

Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
 Data File : BG019517.D
 Acq On : 6 Nov 2015 18:14
 Operator : UM/NP
 Sample : G4245-15
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BCY50

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-BG102815.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	3.128	41	49	58	rBV	73991	152999	1.31%	0.304%
2	3.205	58	62	76	rVB	194882	304067	2.60%	0.604%
3	6.154	559	564	571	rVB	28625	43472	0.37%	0.086%
4	6.642	641	647	657	rVB2	27723	44882	0.38%	0.089%
5	7.323	756	763	769	rBV3	31992	59753	0.51%	0.119%
6	7.494	784	792	798	rBV	120066	200508	1.72%	0.398%
7	7.558	798	803	810	rVB3	27899	46707	0.40%	0.093%
8	7.705	820	828	835	rBV2	125899	206380	1.77%	0.410%
9	7.899	855	861	869	rVB	45827	78536	0.67%	0.156%
10	8.181	902	909	918	rBV	152808	266095	2.28%	0.529%
11	8.293	922	928	937	rVB	55177	91094	0.78%	0.181%
12	8.557	965	973	980	rVB	299705	495078	4.24%	0.983%
13	8.722	995	1001	1009	rVB	159287	267755	2.29%	0.532%
14	8.880	1022	1028	1039	rVB	104073	183203	1.57%	0.364%
15	9.292	1088	1098	1102	rVV3	24877	49331	0.42%	0.098%
16	9.350	1102	1108	1118	rVB2	189160	465787	3.99%	0.925%
17	9.544	1135	1141	1151	rVB	135372	233140	2.00%	0.463%
18	9.668	1151	1162	1169	rVV	169337	396088	3.39%	0.787%
19	9.732	1169	1173	1180	rVV	58963	101910	0.87%	0.202%
20	9.815	1182	1187	1193	rVV	31792	52849	0.45%	0.105%
21	9.879	1193	1198	1204	rVB	25345	42085	0.36%	0.084%
22	10.079	1224	1232	1244	rVB2	162071	289087	2.47%	0.574%
23	10.243	1254	1260	1269	rVB	76266	134812	1.15%	0.268%
24	10.396	1279	1286	1298	rBV3	133540	265862	2.28%	0.528%
25	10.625	1318	1325	1333	rVB	205622	384045	3.29%	0.763%
26	11.013	1384	1391	1393	rVV	198442	362242	3.10%	0.720%
27	11.078	1393	1402	1409	rVV	5730082	11685177	100.00%	23.213%
28	11.184	1412	1420	1427	rVV	1122867	2007327	17.18%	3.988%
29	11.242	1427	1430	1434	rVV3	27285	45574	0.39%	0.091%
30	11.413	1454	1459	1466	rVV4	42739	85674	0.73%	0.170%
31	11.912	1538	1544	1554	rVB2	25158	47201	0.40%	0.094%
32	12.112	1569	1578	1586	rVV2	134241	246046	2.11%	0.489%
33	12.547	1646	1652	1660	rVB	105784	186361	1.59%	0.370%
34	12.764	1684	1689	1693	rVV3	22552	42657	0.37%	0.085%

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35	12.870	1696	1707	1715	rVV	1342361	2301716	19.70%	4.572%
36	13.275	1770	1776	1784	rVB4	24084	47079	0.40%	0.094%
37	13.645	1830	1839	1850	rVB	632662	977175	8.36%	1.941%
38	13.775	1852	1861	1865	rBV4	46136	84594	0.72%	0.168%
39	13.839	1865	1872	1883	rVB	79355	178236	1.53%	0.354%
40	13.939	1883	1889	1890	rBV2	53403	67003	0.57%	0.133%
41	13.980	1890	1896	1904	rVB2	289714	610452	5.22%	1.213%
42	14.133	1914	1922	1928	rBV2	231969	412968	3.53%	0.820%
43	14.204	1928	1934	1940	rVB	491464	683624	5.85%	1.358%
44	14.368	1954	1962	1965	rBV	112644	186407	1.60%	0.370%
45	14.397	1965	1967	1972	rVB	55971	69709	0.60%	0.138%
46	14.503	1977	1985	1995	rBV	477429	873534	7.48%	1.735%
47	14.773	2026	2031	2032	rVV3	33978	45482	0.39%	0.090%
48	14.809	2032	2037	2043	rVV3	383462	650777	5.57%	1.293%
49	14.873	2043	2048	2054	rVV	1514789	2387122	20.43%	4.742%
50	14.938	2054	2059	2063	rVV	159253	251687	2.15%	0.500%
51	14.997	2063	2069	2078	rVV5	56646	124451	1.07%	0.247%
52	15.114	2085	2089	2094	rVV2	47978	84968	0.73%	0.169%
53	15.208	2098	2105	2111	rVV	1444104	2200387	18.83%	4.371%
54	15.273	2111	2116	2125	rVB2	43546	82374	0.70%	0.164%
55	15.408	2130	2139	2145	rBV6	43737	112779	0.97%	0.224%
56	15.473	2145	2150	2158	rVB5	23482	46640	0.40%	0.093%
57	15.631	2172	2177	2181	rVB6	28557	48016	0.41%	0.095%
58	15.796	2199	2205	2210	rVV	654224	1001536	8.57%	1.990%
59	15.855	2210	2215	2220	rVV	759046	1174249	10.05%	2.333%
60	15.907	2220	2224	2228	rVV	243150	362980	3.11%	0.721%
61	15.972	2232	2235	2238	rVV2	58738	92222	0.79%	0.183%
62	16.013	2238	2242	2247	rVV3	76673	144736	1.24%	0.288%
63	16.101	2253	2257	2260	rVV	50528	79392	0.68%	0.158%
64	16.142	2260	2264	2269	rVB	131335	191381	1.64%	0.380%
65	16.248	2273	2282	2284	rBV3	45917	79131	0.68%	0.157%
66	16.283	2284	2288	2297	rVB2	134849	293538	2.51%	0.583%
67	16.518	2323	2328	2336	rVB3	105289	174987	1.50%	0.348%
68	16.718	2356	2362	2371	rVB3	37096	83591	0.72%	0.166%
69	16.824	2371	2380	2388	rBV4	71622	225296	1.93%	0.448%
70	17.000	2405	2410	2415	rVB4	37097	56621	0.48%	0.112%
71	17.288	2452	2459	2464	rBV10	31739	68954	0.59%	0.137%

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72	17.365	2466	2472	2477	rBV	139131	204069	1.75%	0.405%
73	17.429	2477	2483	2487	rBV	105028	161603	1.38%	0.321%
74	17.488	2490	2493	2497	rVB2	47054	54986	0.47%	0.109%
75	17.553	2497	2504	2507	rBV	536047	880761	7.54%	1.750%
76	17.594	2507	2511	2516	rVB	1380250	1970502	16.86%	3.914%
77	17.652	2516	2521	2525	rBV	697678	1013574	8.67%	2.013%
78	17.746	2531	2537	2547	rVB3	41877	87290	0.75%	0.173%
79	17.952	2562	2572	2578	rBV	1213594	1863381	15.95%	3.702%
80	18.052	2583	2589	2594	rVV8	25192	57697	0.49%	0.115%
81	18.240	2617	2621	2627	rVB3	37788	62460	0.53%	0.124%
82	18.340	2634	2638	2642	rBV4	42497	63555	0.54%	0.126%
83	18.422	2648	2652	2655	rBV	51190	68068	0.58%	0.135%
84	18.463	2655	2659	2666	rVB3	197034	303866	2.60%	0.604%
85	18.622	2677	2686	2690	rBV	106973	206105	1.76%	0.409%
86	18.716	2695	2702	2705	rBV	64904	106456	0.91%	0.211%
87	18.757	2705	2709	2714	rVB	73472	99836	0.85%	0.198%
88	18.851	2720	2725	2731	rVB4	89424	131905	1.13%	0.262%
89	19.339	2797	2808	2811	rVB8	28574	71040	0.61%	0.141%
90	19.591	2844	2851	2857	rVB	255871	369684	3.16%	0.734%
91	19.926	2901	2908	2918	rBV2	981651	1668067	14.28%	3.314%
92	20.320	2968	2975	2980	rBV2	73019	146315	1.25%	0.291%
93	20.384	2980	2986	2991	rVV	269288	418938	3.59%	0.832%
94	20.990	3083	3089	3093	rBV	76206	125980	1.08%	0.250%
95	21.037	3093	3097	3105	rVB	100871	162341	1.39%	0.322%
96	21.830	3223	3232	3239	rBV	702682	1234532	10.56%	2.452%
97	22.253	3299	3304	3310	rVB2	22064	47704	0.41%	0.095%
98	22.470	3336	3341	3348	rVB2	49420	96805	0.83%	0.192%
99	24.950	3752	3763	3774	rVB	500830	1423301	12.18%	2.827%
100	25.185	3790	3803	3816	rVB2	368392	1119143	9.58%	2.223%

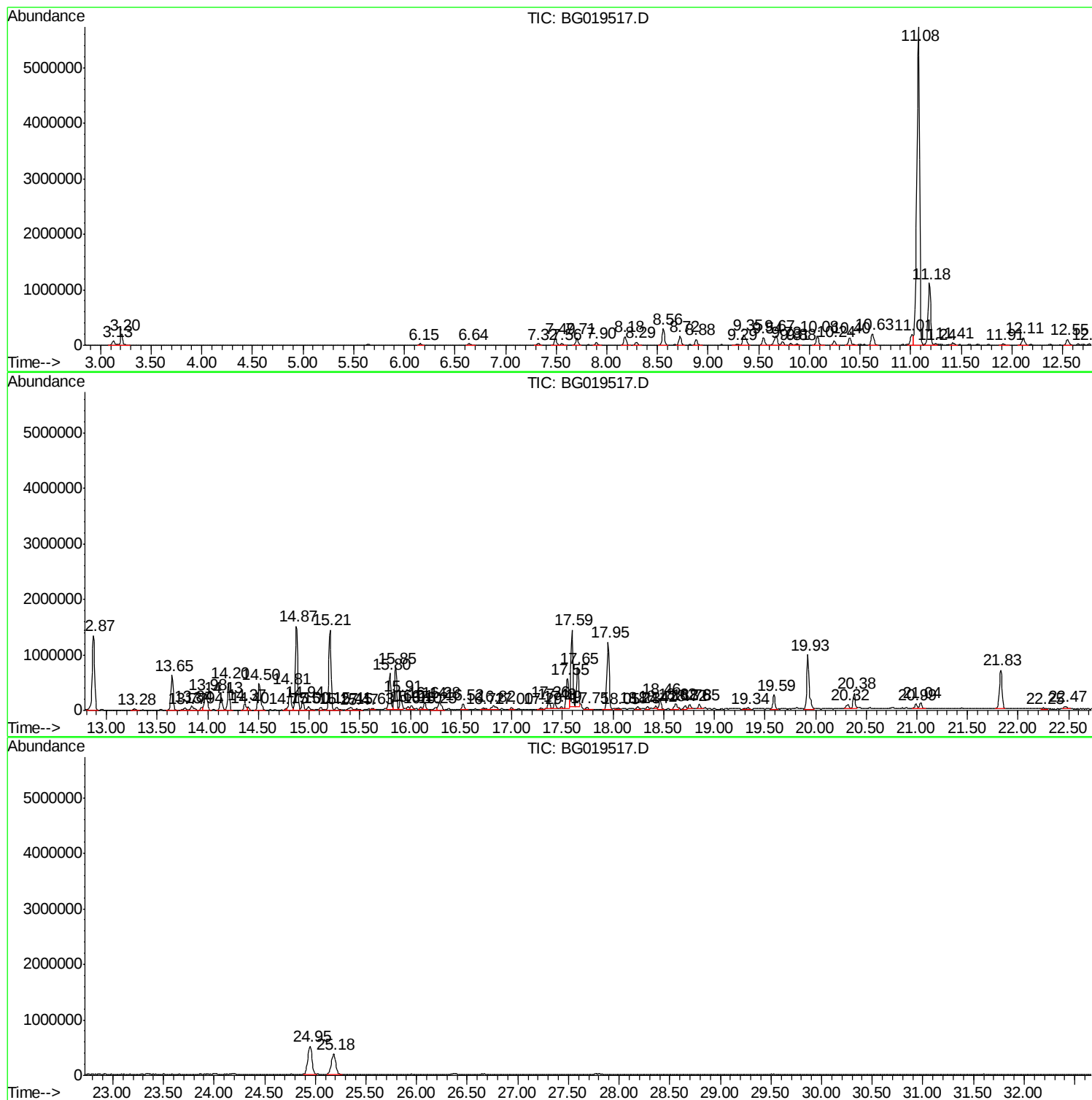
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102815.M
Quant Title : SVOA CALIBRATION

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



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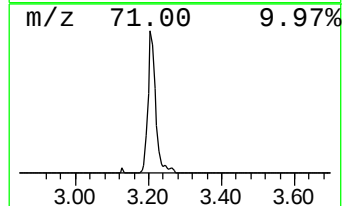
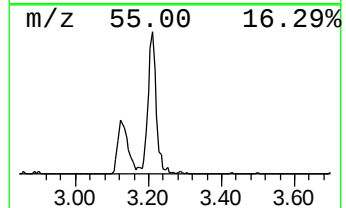
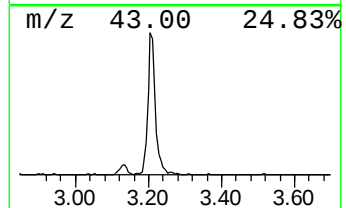
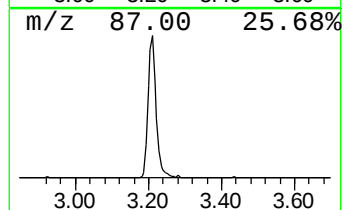
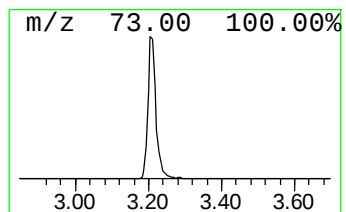
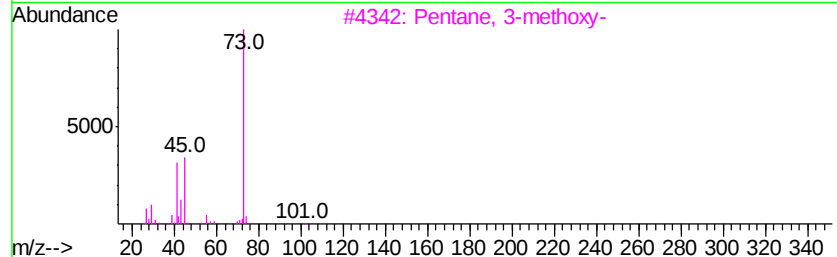
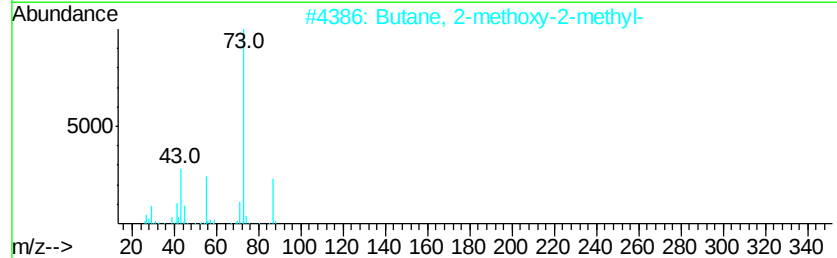
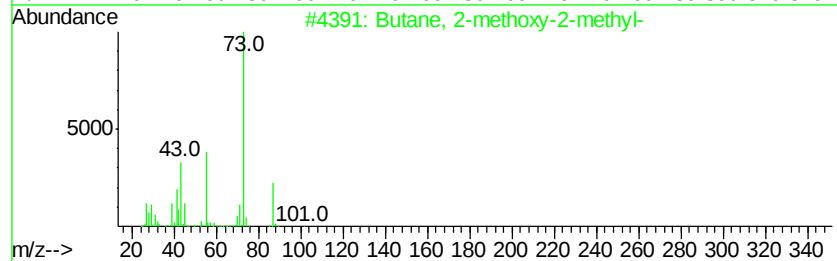
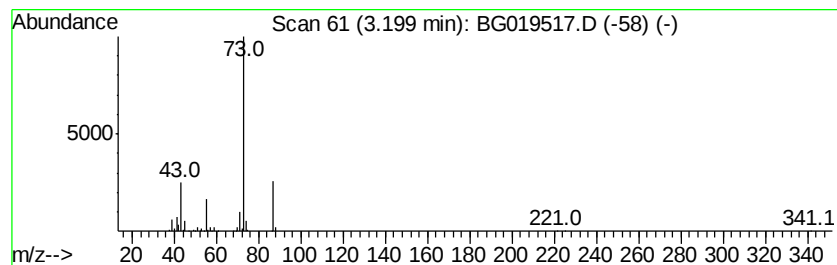
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102815.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.20	22.85 ng/ul	304067	1,4-Dichlorobenzene-d4	8.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	50
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	47
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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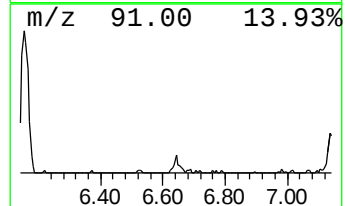
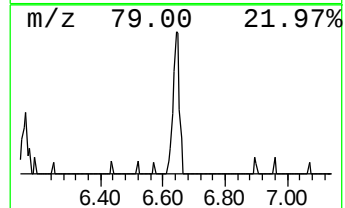
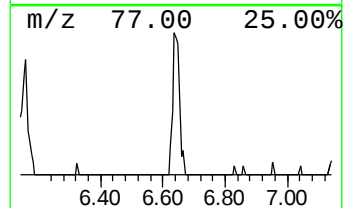
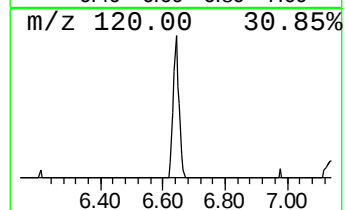
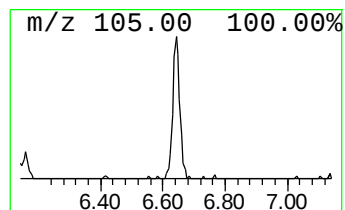
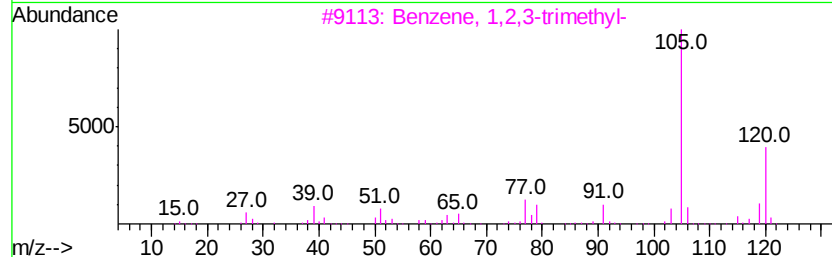
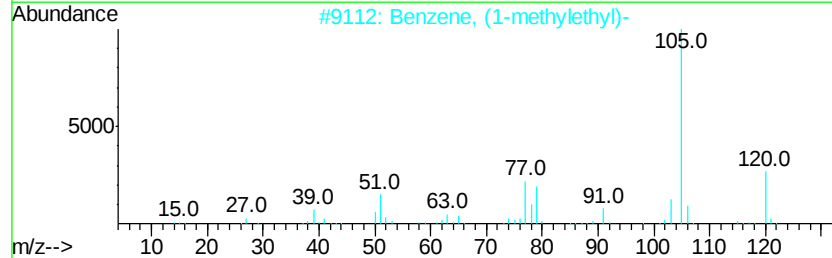
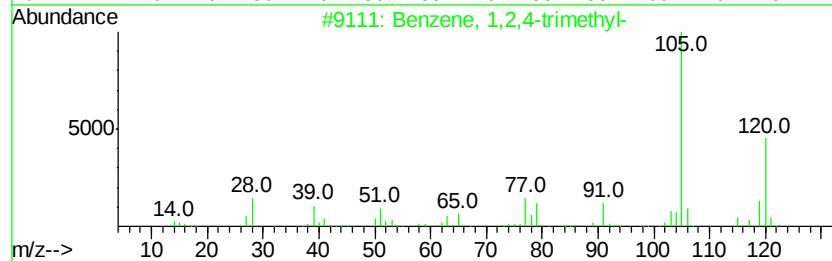
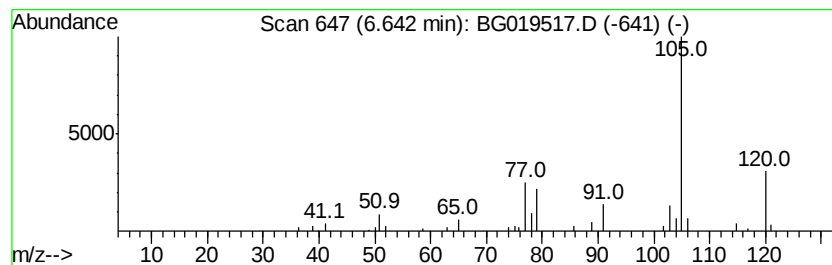
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1,2,4-trimethyl- Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.64	3.37 ng/ul	44882	1,4-Dichlorobenzene-d4	8.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
2		Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	86
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	86
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	80
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	72



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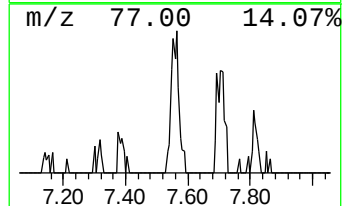
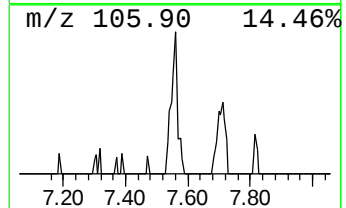
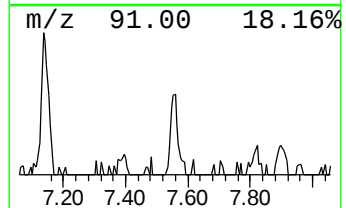
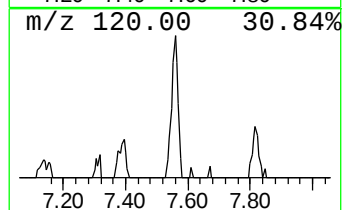
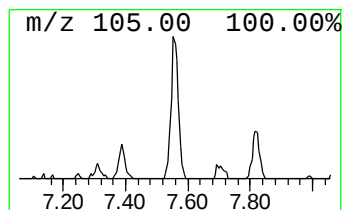
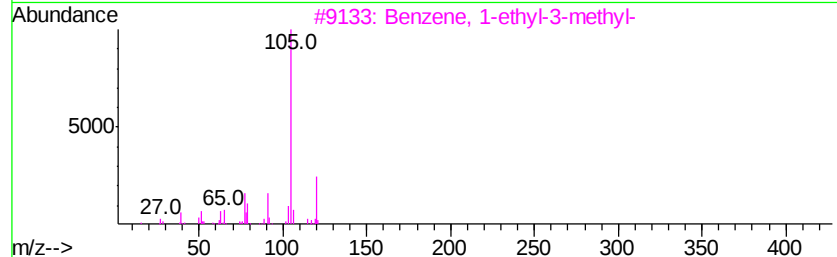
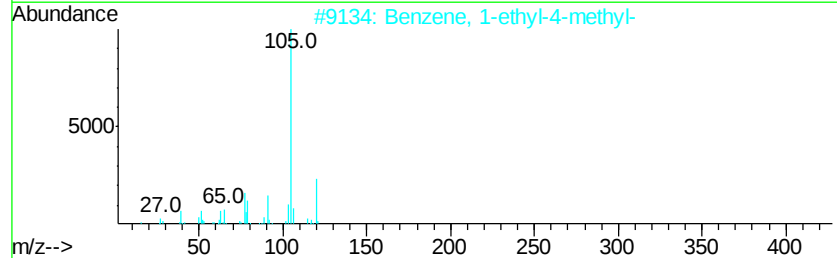
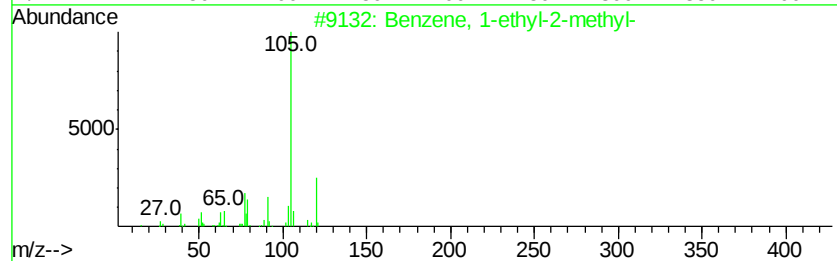
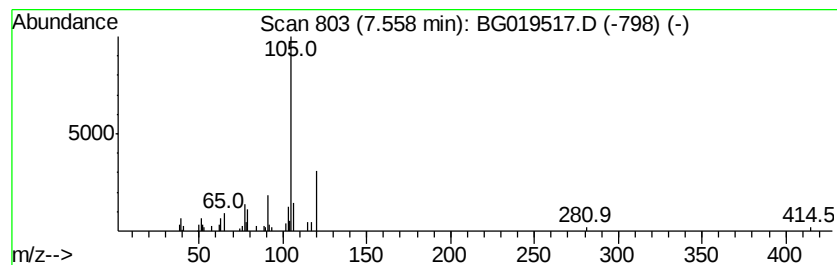
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Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	3.51 ng/ul	46707	1,4-Dichlorobenzene-d4	8.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	87
2		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	83
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	83
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	72
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	72



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
Operator : UM/NP
Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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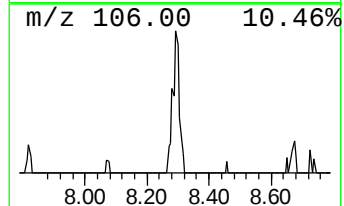
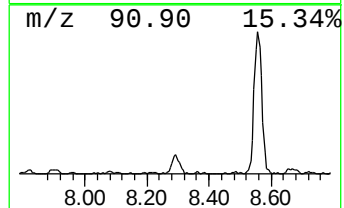
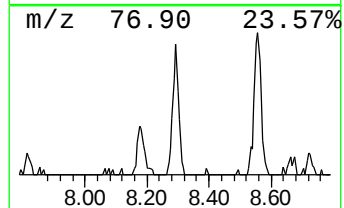
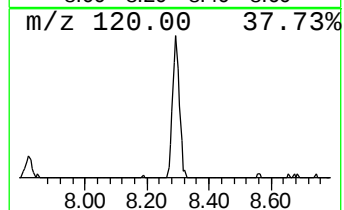
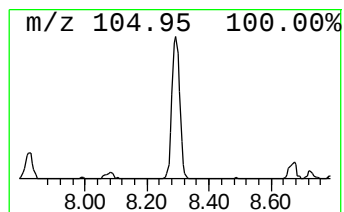
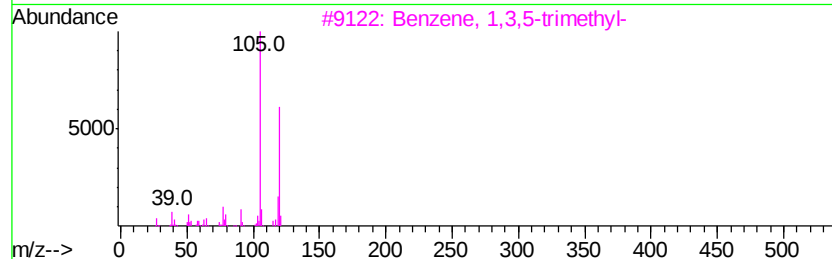
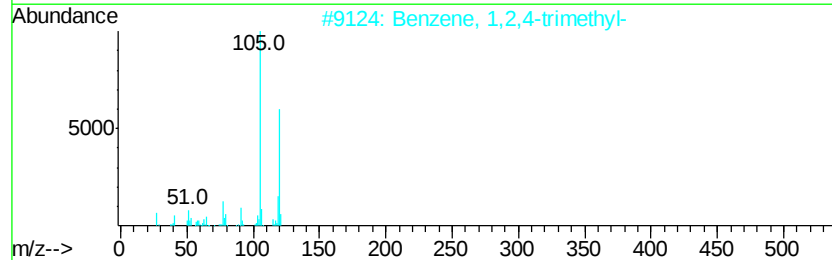
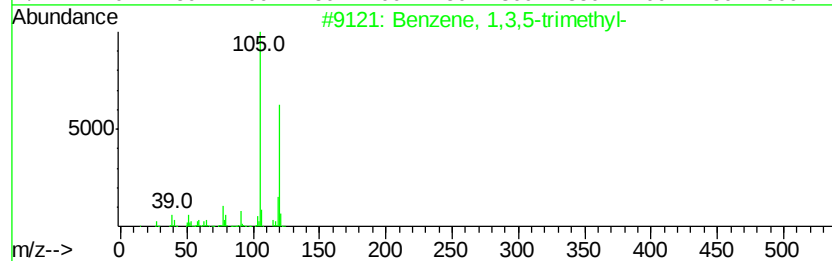
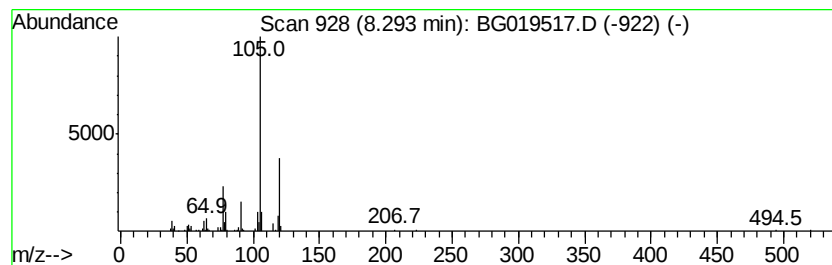
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1,3,5-trimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.29	6.85 ng/ul	91094	1,4-Dichlorobenzene-d4	8.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
3		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	87
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	87



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
Operator : UM/NP
Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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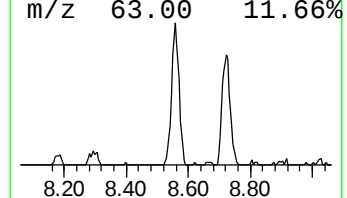
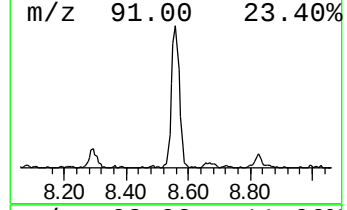
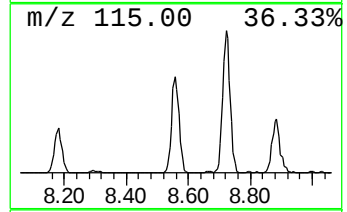
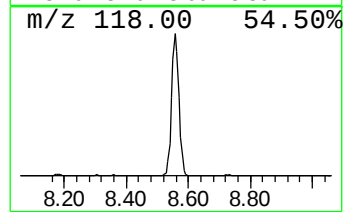
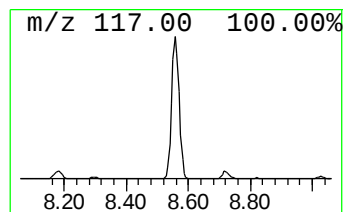
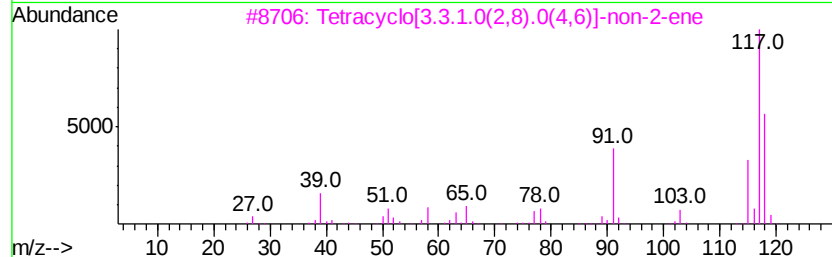
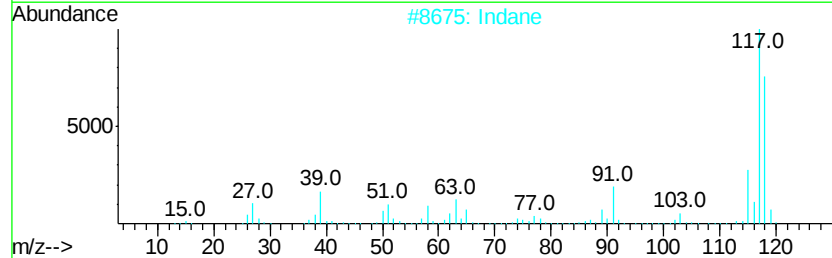
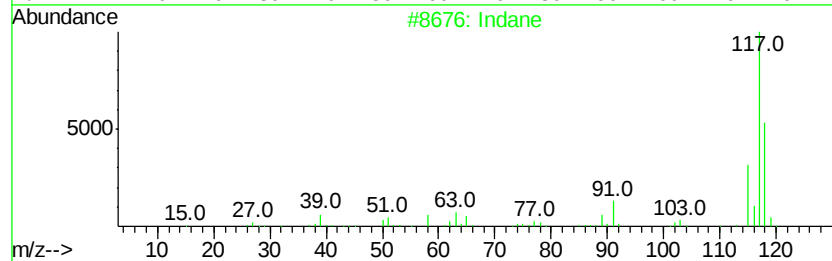
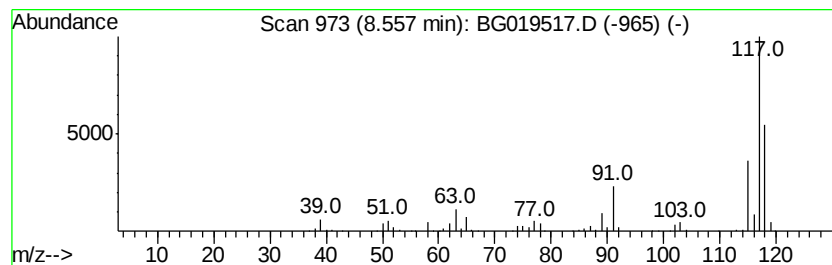
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.56	37.21 ng/ul	495078	1,4-Dichlorobenzene-d4	8.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	94
2	Indane		118	C9H10	000496-11-7	90
3	Tetracyclo[3.3.1.0(2,8).0(4,6)]-	...	118	C9H10	1000191-13-7	76
4	Benzene, cyclopropyl-		118	C9H10	000873-49-4	68
5	Benzene, 2-propenyl-		118	C9H10	000300-57-2	58



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
Operator : UM/NP
Sample : G4245-15
Misc :
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Instrument :
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ClientSampleId :
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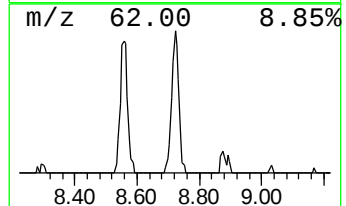
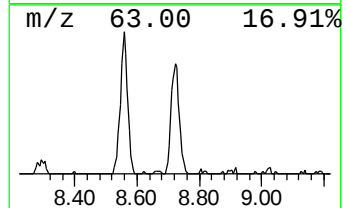
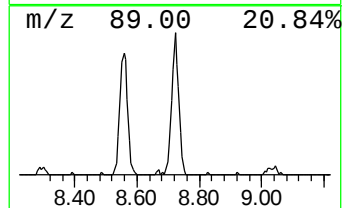
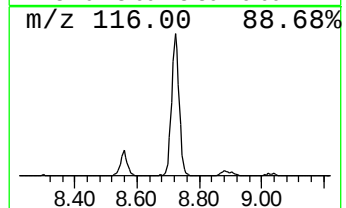
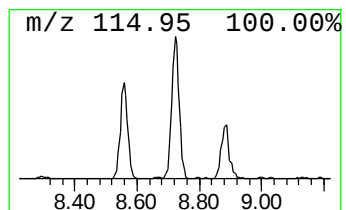
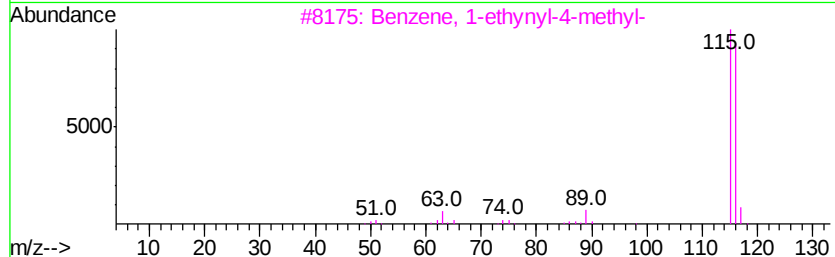
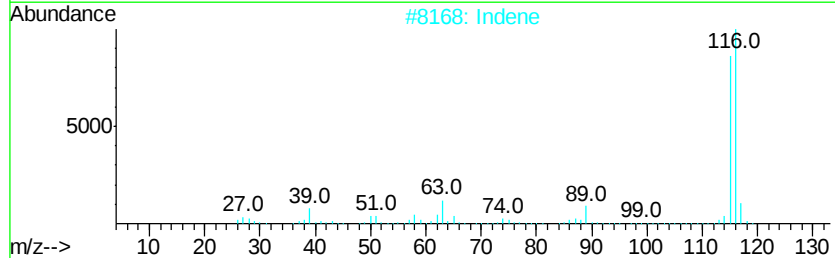
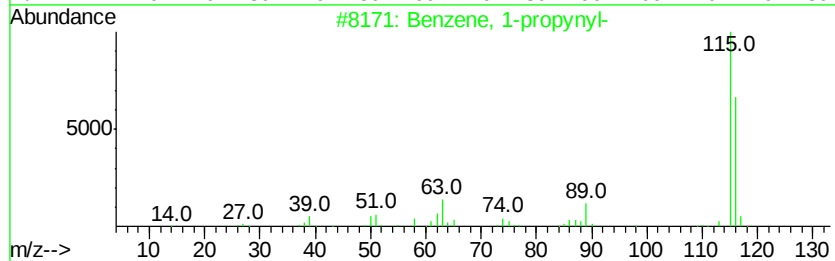
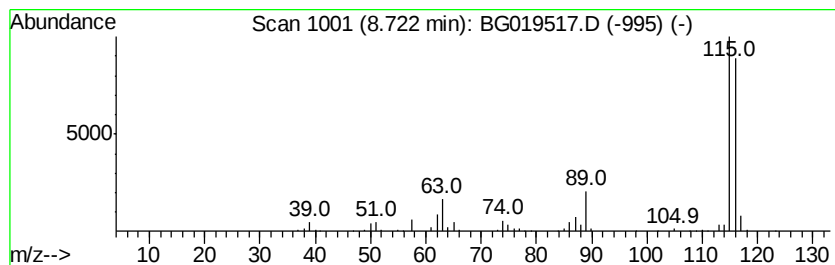
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1-propynyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.72	20.12 ng/ul	267755	1,4-Dichlorobenzene-d4	8.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
2		Indene	116	C9H8	000095-13-6	93
3		Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91
4		Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
5		Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	91



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
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Sample : G4245-15
Misc :
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Instrument :
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ClientSampleId :
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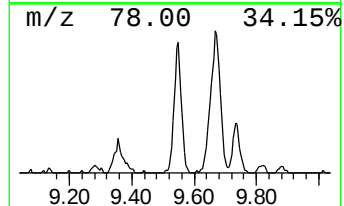
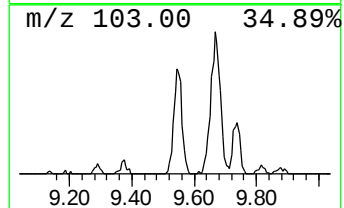
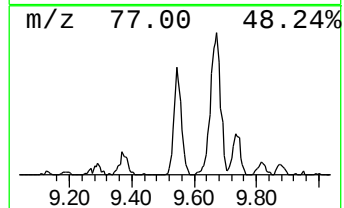
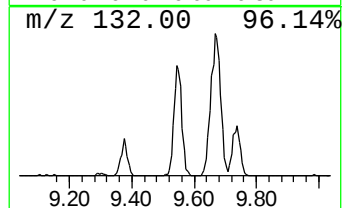
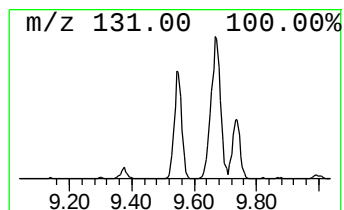
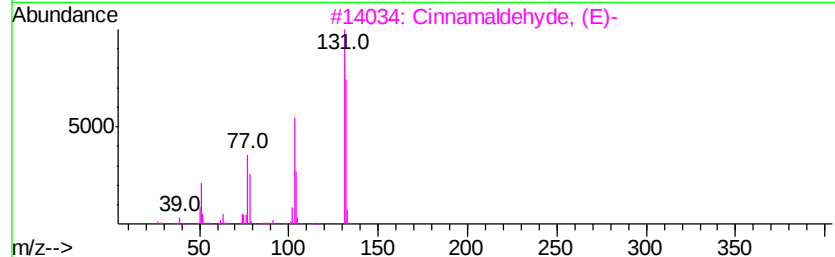
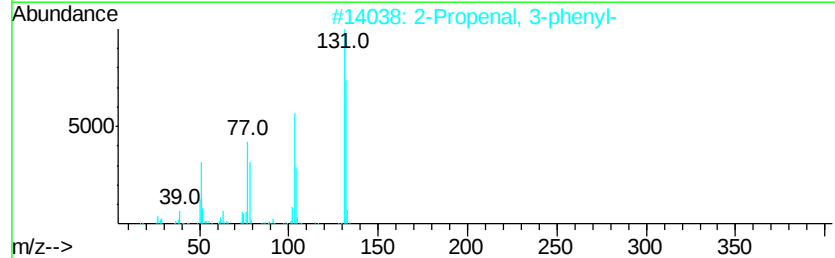
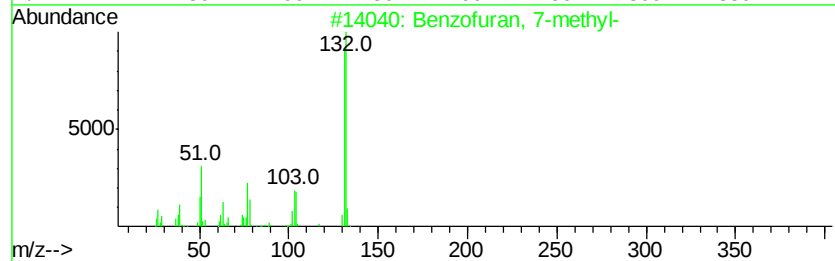
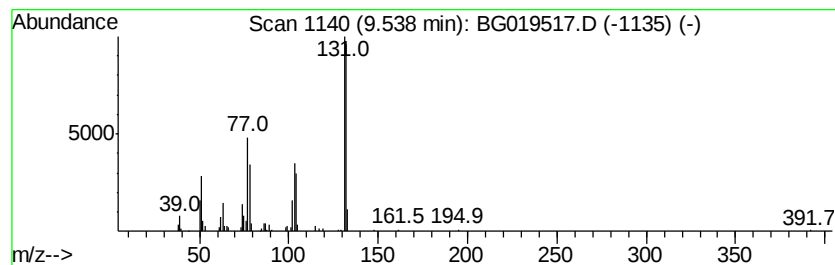
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzofuran, 7-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.54	17.52 ng/ul	233140	1,4-Dichlorobenzene-d4	8.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
2		2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	93
3		Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	93
4		Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	93
5		Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
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Sample : G4245-15
Misc :
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Instrument :
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ClientSampleId :
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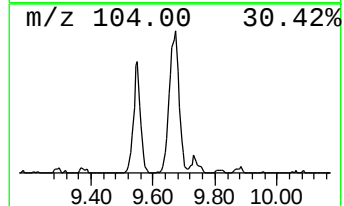
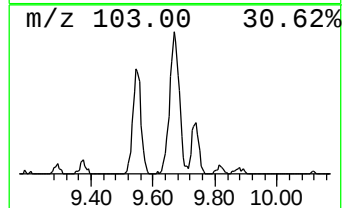
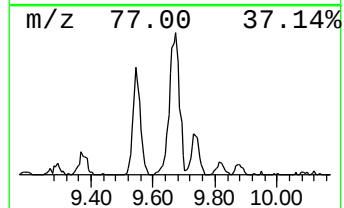
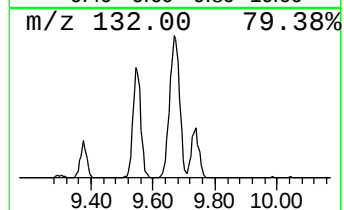
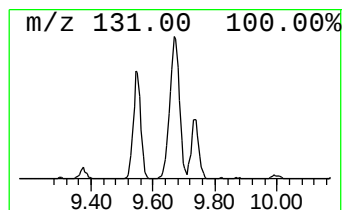
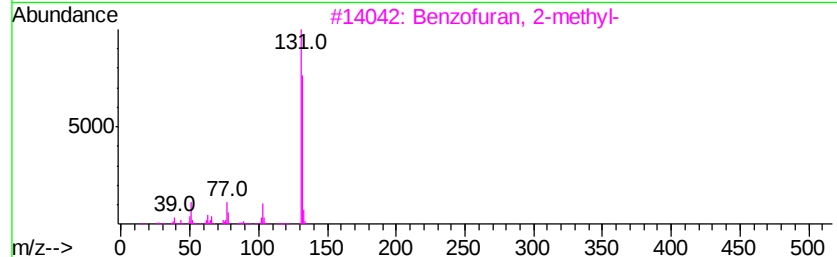
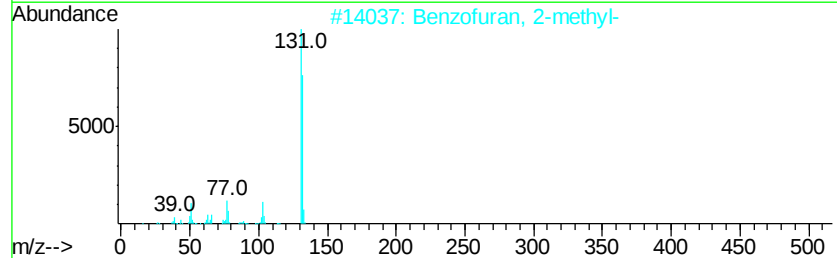
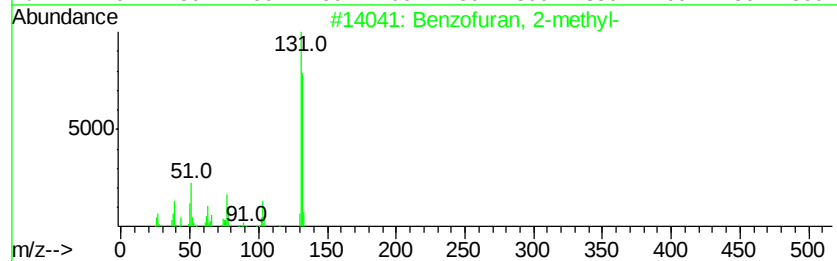
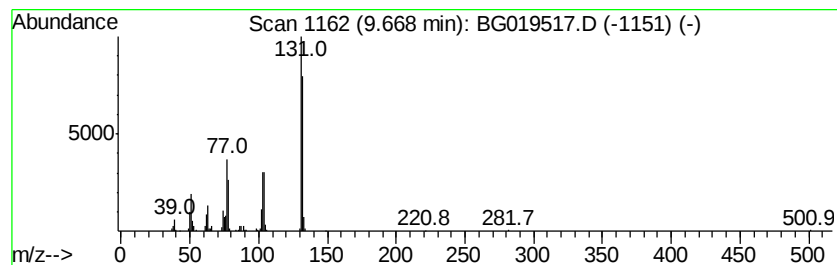
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzofuran, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.67	21.87 ng/ul	396088	Naphthalene-d8	11.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2		Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
3		Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
4		Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	93
5		Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	93



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
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Sample : G4245-15
Misc :
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ClientSampleId :
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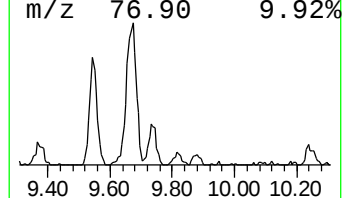
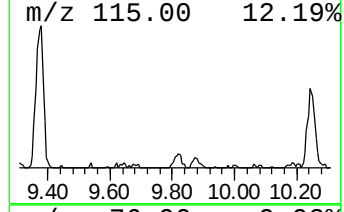
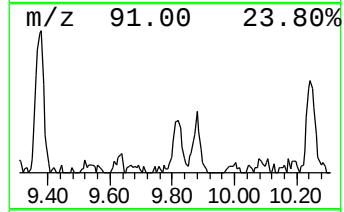
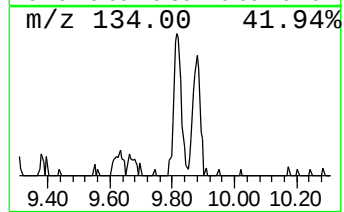
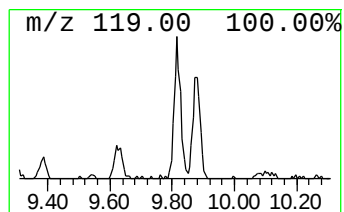
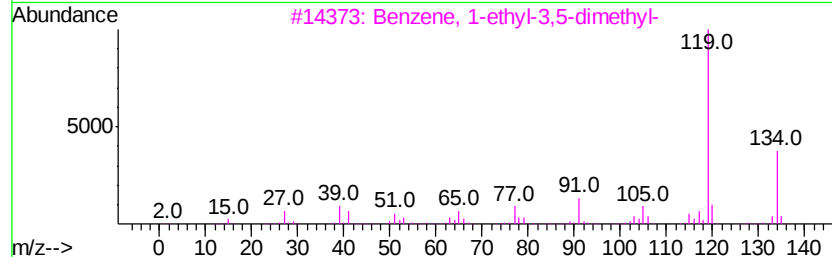
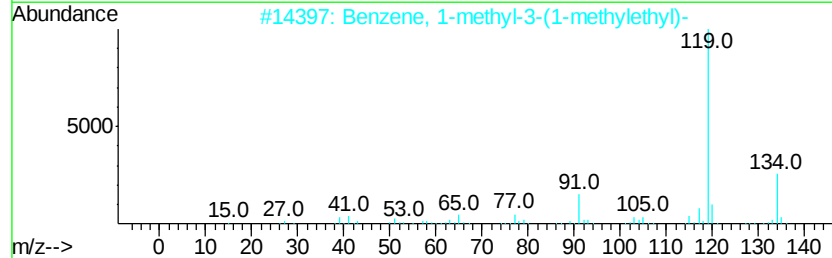
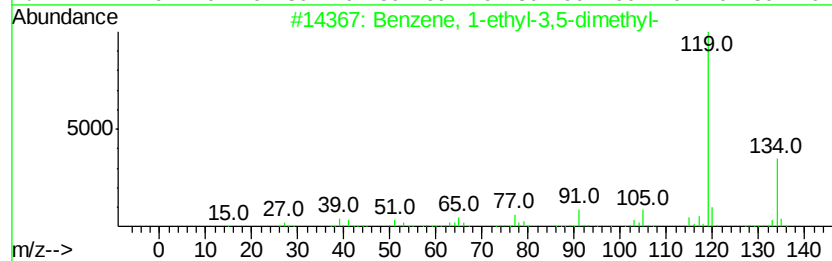
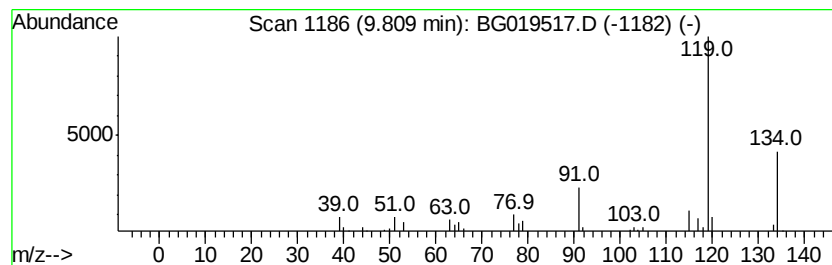
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.81	2.92 ng/ul	52849	Naphthalene-d8	11.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
2		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91
3		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
4		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
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Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

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ClientSampleId :
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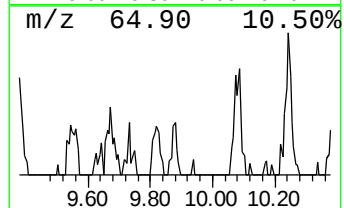
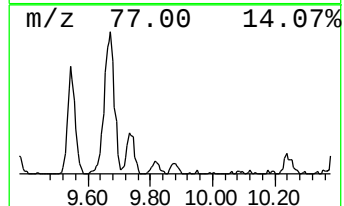
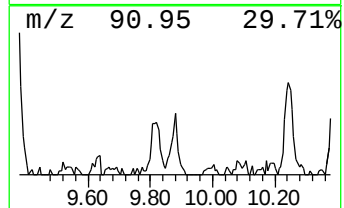
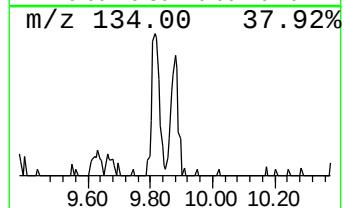
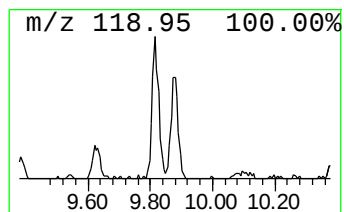
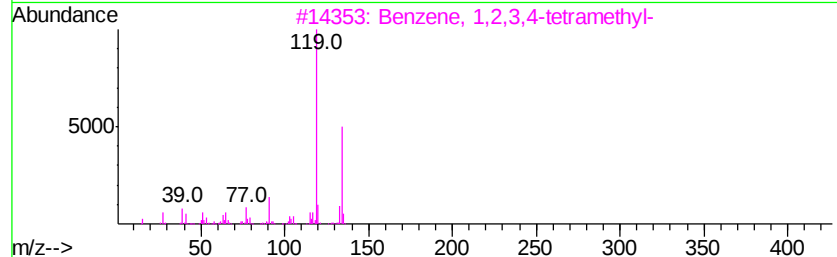
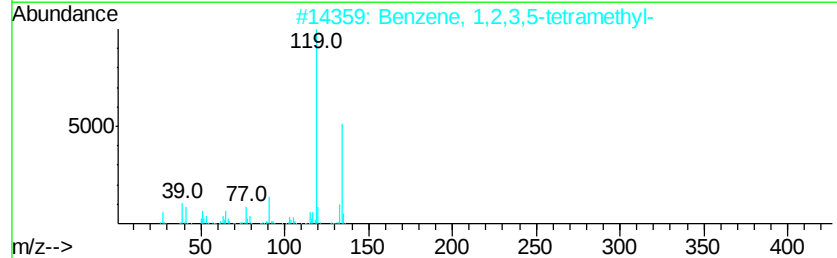
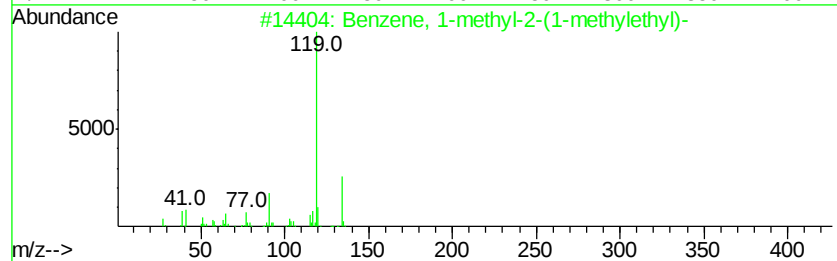
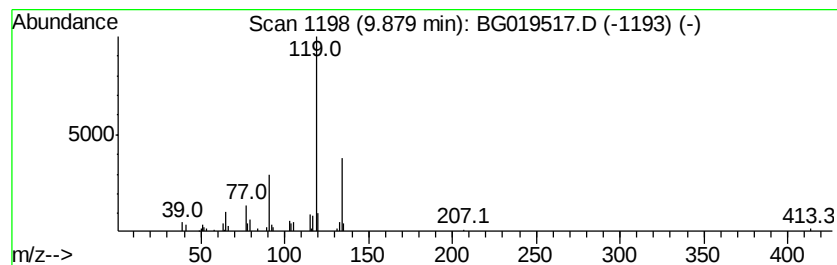
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Benzene, 1-methyl-2-(1-meth... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.88	2.32 ng/ul	42085	Napthalene-d8	11.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	94
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
5		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
Operator : UM/NP
Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BCY50

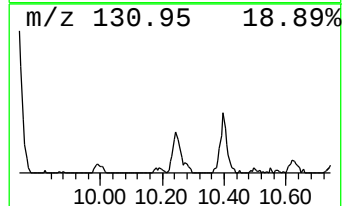
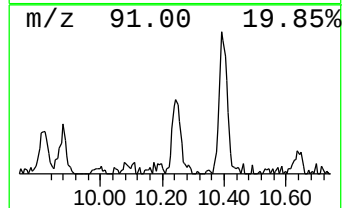
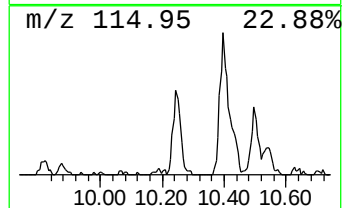
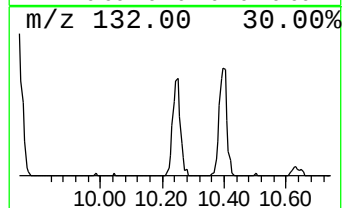
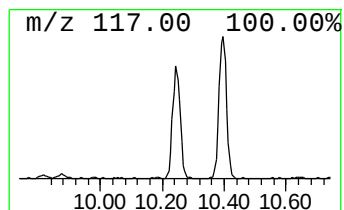
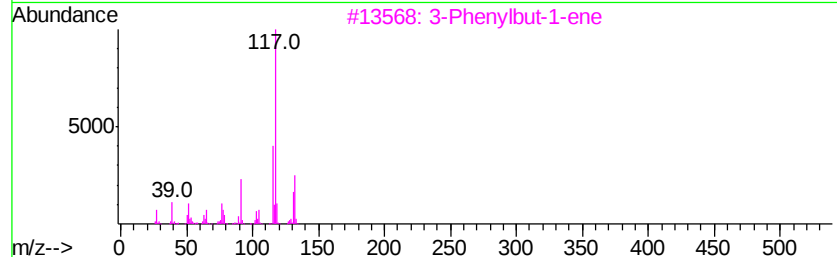
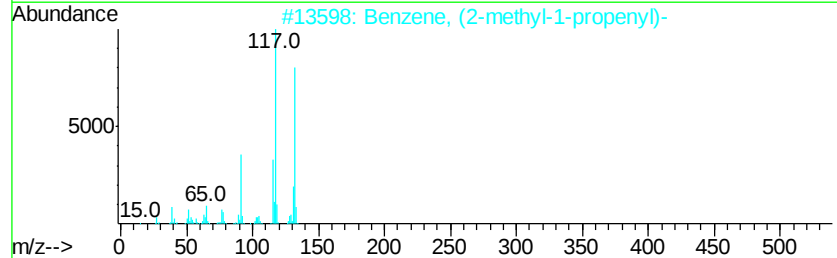
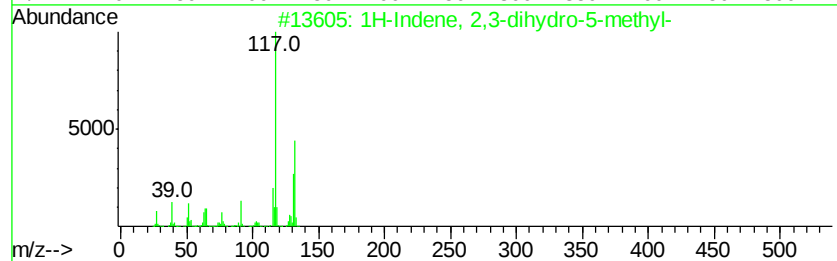
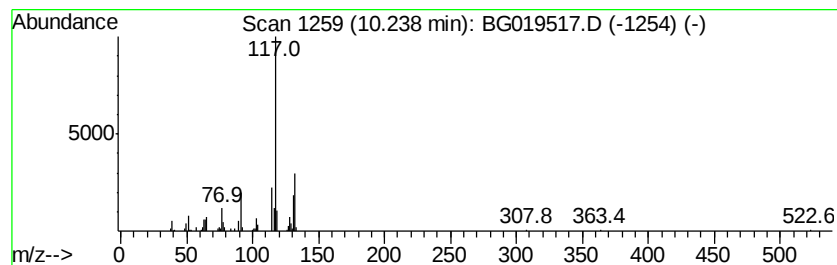
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.24	7.44 ng/ul	134812	Napthalene-d8	11.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
2		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	91
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
4		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	91
5		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	91



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
Operator : UM/NP
Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BCY50

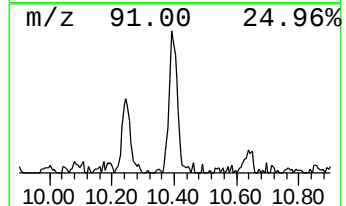
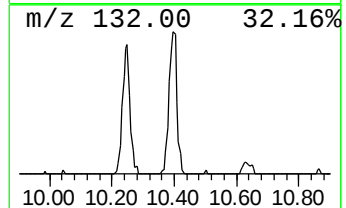
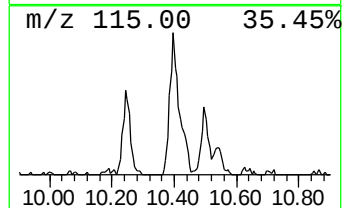
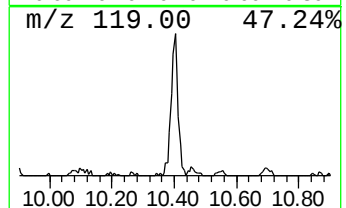
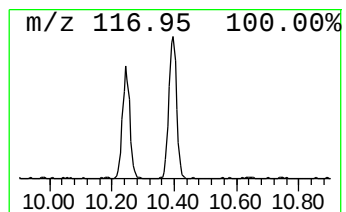
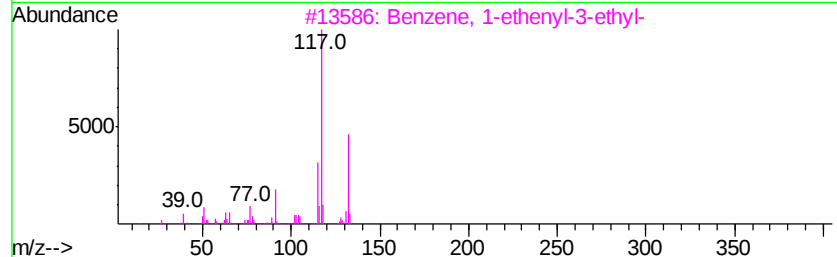
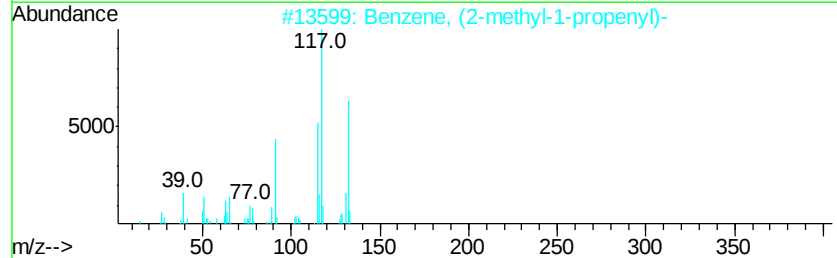
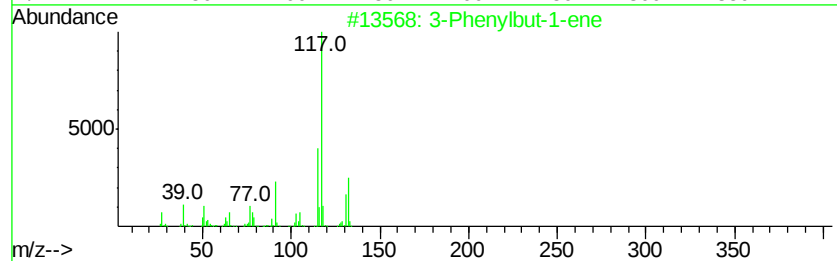
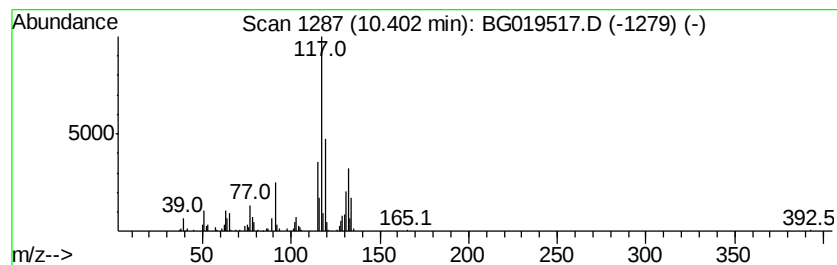
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 3-Phenylbut-1-ene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.40	14.68 ng/ul	265862	Naphthalene-d8	11.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
2		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	76
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	68
4		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	68
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	68



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
Operator : UM/NP
Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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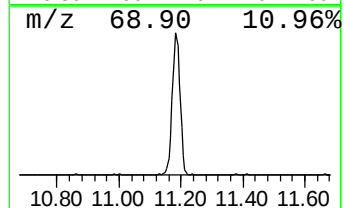
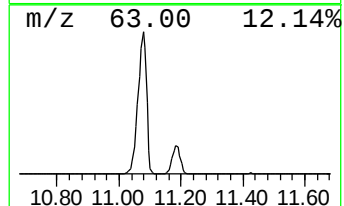
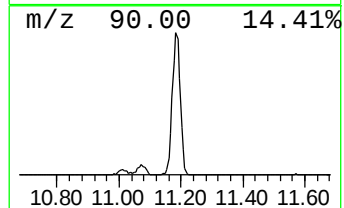
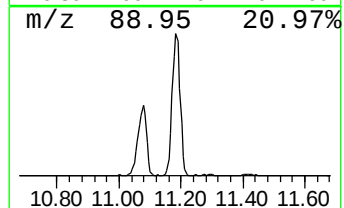
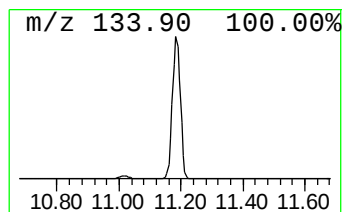
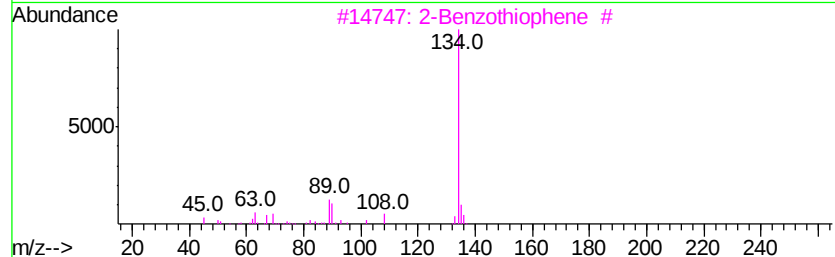
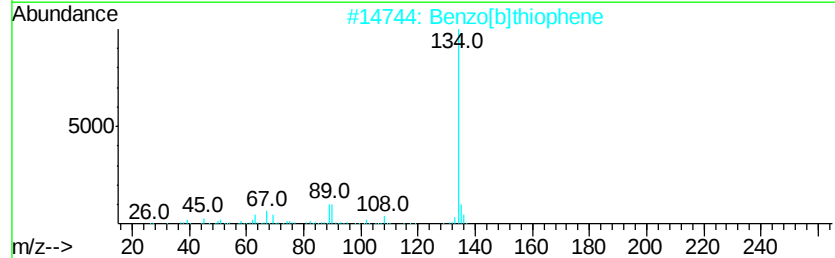
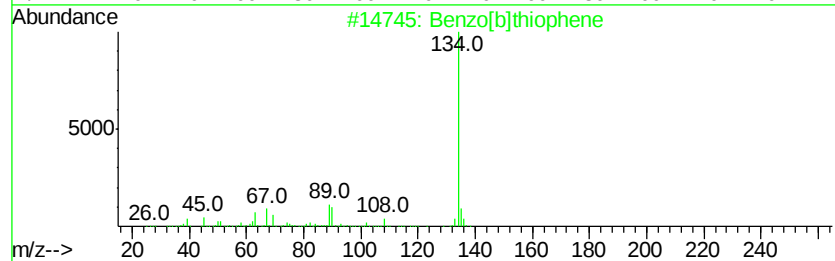
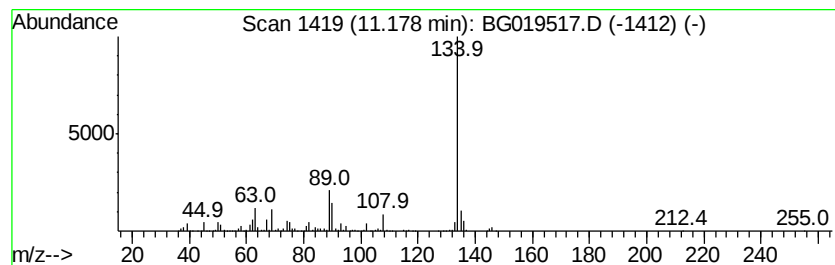
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Benzo[b]thiophene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.18	110.83 ng/ul	2007330	Napthalene-d8	11.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]thiophene	134	C8H6S	000095-15-8	94
2		Benzo[b]thiophene	134	C8H6S	000095-15-8	91
3		2-Benzothiophene #	134	C8H6S	000270-82-6	91
4		Benzo[b]thiophene	134	C8H6S	000095-15-8	90
5		Cyclopenta[b]thiapyran	134	C8H6S	000271-17-0	87



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
Operator : UM/NP
Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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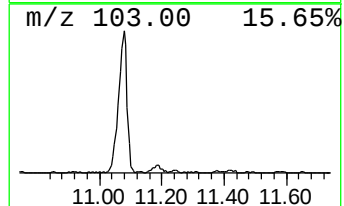
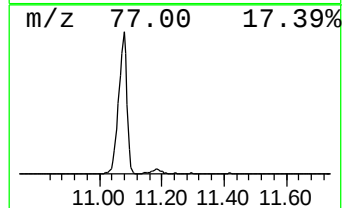
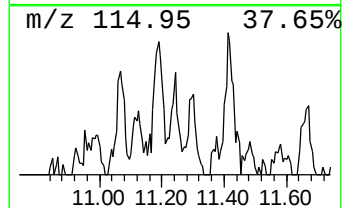
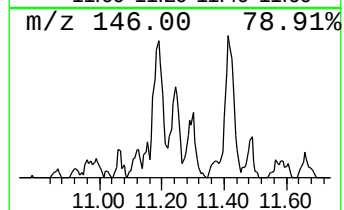
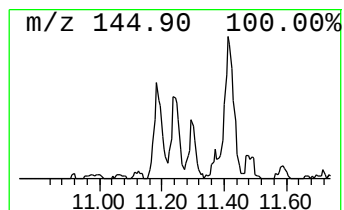
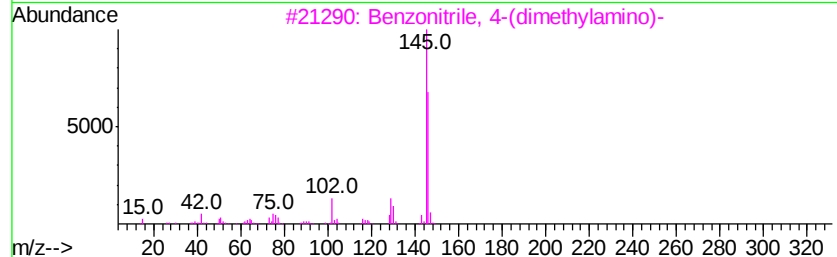
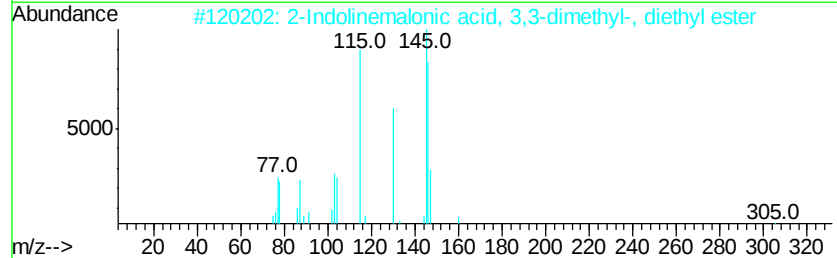
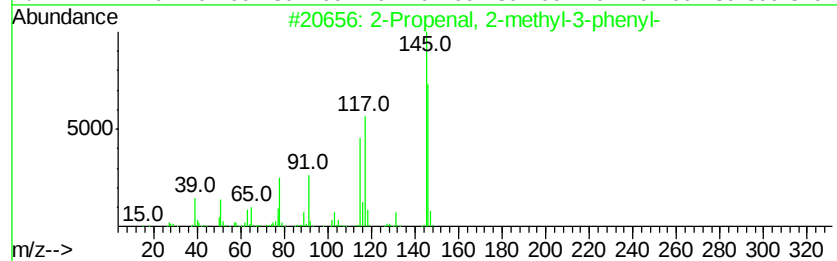
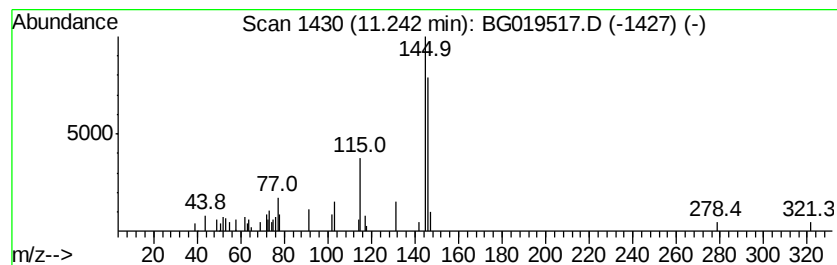
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 2-Propenal, 2-methyl-3-phenyl- Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.24	2.52 ng/ul	45574	Napthalene-d8	11.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propenal, 2-methyl-3-phenyl-	146	C10H10O	000101-39-3	72
2		2-Indolinemalonic acid, 3,3-dime...	305	C17H23NO4	018781-64-1	72
3		Benzonitrile, 4-(dimethylamino)-	146	C9H10N2	001197-19-9	59
4		Benzonitrile, 4-(dimethylamino)-	146	C9H10N2	001197-19-9	53
5		Cinnoline, 1,2-dihydro-2-methyl-	146	C9H10N2	054063-10-4	49



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
Operator : UM/NP
Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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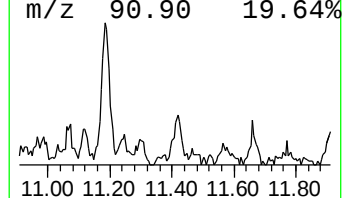
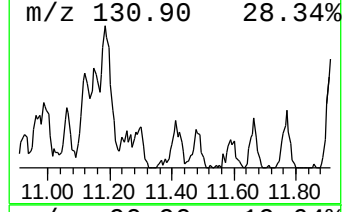
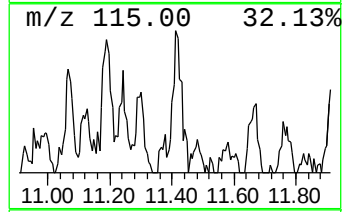
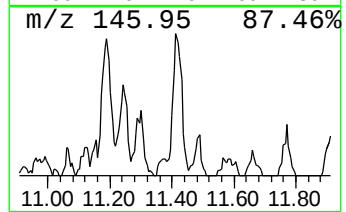
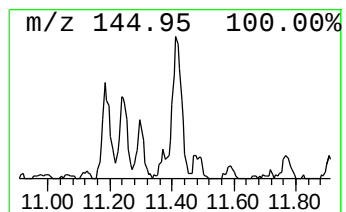
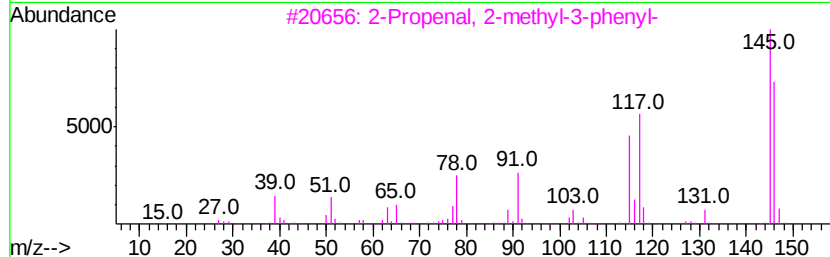
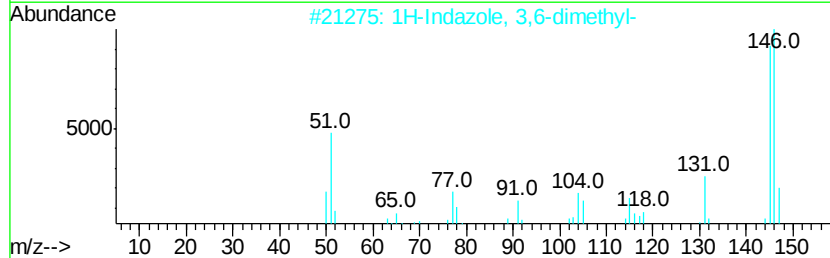
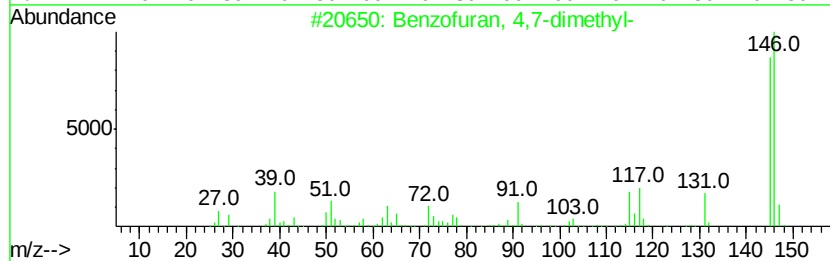
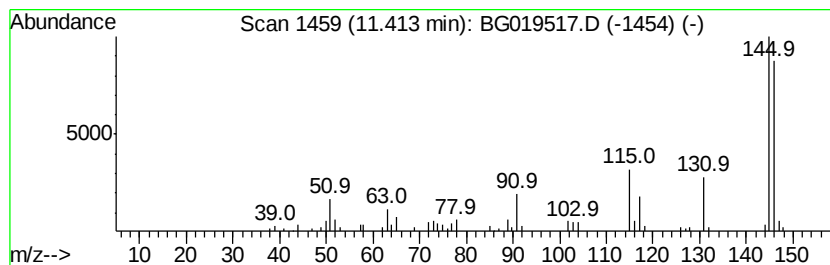
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Benzofuran, 4,7-dimethyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.41	4.73 ng/ul	85674	Napthalene-d8	11.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzofuran, 4,7-dimethyl-	146	C10H10O	028715-26-6	83
2		1H-Indazole, 3,6-dimethyl-	146	C9H10N2	007746-28-3	72
3		2-Propenal, 2-methyl-3-phenyl-	146	C10H10O	000101-39-3	50
4		Benzonitrile, 4-(dimethylamino)-	146	C9H10N2	001197-19-9	45
5		1H-Benzimidazole, 2-ethyl-	146	C9H10N2	001848-84-6	45



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
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Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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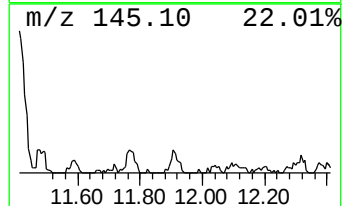
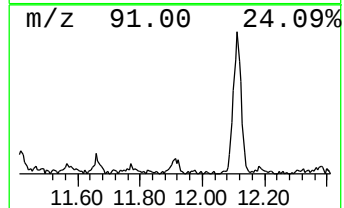
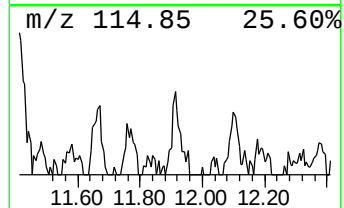
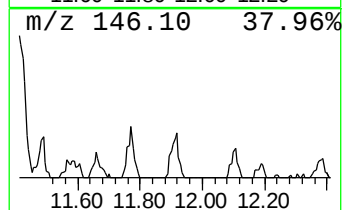
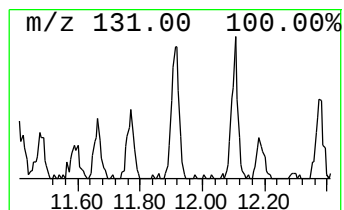
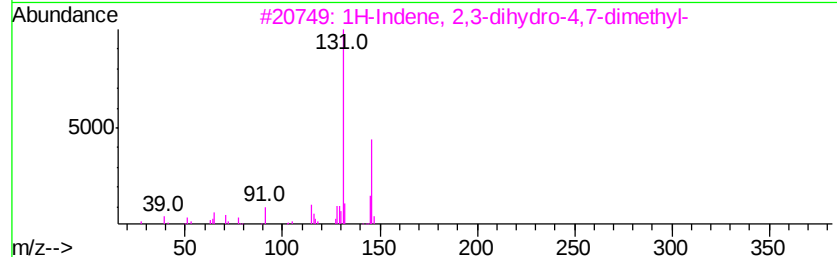
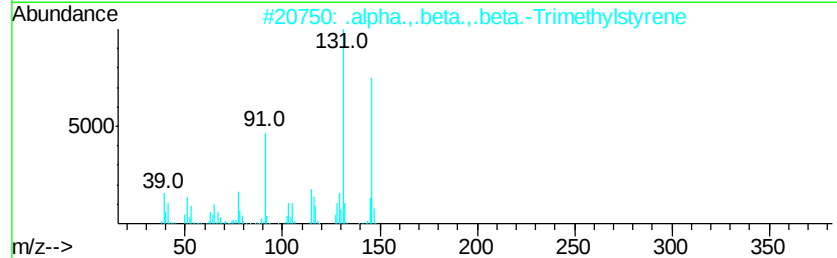
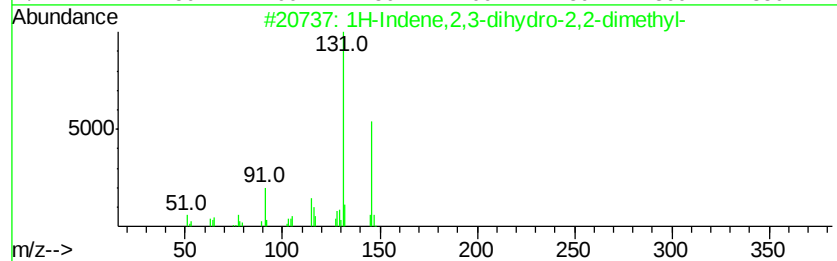
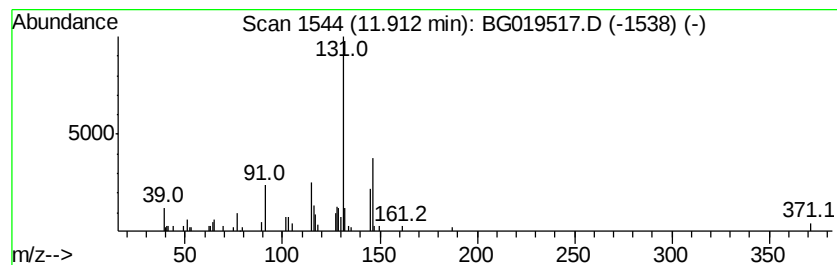
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 1H-Indene,2,3-dihydro-2,2-d... Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.91	2.61 ng/ul	47201	Napthalene-d8	11.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene,2,3-dihydro-2,2-dimethyl-	146	C11H14	020836-11-7	87
2		.alpha.,.beta.,.beta.-Trimethyls...	146	C11H14	000769-57-3	87
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	86
4		1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	83
5		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	81



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
Operator : UM/NP
Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
BCY50

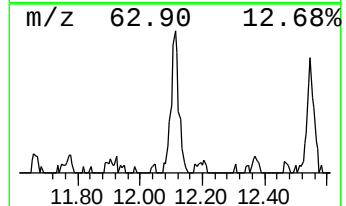
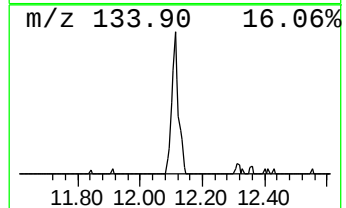
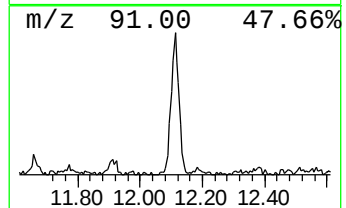
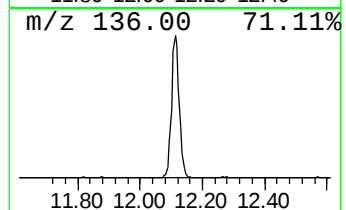
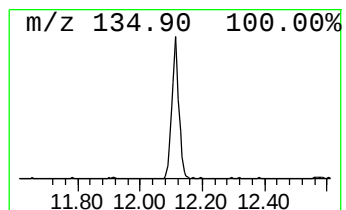
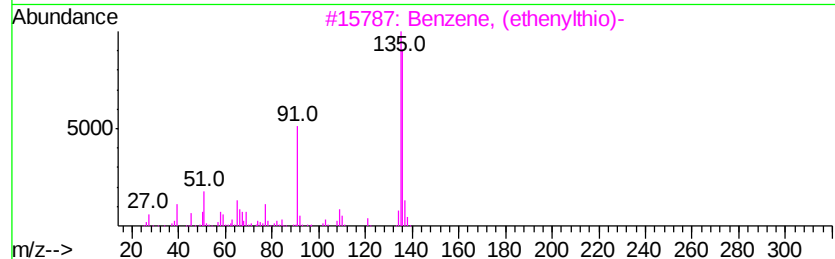
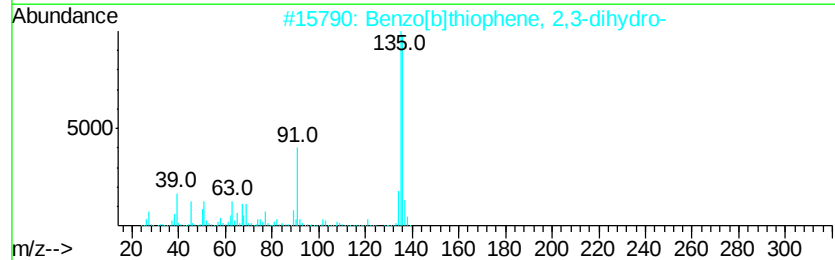
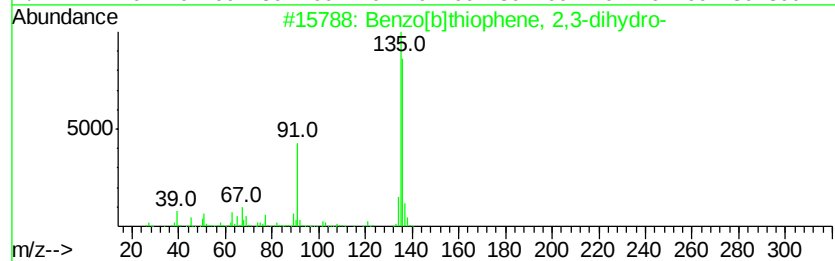
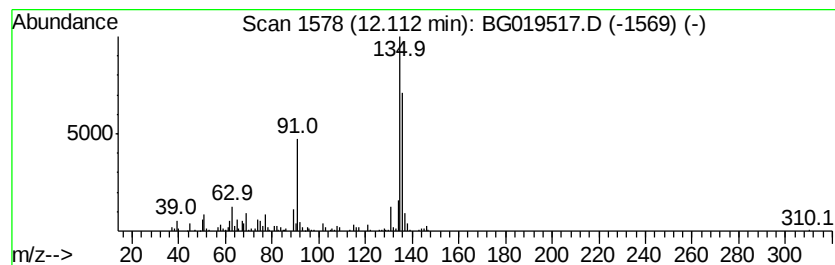
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Benzo[b]thiophene, 2,3-dihy... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.11	13.58 ng/ul	246046	Napthalene-d8	11.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	95
2		Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	93
3		Benzene, (ethenylthio)-	136	C8H8S	001822-73-7	74
4		Benzo[c]thiophene, 1,3-dihydro-	136	C8H8S	002471-92-3	72
5		4-Hydroxy-3-methylbenzaldehyde	136	C8H8O2	015174-69-3	59



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
Operator : UM/NP
Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BCY50

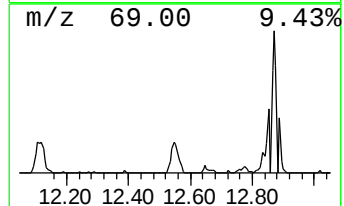
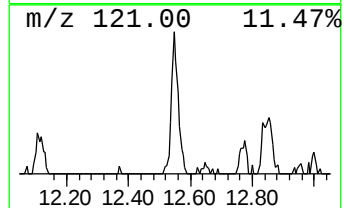
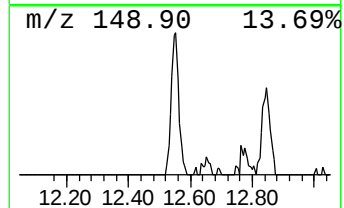
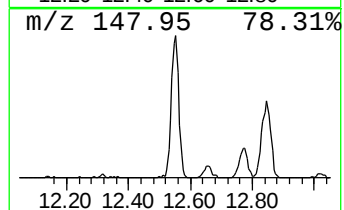
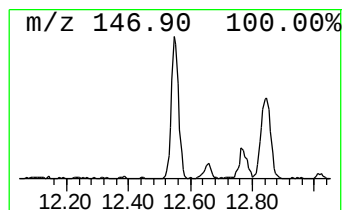
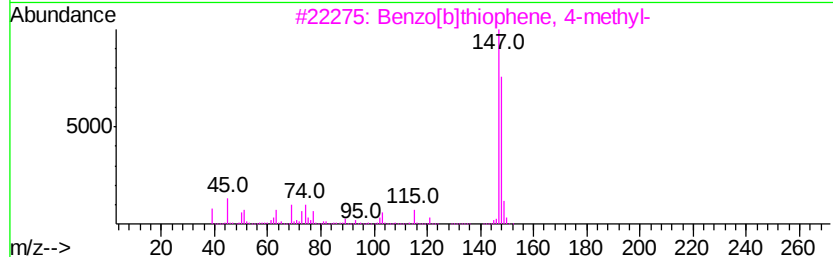
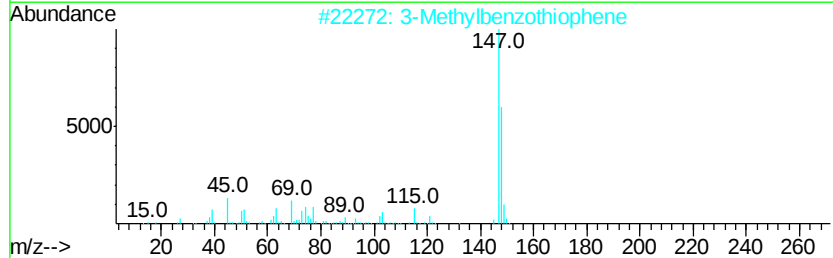
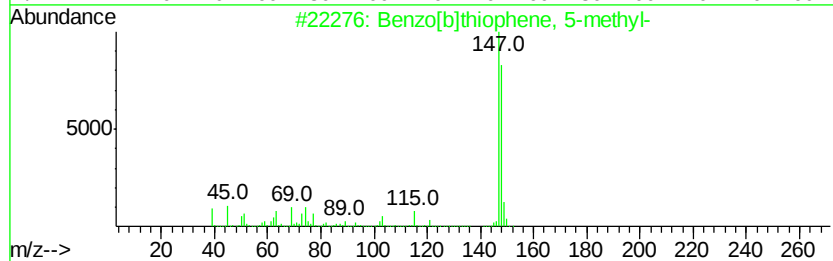
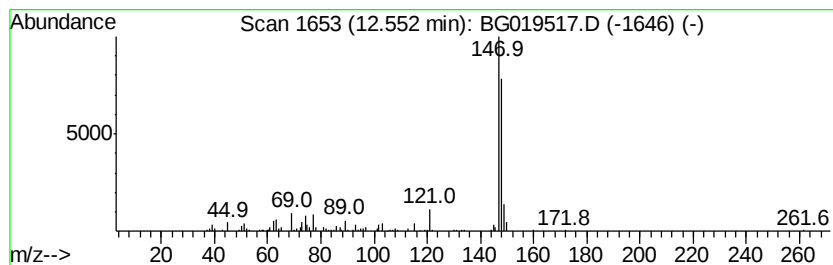
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Benzo[b]thiophene, 5-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.55	10.29 ng/ul	186361	Napthalene-d8	11.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]thiophene, 5-methyl-	148	C9H8S	014315-14-1	87
2		3-Methylbenzothiophene	148	C9H8S	001455-18-1	87
3		Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	83
4		Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	74
5		Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	74



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
Operator : UM/NP
Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BCY50

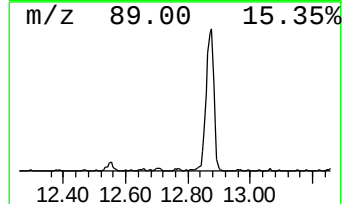
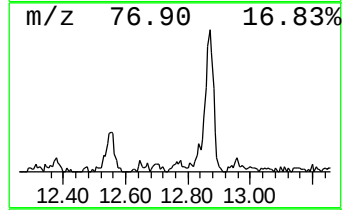
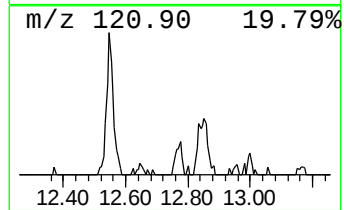
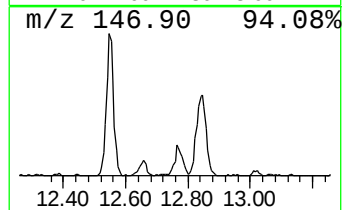
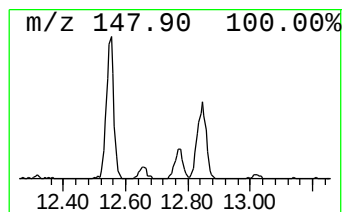
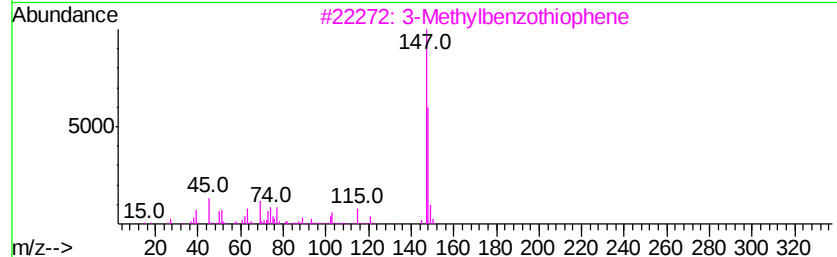
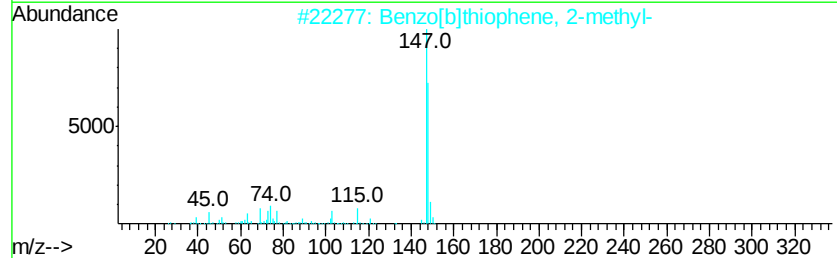
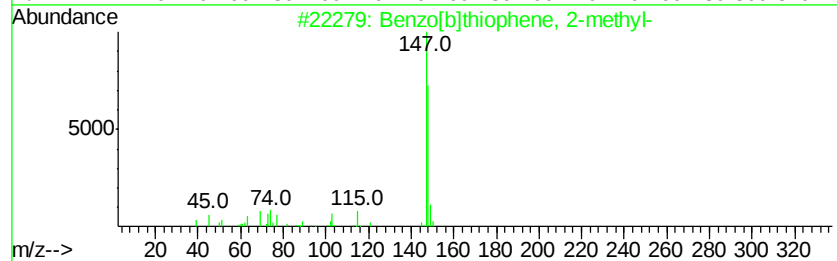
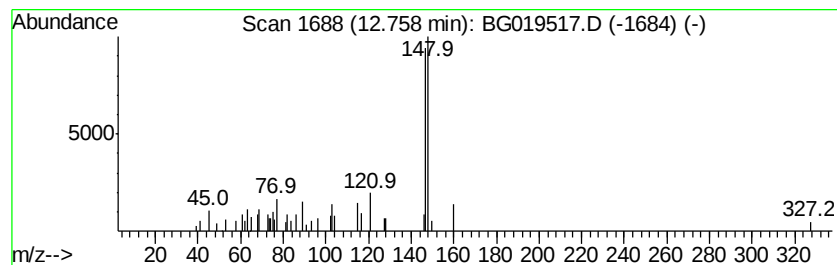
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 Benzo[b]thiophene, 2-methyl- Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	2.36 ng/ul	42657	Napthalene-d8	11.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	87
2		Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	87
3		3-Methylbenzothiophene	148	C9H8S	001455-18-1	80
4		Benzo[b]thiophene, 5-methyl-	148	C9H8S	014315-14-1	80
5		Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	72



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
Operator : UM/NP
Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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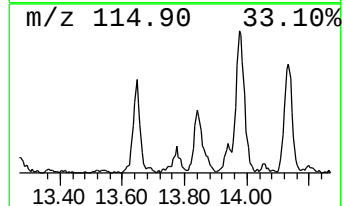
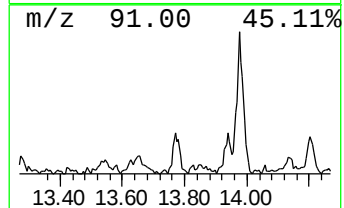
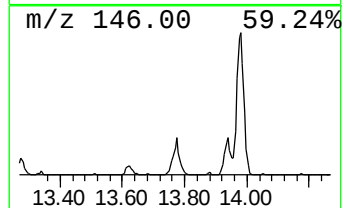
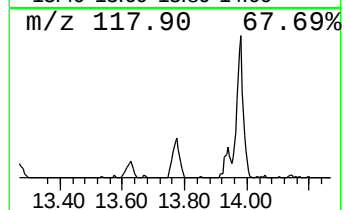
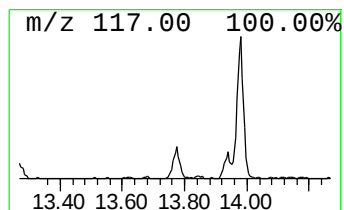
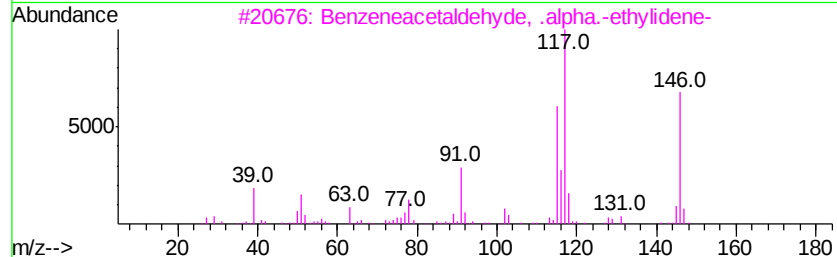
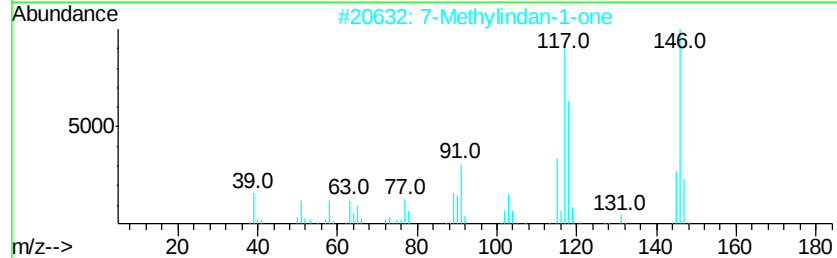
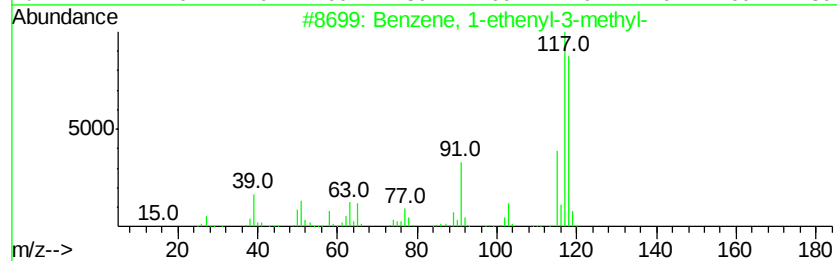
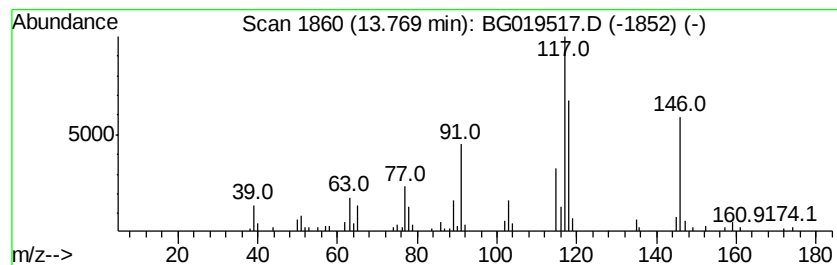
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 Benzene, 1-ethenyl-3-methyl- Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	2.60 ng/ul	84594	Acenaphthene-d10	14.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	90
2			7-Methylindan-1-one	146	C10H10O	039627-61-7	86
3			Benzeneacetaldehyde, .alpha.-eth...	146	C10H10O	004411-89-6	76
4			cis-.beta.-Methylstyrene	118	C9H10	000766-90-5	60
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
Operator : UM/NP
Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

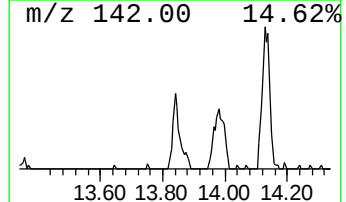
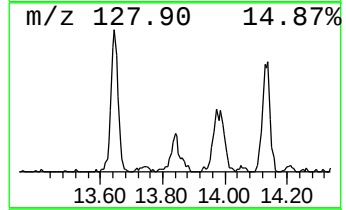
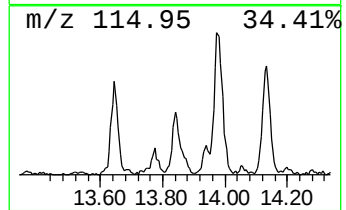
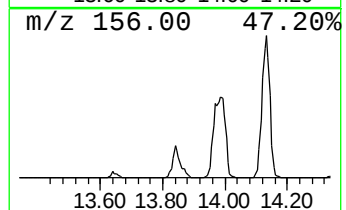
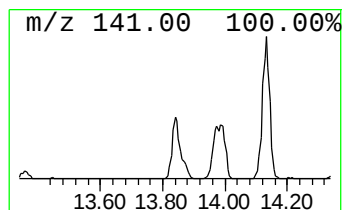
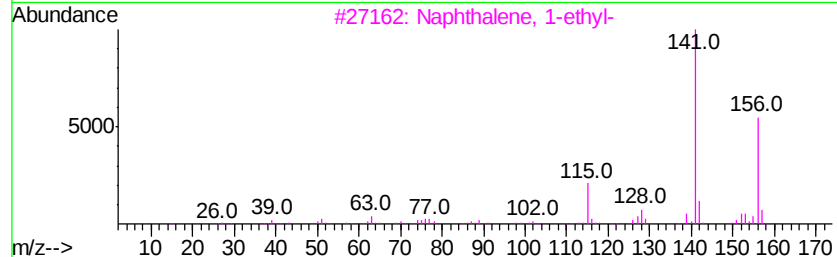
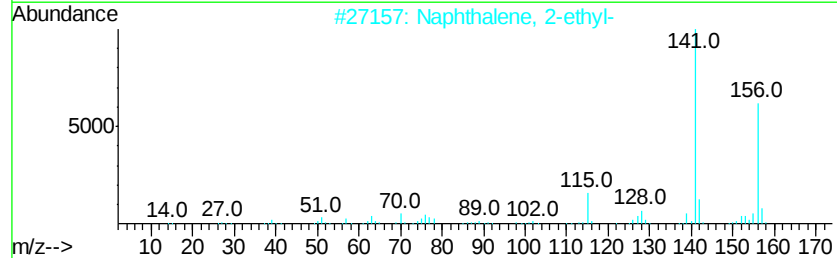
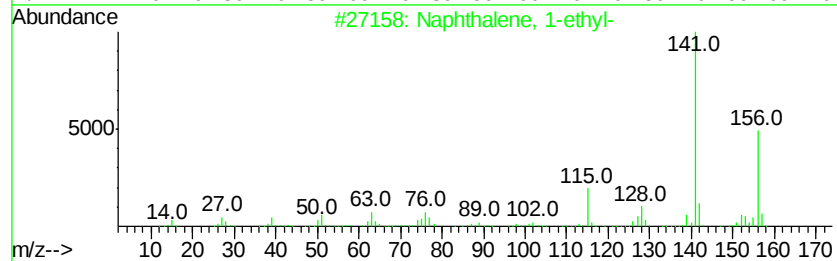
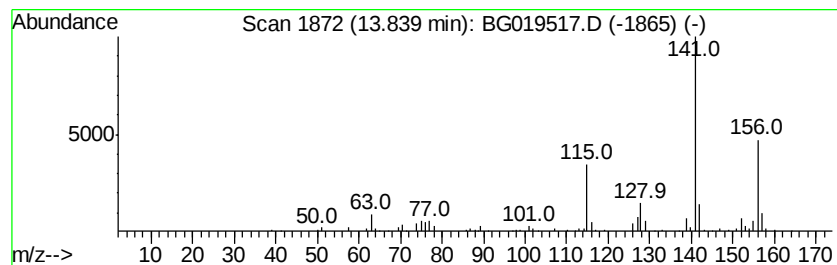
Instrument :
BNA_G
ClientSampleId :
BCY50

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102815.M
Quant Title : SVOA CALIBRATION

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

Peak Number 26 Naphthalene, 1-ethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.84	5.48 ng/ul	178236	Acenaphthene-d10	14.81
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1-ethyl-	156 C12H12	001127-76-0 94
2		Naphthalene, 2-ethyl-	156 C12H12	000939-27-5 94
3		Naphthalene, 1-ethyl-	156 C12H12	001127-76-0 94
4		Naphthalene, 1-ethyl-	156 C12H12	001127-76-0 91
5		Naphthalene, 2-ethyl-	156 C12H12	000939-27-5 91



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
Operator : UM/NP
Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BCY50

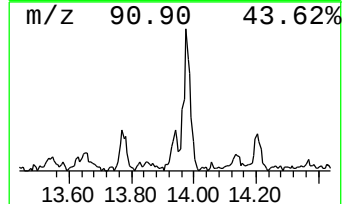
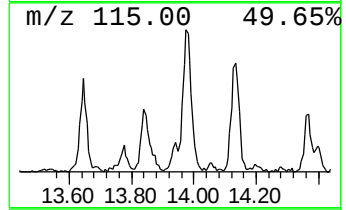
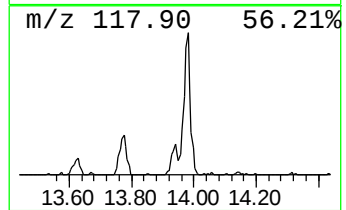
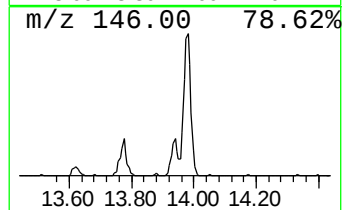
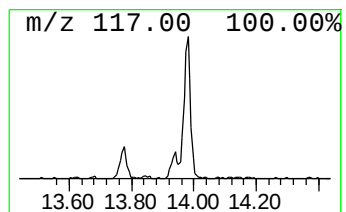
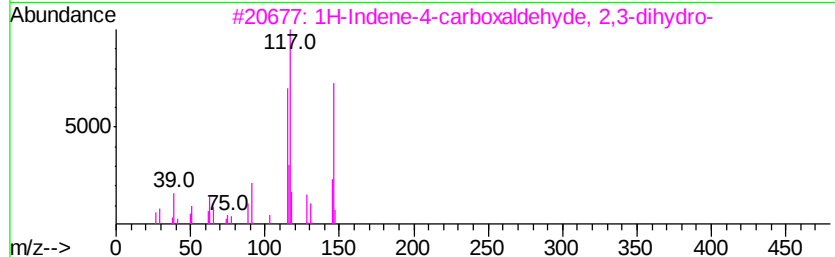
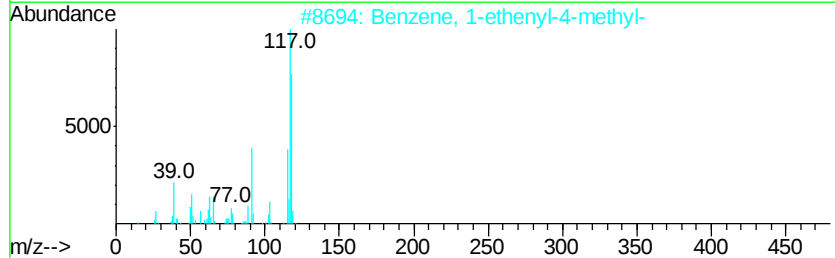
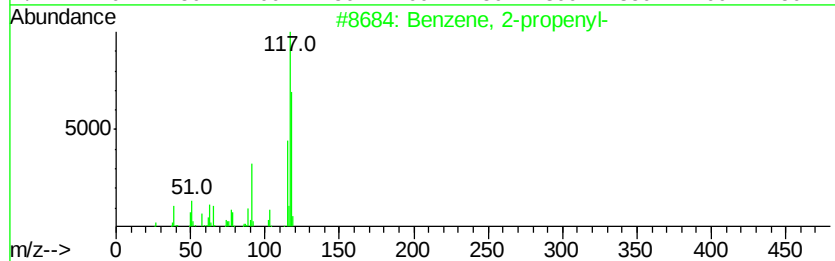
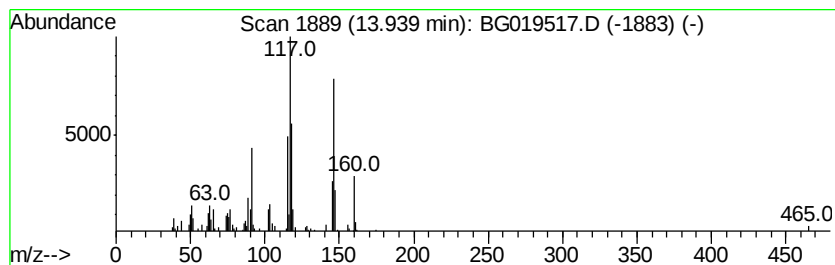
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 Benzene, 2-propenyl- Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.94	2.06 ng/ul	67003	Acenaphthene-d10	14.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	86
2		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	83
3		1H-Indene-4-carboxaldehyde, 2,3-...	146	C10H10O	051932-70-8	52
4		1H-Indene-4-carboxaldehyde, 2,3-...	146	C10H10O	051932-70-8	52
5		cis-.beta.-Methylstyrene	118	C9H10	000766-90-5	46



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
Operator : UM/NP
Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BCY50

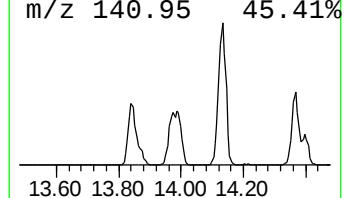
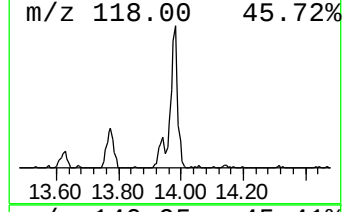
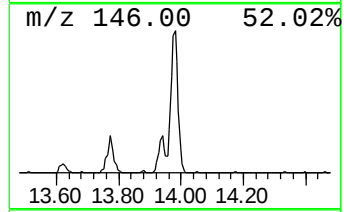
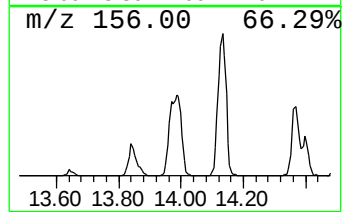
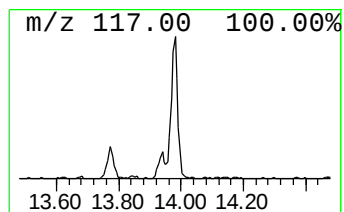
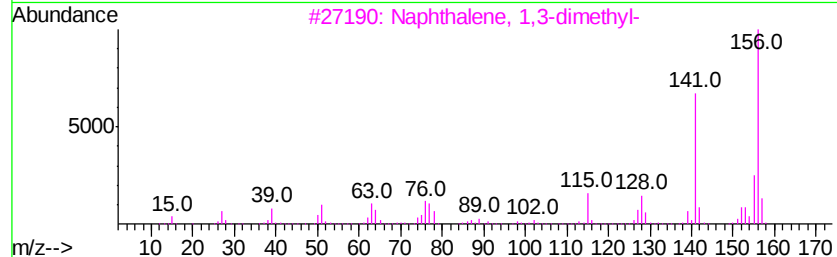
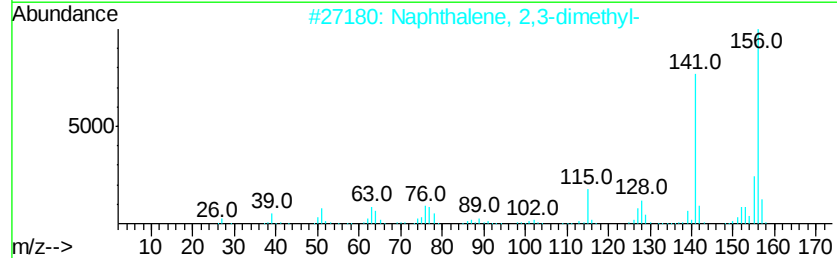
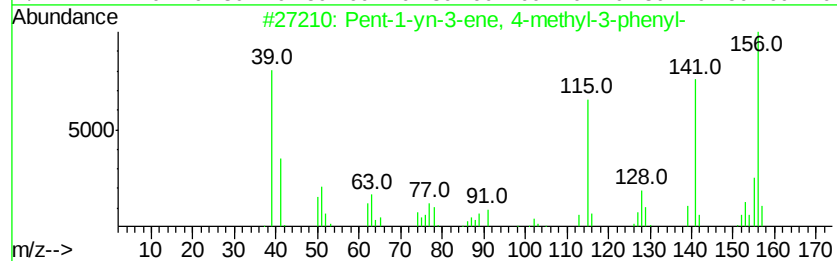
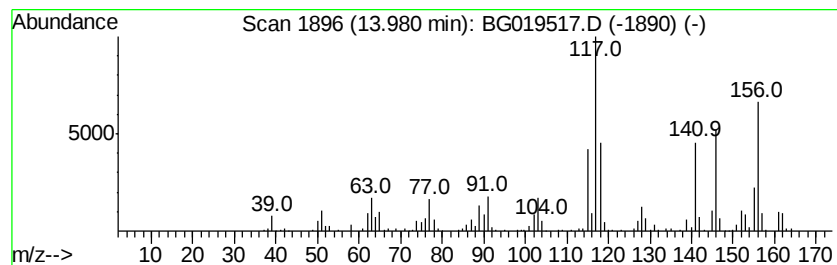
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 Pent-1-yn-3-ene, 4-methyl-3... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.98	18.76 ng/ul	610452	Acenaphthene-d10	14.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pent-1-yn-3-ene, 4-methyl-3-phenyl-	156	C12H12	065050-80-8	60
2		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	56
3		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	53
4		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	50
5		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	47



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
Operator : UM/NP
Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BCY50

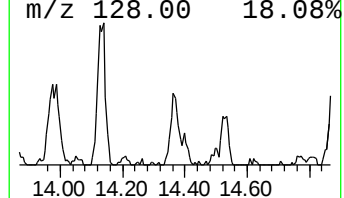
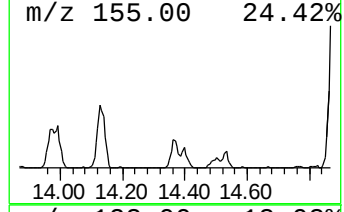
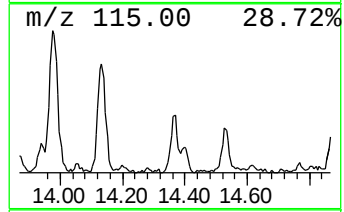
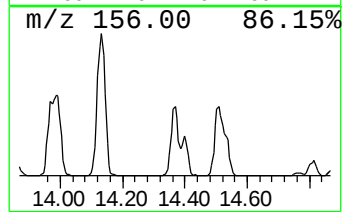
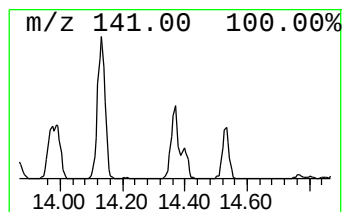
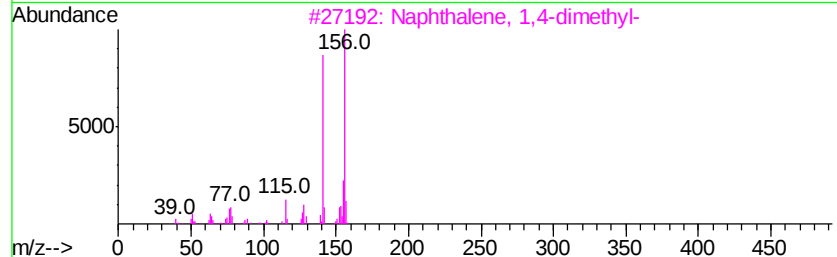
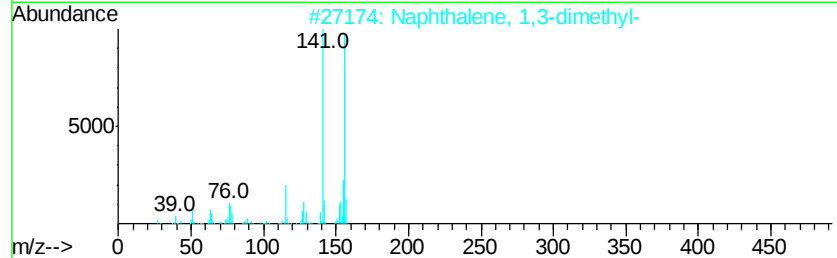
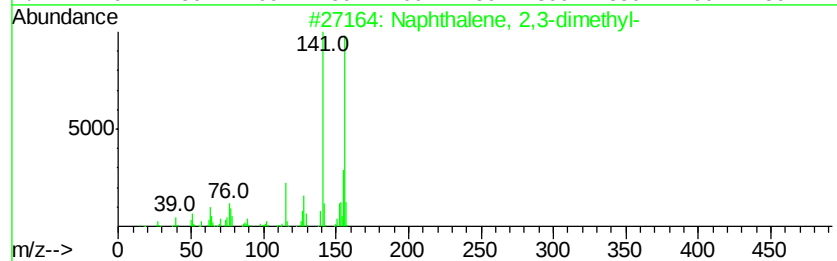
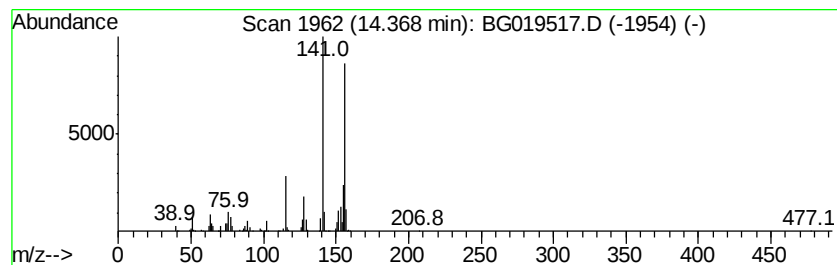
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102815.M
Quant Title : SVOA CALIBRATION

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

Peak Number 30 Naphthalene, 2,3-dimethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.37	5.73 ng/ul	186407	Acenaphthene-d10	14.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
3		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
4		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
5		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95



Data Path : Z:\HPCHEM1\BNA G\DATA\BG110615\
Data File : BG019517.D
Acq On : 6 Nov 2015 18:14
Operator : UM/NP
Sample : G4245-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BCY50

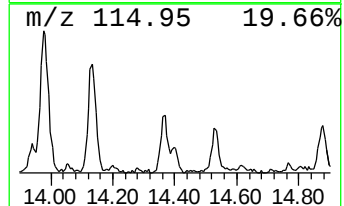
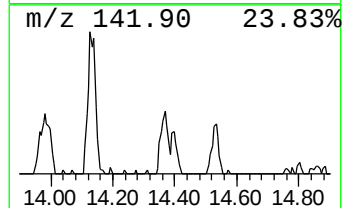
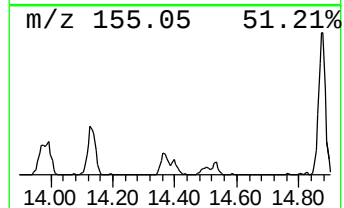
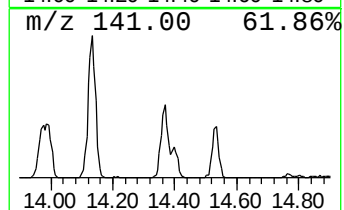
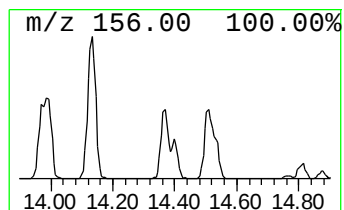
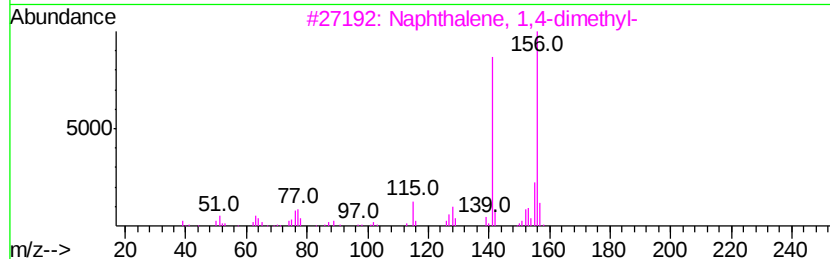
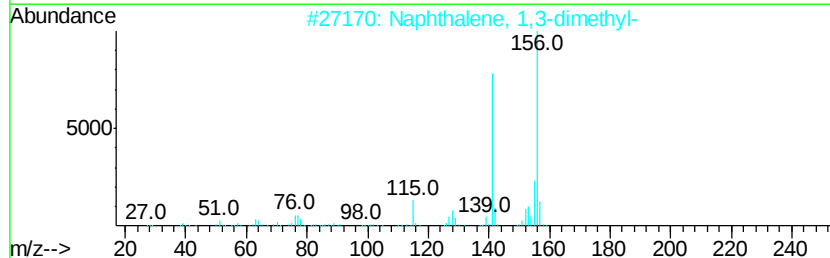
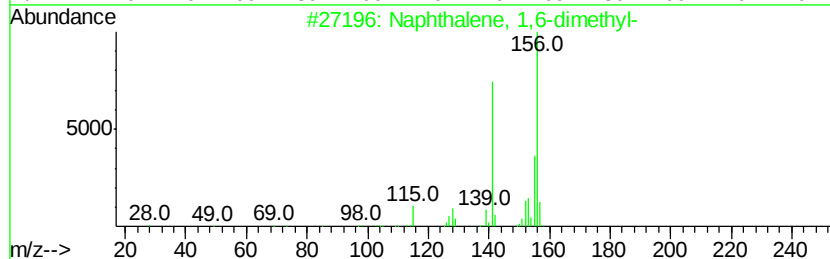
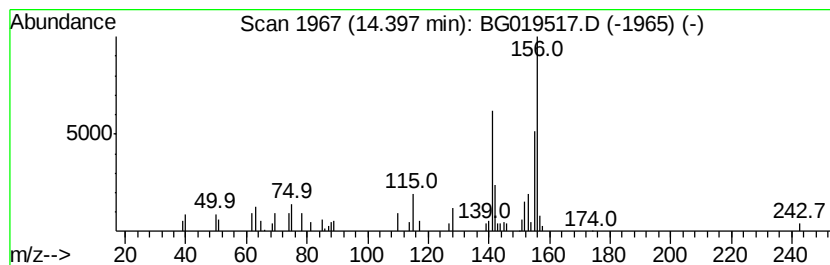
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102815.M
Quant Title : SVOA CALIBRATION

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

Peak Number 31 Naphthalene, 1,6-dimethyl- Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.40	2.14 ng/ul	69709	Acenaphthene-d10	14.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	72
2		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	47
3		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	46
4		Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	43
5		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	43



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG110615\
 Data File : BG019517.D
 Acq On : 6 Nov 2015 18:14
 Operator : UM/NP
 Sample : G4245-15
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BCY50

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102815.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
Butane, 2-methoxy...	3.20	22.9	ng/ul	304067	1	8.18	266095	20.0
Benzene, 1,2,4-tr...	6.64	3.4	ng/ul	44882	1	8.18	266095	20.0
Benzene, 1-ethyl-...	7.56	3.5	ng/ul	46707	1	8.18	266095	20.0
Benzene, 1,3,5-tr...	8.29	6.8	ng/ul	91094	1	8.18	266095	20.0
Indane	8.56	37.2	ng/ul	495078	1	8.18	266095	20.0
Benzene, 1-propynyl-	8.72	20.1	ng/ul	267755	1	8.18	266095	20.0
Benzofuran, 7-met...	9.54	17.5	ng/ul	233140	1	8.18	266095	20.0
Benzofuran, 2-met...	9.67	21.9	ng/ul	396088	2	11.01	362242	20.0
Benzene, 1-ethyl-...	9.81	2.9	ng/ul	52849	2	11.01	362242	20.0
Benzene, 1-methyl...	9.88	2.3	ng/ul	42085	2	11.01	362242	20.0
1H-Indene, 2,3-di...	10.24	7.4	ng/ul	134812	2	11.01	362242	20.0
3-Phenylbut-1-ene	10.40	14.7	ng/ul	265862	2	11.01	362242	20.0
Benzo[b]thiophene	11.18	110.8	ng/ul	2007330	2	11.01	362242	20.0
2-Propenal, 2-met...	11.24	2.5	ng/ul	45574	2	11.01	362242	20.0
Benzofuran, 4,7-d...	11.41	4.7	ng/ul	85674	2	11.01	362242	20.0
1H-Indene, 2,3-dih...	11.91	2.6	ng/ul	47201	2	11.01	362242	20.0
Benzo[b]thiophene...	12.11	13.6	ng/ul	246046	2	11.01	362242	20.0
Benzo[b]thiophene...	12.55	10.3	ng/ul	186361	2	11.01	362242	20.0
Benzo[b]thiophene...	12.76	2.4	ng/ul	42657	2	11.01	362242	20.0
Benzene, 1-etheny...	13.77	2.6	ng/ul	84594	3	14.81	650777	20.0
Naphthalene, 1-et...	13.84	5.5	ng/ul	178236	3	14.81	650777	20.0
Benzene, 2-propenyl-	13.94	2.1	ng/ul	67003	3	14.81	650777	20.0
Pent-1-yn-3-ene, ...	13.98	18.8	ng/ul	610452	3	14.81	650777	20.0
Naphthalene, 2,3-...	14.37	5.7	ng/ul	186407	3	14.81	650777	20.0
Naphthalene, 1,6-...	14.40	2.1	ng/ul	69709	3	14.81	650777	20.0