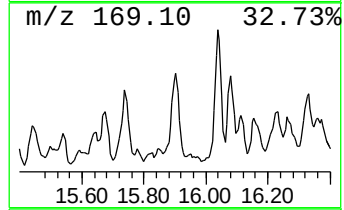
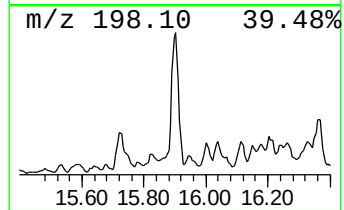
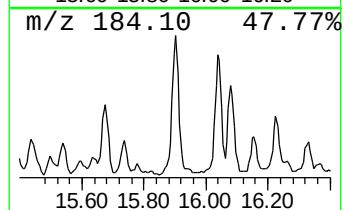
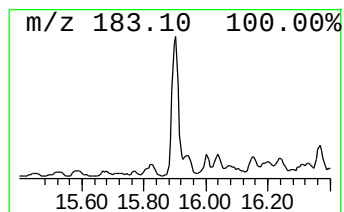
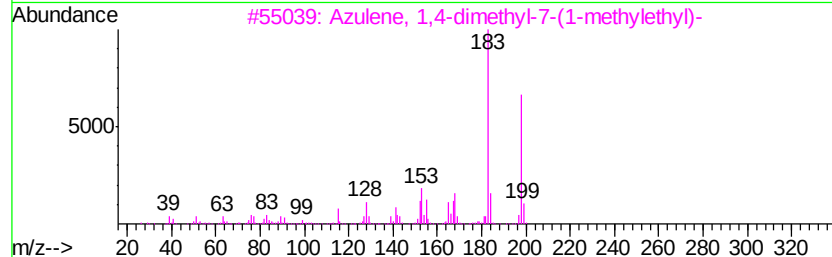
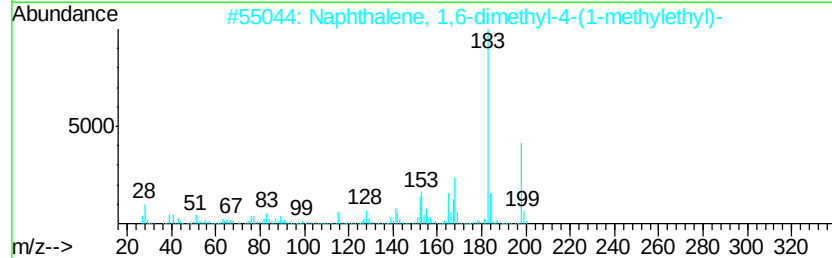
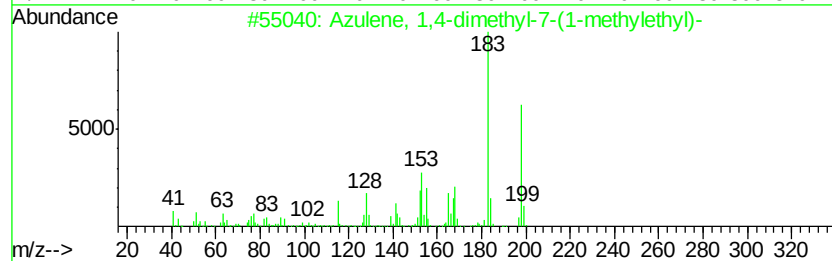
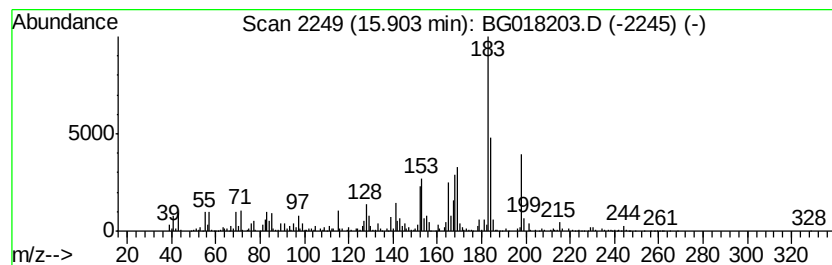
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Quant Title : SVOA CALIBRATION

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

Peak Number 37 Azulene, 1,4-dimethyl-7-(1-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.90	13.46 ng/ul	1161480	Phenanthrene-d10	16.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Azulene, 1,4-dimethyl-7-(1-methyl-...	198	C15H18	000489-84-9	95
2		Naphthalene, 1,6-dimethyl-4-(1-methyl-...	198	C15H18	000483-78-3	92
3		Azulene, 1,4-dimethyl-7-(1-methyl-...	198	C15H18	000489-84-9	89
4		Naphthalene, 1,6-dimethyl-4-(1-methyl-...	198	C15H18	000483-78-3	87
5		Azulene, 1,4-dimethyl-7-(1-methyl-...	198	C15H18	000489-84-9	70



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C02A6

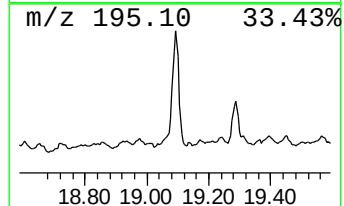
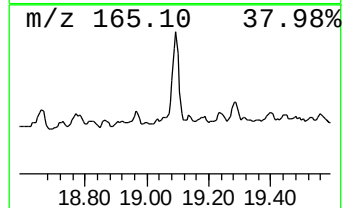
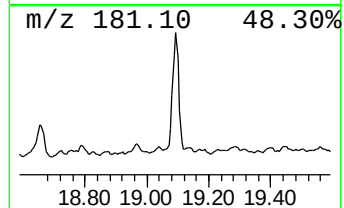
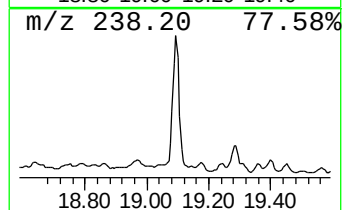
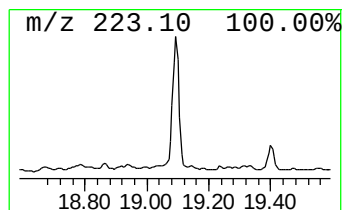
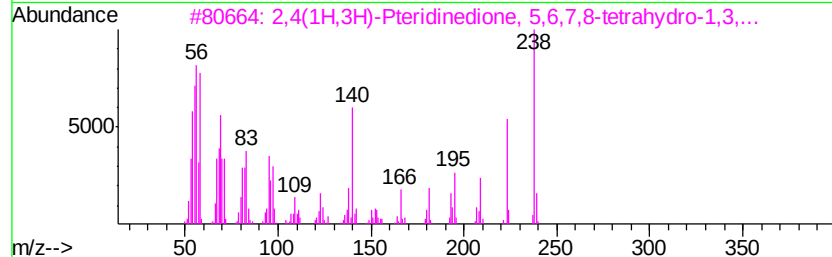
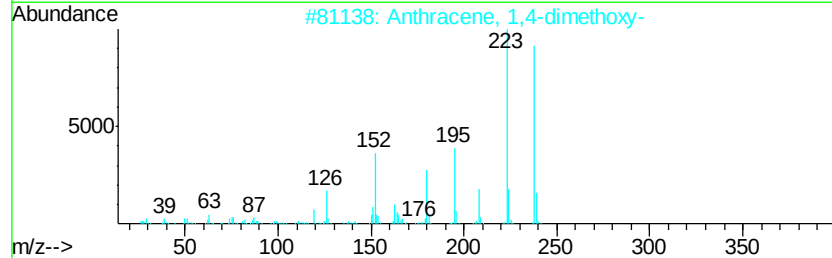
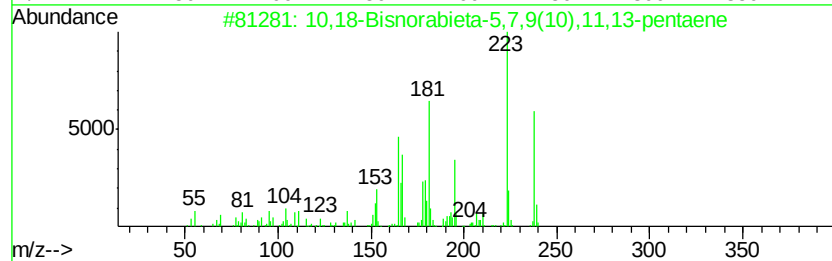
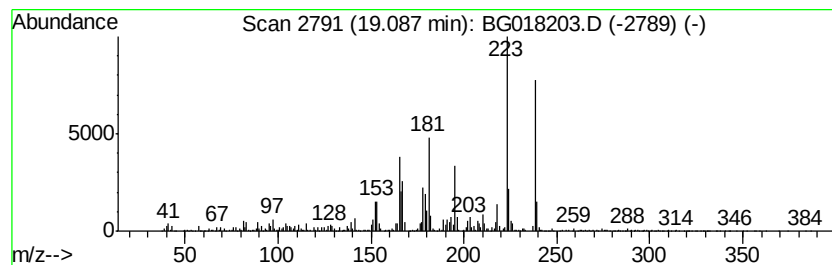
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Quant Title : SVOA CALIBRATION

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

Peak Number 38 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.09	20.00 ng/ul	6533550	Chrysene-d12	21.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99
2		Anthracene, 1,4-dimethoxy-	238	C16H14O2	013076-29-4	60
3		2,4(1H,3H)-Pteridinedione, 5,6,7...	238	C11H18N4O2	037921-31-6	56
4		4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	53
5		Cyclopent[a]indene, 3,8-dihydro-...	238	C18H22	017384-72-4	50



Data Path : Z:\HPCHEM1\BNA G\Data\BG073115\
Data File : BG018203.D
Acq On : 1 Aug 2015 4:38
Operator : TP/UM
Sample : G3065-12
Misc :
ALS Vial : 23 Sample Multiplier: 1

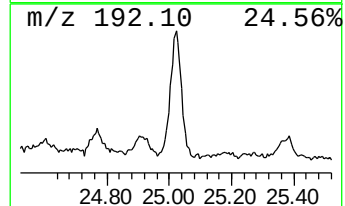
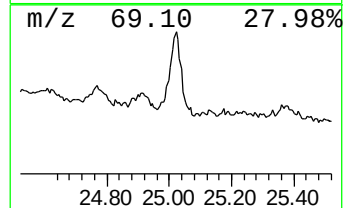
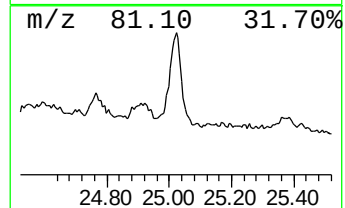
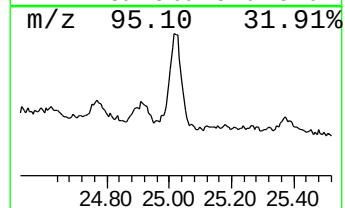
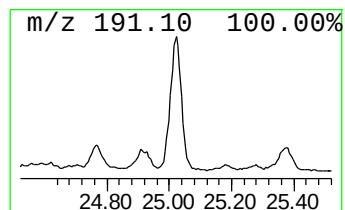
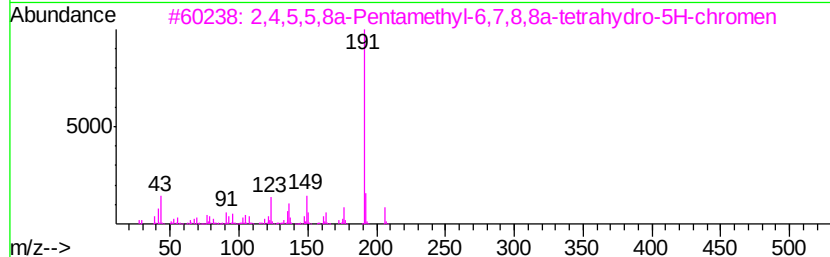
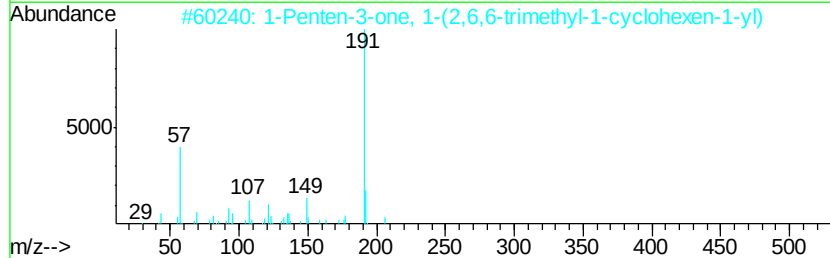
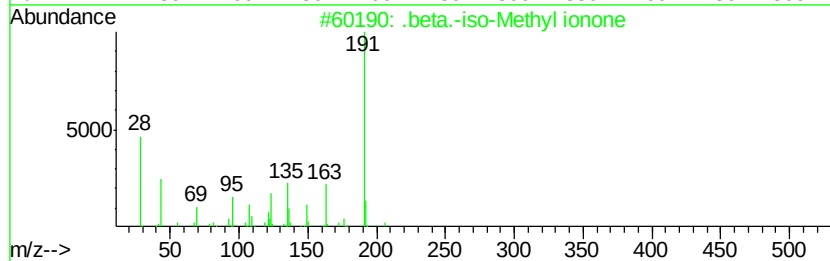
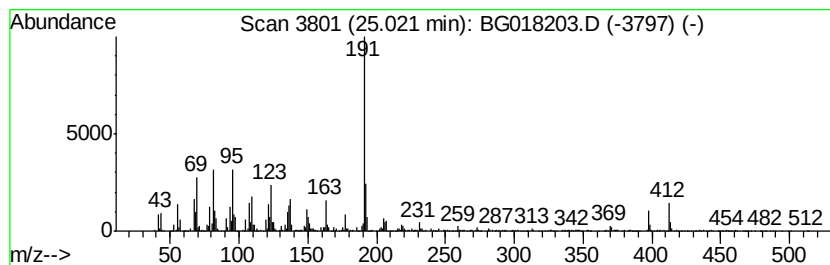
Instrument :
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG073115.M
Quant Title : SVOA CALIBRATION

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

Peak Number 39 .beta.-iso-Methyl ionone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.02	20.00 ng/ul	3320430	Perylene-d12	23.36
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		.beta.-iso-Methyl ionone	206 C14H22O	1000285-40-2 76
2		1-Penten-3-one, 1-(2,6,6-trimeth...	206 C14H22O	000127-43-5 58
3		2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206 C14H22O	1000195-40-9 50
4		5-(p-Aminophenyl)-2-thiazolamine	191 C9H9N3S	090349-87-4 43
5		4H-1,2,4-Triazole, 3-ethyl-5-(3,...	191 C9H13N5	328532-29-2 43



Data Path : Z:\HPCHEM1\BNA_G\Data\BG073115\
 Data File : BG018203.D
 Acq On : 1 Aug 2015 4:38
 Operator : TP/UM
 Sample : G3065-12
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C02A6

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG073115.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Ethyl-1-dodecanol	6.48	2.1	ng/ul	38973	1	7.48	371179	20.0
(DEL) Alkane: Cyc...	6.62	2.0	ng/ul	37838	1	7.48	371179	20.0
(DEL) Alkane: Str...	7.11	4.0	ng/ul	74642	1	7.48	371179	20.0
(DEL) Alkane: Cyc...	7.75	2.7	ng/ul	50651	1	7.48	371179	20.0
unknown-01	8.02	3.6	ng/ul	66098	1	7.48	371179	20.0
Benzene, 2-ethyl-...	8.12	2.0	ng/ul	37548	1	7.48	371179	20.0
1,2,4,4-Tetrameth...	8.48	4.6	ng/ul	86036	1	7.48	371179	20.0
unknown-02	8.58	6.6	ng/ul	121832	1	7.48	371179	20.0
(DEL) Alkane: Str...	8.71	8.0	ng/ul	147852	1	7.48	371179	20.0
unknown-03	8.82	7.3	ng/ul	135821	1	7.48	371179	20.0
Naphthalene, deca...	9.18	2.9	ng/ul	87789	2	10.26	614510	20.0
trans-Decalin, 2-...	9.44	3.1	ng/ul	96689	2	10.26	614510	20.0
Benzene, 1-ethyl-...	9.67	2.0	ng/ul	62003	2	10.26	614510	20.0
(DEL) Alkane: Str...	10.44	2.3	ng/ul	72019	2	10.26	614510	20.0
unknown-04	10.94	2.7	ng/ul	82340	2	10.26	614510	20.0
(DEL) Alkane: Str...	11.30	4.4	ng/ul	134204	2	10.26	614510	20.0
unknown-05	11.35	4.2	ng/ul	128044	2	10.26	614510	20.0
unknown-06	11.55	2.0	ng/ul	61667	2	10.26	614510	20.0
(DEL) Alkane: Str...	11.71	3.1	ng/ul	95523	2	10.26	614510	20.0
unknown-07	12.01	4.1	ng/ul	125507	2	10.26	614510	20.0
2(1H)-Benzocycloo...	12.49	2.8	ng/ul	109743	3	14.16	777032	20.0
cis,trans-2,9-Dim...	12.57	5.7	ng/ul	220716	3	14.16	777032	20.0
unknown-08	12.63	2.6	ng/ul	102282	3	14.16	777032	20.0
unknown-09	12.74	5.6	ng/ul	219283	3	14.16	777032	20.0
Decahydro-4,4,8,9...	12.92	10.9	ng/ul	422613	3	14.16	777032	20.0
unknown-10	13.08	2.5	ng/ul	98508	3	14.16	777032	20.0
Naphthalene, 1,8-...	13.47	6.3	ng/ul	245317	3	14.16	777032	20.0
unknown-11	13.95	5.2	ng/ul	201212	3	14.16	777032	20.0
unknown-12	14.06	8.3	ng/ul	324286	3	14.16	777032	20.0
1H-Inden-1-one, 2...	14.45	2.9	ng/ul	111267	3	14.16	777032	20.0
Naphthalene, 1,2,...	14.49	9.7	ng/ul	375992	3	14.16	777032	20.0
unknown-13	14.70	14.1	ng/ul	546262	3	14.16	777032	20.0
3-(2-Methyl-prope...	14.79	11.5	ng/ul	446677	3	14.16	777032	20.0
Azulene, 1,4-dime...	15.90	13.5	ng/ul	1161480	4	16.92	1726310	20.0
10,18-Bisnorabiet...	19.09	20.0	ng/ul	6533550	5	21.16	6533550	20.0
.beta.-iso-Methyl...	25.02	20.0	ng/ul	3320430	6	23.36	3320430	20.0