

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG122615\
 Data File : BG020564.D
 Acq On : 27 Dec 2015 21:13
 Operator : UM/SJ
 Sample : G4846-11
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 D9N57

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-BG122415.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	2.875	3	6	10	rVB2	2638	3411	0.28%	0.031%
2	3.093	39	43	53	rBV	39956	72310	5.98%	0.656%
3	3.175	53	57	64	rVB	27367	38383	3.18%	0.348%
4	5.555	454	462	467	rBV	4399	7588	0.63%	0.069%
5	5.696	481	486	494	rVB3	4068	8991	0.74%	0.082%
6	6.066	543	549	555	rVB2	4531	6937	0.57%	0.063%
7	7.153	728	734	740	rBB2	3052	5520	0.46%	0.050%
8	7.235	740	748	753	rBV3	20045	36551	3.02%	0.332%
9	7.282	753	756	762	rVB2	4737	6747	0.56%	0.061%
10	7.394	769	775	782	rBV	61785	101210	8.37%	0.919%
11	7.452	782	785	792	rVB	3152	4629	0.38%	0.042%
12	7.605	804	811	820	rVB	66825	112608	9.32%	1.022%
13	7.717	825	830	835	rVB	11692	17734	1.47%	0.161%
14	8.075	883	891	898	rBV	67466	112053	9.27%	1.017%
15	8.187	905	910	916	rVB2	5211	8565	0.71%	0.078%
16	8.451	949	955	962	rVB	22516	36844	3.05%	0.334%
17	8.616	977	983	991	rVB2	13653	23267	1.92%	0.211%
18	8.710	995	999	1004	rVV	2935	3832	0.32%	0.035%
19	8.786	1004	1012	1025	rVB	58969	104581	8.65%	0.949%
20	9.174	1076	1078	1084	rVB3	2294	3531	0.29%	0.032%
21	9.244	1084	1090	1100	rBV	83915	150229	12.43%	1.364%
22	9.703	1163	1168	1173	rBV2	1808	2879	0.24%	0.026%
23	9.767	1173	1179	1186	rVB2	2952	4820	0.40%	0.044%
24	9.967	1207	1213	1222	rBB	72968	132537	10.96%	1.203%
25	10.132	1236	1241	1247	rBB2	2843	4565	0.38%	0.041%
26	10.290	1260	1268	1278	rBB6	8255	19663	1.63%	0.178%
27	10.390	1279	1285	1289	rBV3	2845	6513	0.54%	0.059%
28	10.419	1289	1290	1295	rVB2	1941	1951	0.16%	0.018%
29	10.513	1299	1306	1314	rBV	112202	196474	16.25%	1.784%
30	10.895	1363	1371	1374	rBV	99305	181372	15.00%	1.646%
31	10.948	1374	1380	1391	rVV	670473	1208864	100.00%	10.974%
32	11.066	1393	1400	1409	rVB	16707	29338	2.43%	0.266%
33	12.441	1629	1634	1637	rBV2	2753	4542	0.38%	0.041%
34	12.546	1644	1652	1661	rBV	193852	327116	27.06%	2.969%

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35	12.664	1665	1672	1678	rBB2	3745	6645	0.55%	0.060%
36	12.758	1679	1688	1698	rBV	122481	207052	17.13%	1.880%
37	13.545	1816	1822	1828	rBV	59804	89698	7.42%	0.814%
38	13.739	1848	1855	1864	rVB2	11893	20775	1.72%	0.189%
39	13.886	1871	1880	1886	rBB5	10821	27582	2.28%	0.250%
40	14.027	1897	1904	1909	rBV	22581	39567	3.27%	0.359%
41	14.109	1909	1918	1925	rVB	250530	368293	30.47%	3.343%
42	14.268	1940	1945	1948	rBV	9999	13253	1.10%	0.120%
43	14.297	1948	1950	1955	rVB2	2731	3411	0.28%	0.031%
44	14.350	1955	1959	1962	rBV2	4662	6187	0.51%	0.056%
45	14.409	1962	1969	1979	rVB	254297	440440	36.43%	3.998%
46	14.714	2010	2021	2026	rBV2	183924	316443	26.18%	2.873%
47	14.773	2026	2031	2038	rVV	360124	565060	46.74%	5.129%
48	14.838	2039	2042	2050	rVB	7000	10595	0.88%	0.096%
49	14.914	2050	2055	2063	rBV3	17481	31506	2.61%	0.286%
50	15.020	2068	2073	2077	rVB	5153	7422	0.61%	0.067%
51	15.108	2081	2088	2096	rVV	170441	264442	21.88%	2.401%
52	15.179	2096	2100	2105	rVB3	1387	1886	0.16%	0.017%
53	15.308	2119	2122	2129	rVB3	1739	3216	0.27%	0.029%
54	15.372	2130	2133	2137	rBV	1172	1848	0.15%	0.017%
55	15.490	2149	2153	2156	rVV2	1245	1739	0.14%	0.016%
56	15.578	2164	2168	2173	rBB2	1987	3048	0.25%	0.028%
57	15.701	2179	2189	2194	rBV	369190	563008	46.57%	5.111%
58	15.754	2194	2198	2204	rVV	213961	321272	26.58%	2.916%
59	15.825	2204	2210	2216	rVV	102951	153530	12.70%	1.394%
60	15.878	2216	2219	2222	rVV2	6546	9139	0.76%	0.083%
61	15.919	2222	2226	2231	rVB3	8536	14727	1.22%	0.134%
62	16.007	2237	2241	2244	rBV	4117	5042	0.42%	0.046%
63	16.048	2244	2248	2253	rVV3	8345	11157	0.92%	0.101%
64	16.189	2261	2272	2281	rBB3	10149	24006	1.99%	0.218%
65	16.283	2282	2288	2294	rBB2	2196	3684	0.30%	0.033%
66	16.430	2310	2313	2319	rVB4	1817	2491	0.21%	0.023%
67	16.548	2328	2333	2337	rBV2	8894	12386	1.02%	0.112%
68	16.712	2357	2361	2362	rBV	1937	2013	0.17%	0.018%
69	16.759	2365	2369	2373	rVV3	5431	8843	0.73%	0.080%
70	16.806	2373	2377	2382	rVB2	2204	3363	0.28%	0.031%
71	16.906	2391	2394	2399	rVB2	3040	3515	0.29%	0.032%

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72	17.106	2420	2428	2432	rBV4	2902	6335	0.52%	0.058%
73	17.276	2452	2457	2464	rVB	25226	37149	3.07%	0.337%
74	17.458	2481	2488	2491	rBV	258280	396760	32.82%	3.602%
75	17.499	2491	2495	2500	rVV	367292	541193	44.77%	4.913%
76	17.558	2500	2505	2517	rVB	406829	660695	54.65%	5.998%
77	17.858	2546	2556	2564	rVB	38450	57203	4.73%	0.519%
78	17.964	2568	2574	2579	rBV3	1944	3844	0.32%	0.035%
79	18.334	2628	2637	2640	rBV4	6473	12205	1.01%	0.111%
80	18.381	2640	2645	2654	rVB2	12713	23439	1.94%	0.213%
81	18.451	2654	2657	2664	rVB3	2568	3829	0.32%	0.035%
82	18.528	2664	2670	2675	rBV	23201	38283	3.17%	0.348%
83	18.627	2679	2687	2692	rBV3	3170	5835	0.48%	0.053%
84	18.674	2692	2695	2699	rVB	4201	4659	0.39%	0.042%
85	18.769	2706	2711	2717	rVB3	2767	5413	0.45%	0.049%
86	18.821	2717	2720	2724	rBV2	3908	5445	0.45%	0.049%
87	18.868	2724	2728	2732	rVB2	3788	5409	0.45%	0.049%
88	19.503	2825	2836	2842	rBV	54972	81879	6.77%	0.743%
89	19.726	2871	2874	2877	rBV	3061	3640	0.30%	0.033%
90	19.838	2886	2893	2907	rBV	518319	802628	66.40%	7.286%
91	20.243	2957	2962	2968	rVB2	5346	7952	0.66%	0.072%
92	20.355	2979	2981	2985	rVB2	1593	1951	0.16%	0.018%
93	20.455	2995	2998	3002	rBV2	1738	2476	0.20%	0.022%
94	20.731	3043	3045	3049	rVB5	1403	2031	0.17%	0.018%
95	21.436	3163	3165	3174	rVB6	2021	2808	0.23%	0.025%
96	21.595	3190	3192	3202	rVB4	1948	3182	0.26%	0.029%
97	21.730	3208	3215	3226	rBV	308034	521043	43.10%	4.730%
98	24.009	3600	3603	3604	rBV3	2184	2176	0.18%	0.020%
99	24.779	3722	3734	3753	rBV	247091	752986	62.29%	6.835%
100	25.002	3761	3772	3793	rVB2	158327	469230	38.82%	4.259%

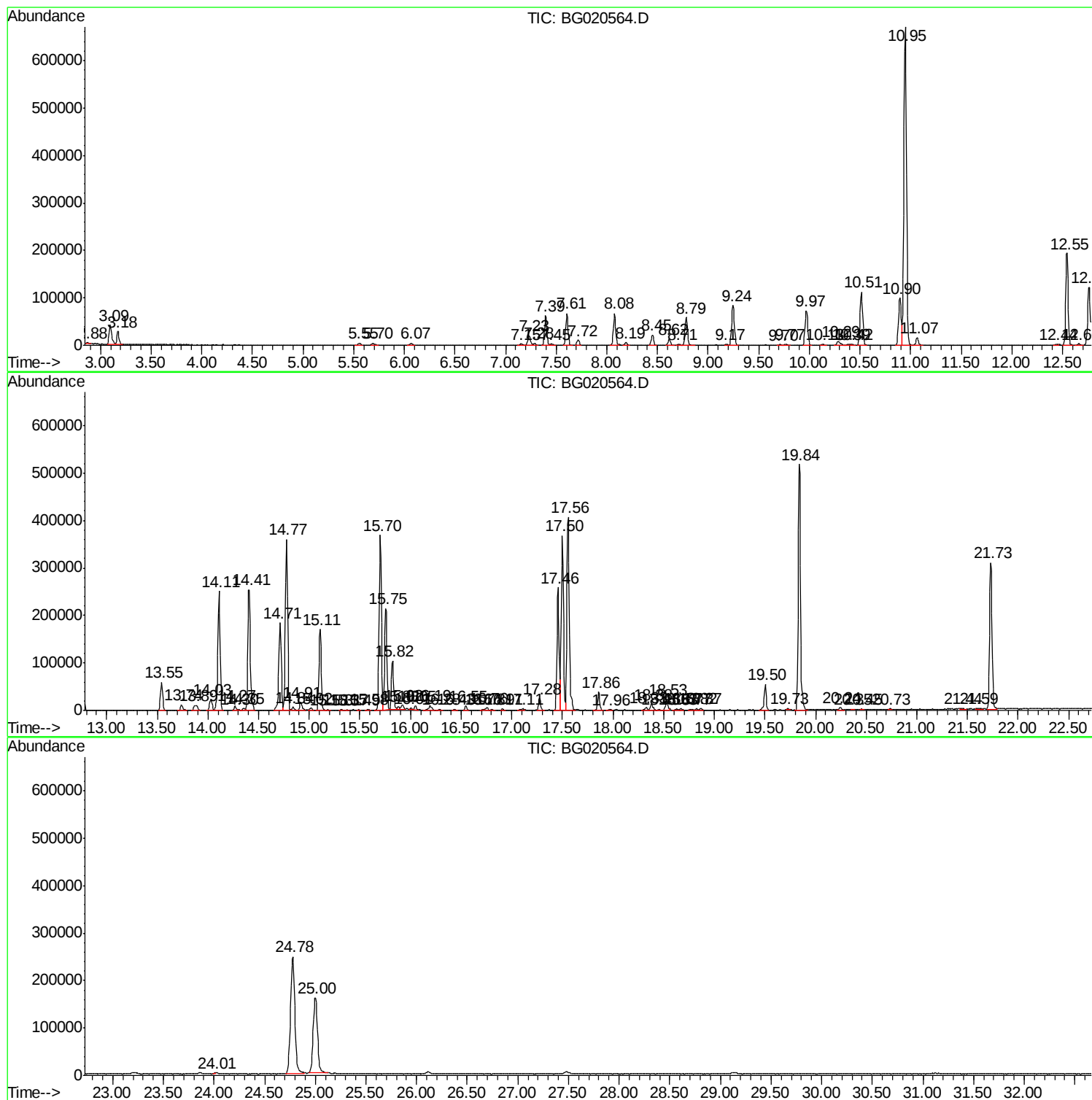
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG122415.M
Quant Title : SVOA CALIBRATION

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



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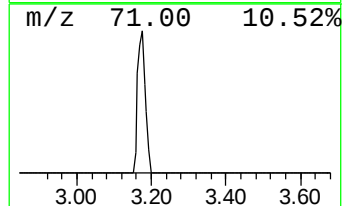
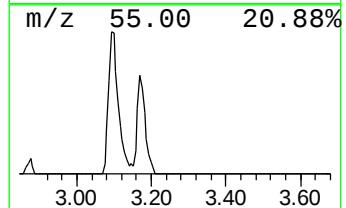
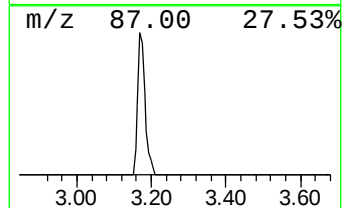
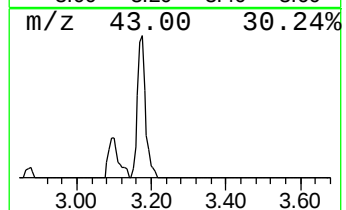
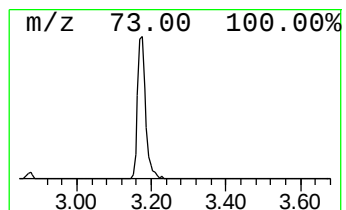
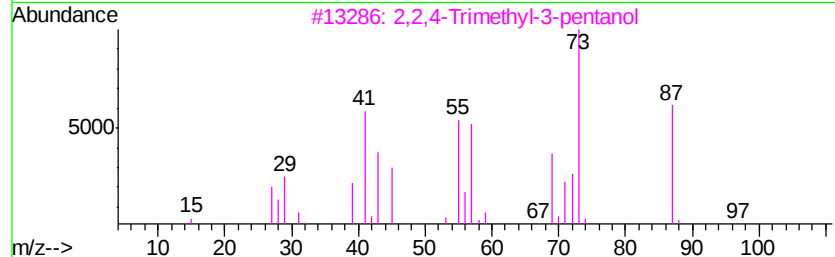
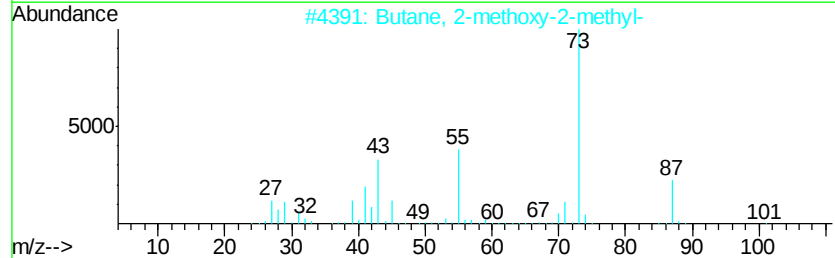
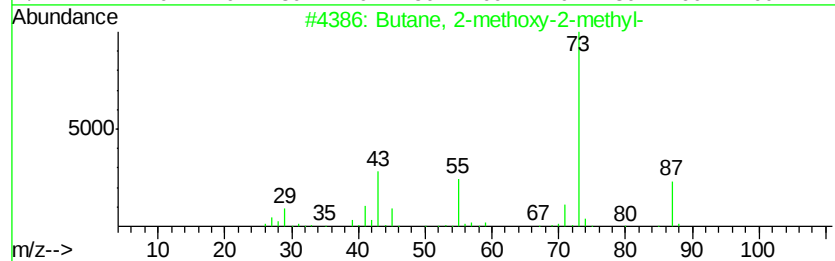
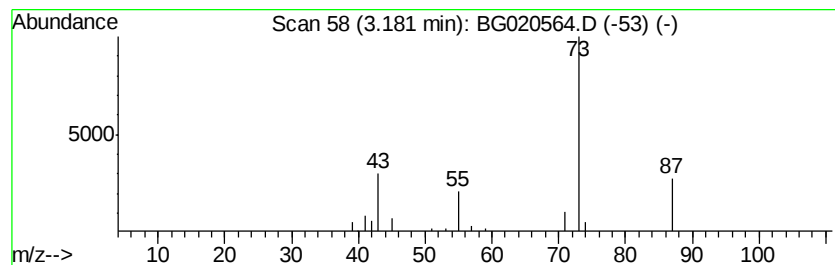
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG122415.M
Quant Title : SVOA CALIBRATION

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.
3.18	6.85 ng/ul	38383	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	56
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43
4		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	12
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	12



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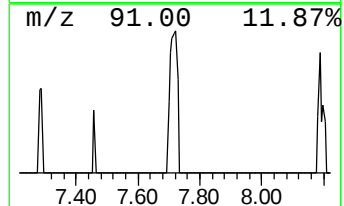
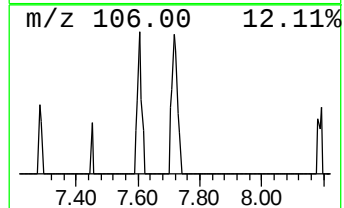
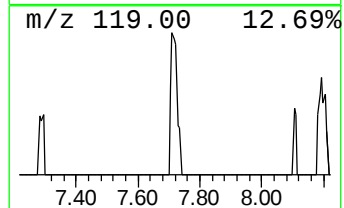
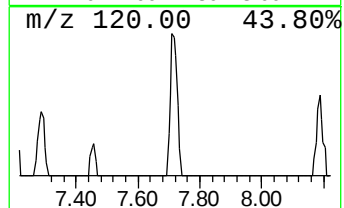
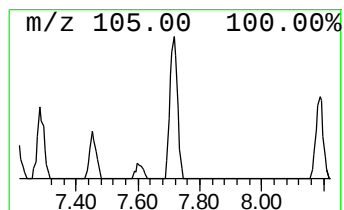
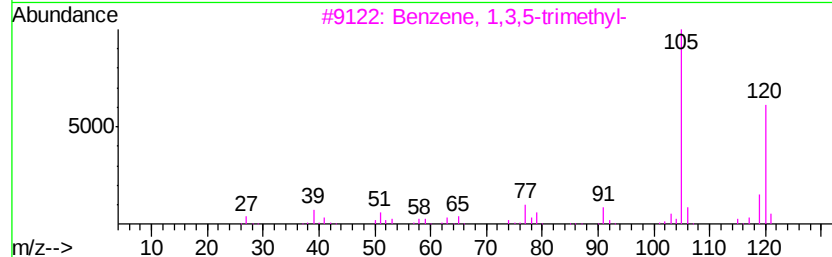
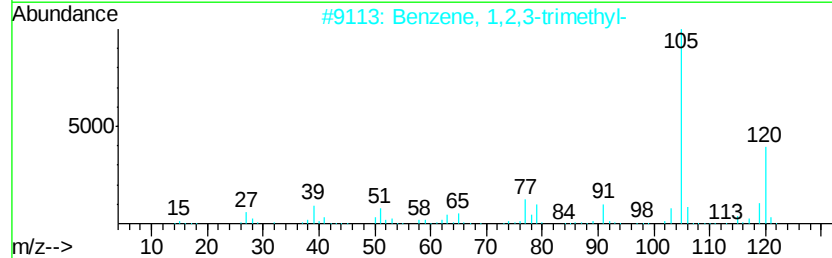
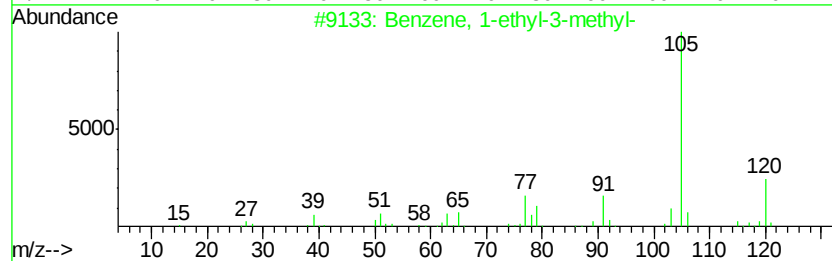
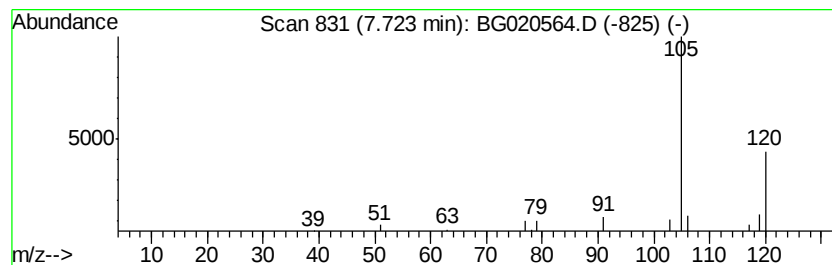
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1-ethyl-3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.72	3.17 ng/ul	17734	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91



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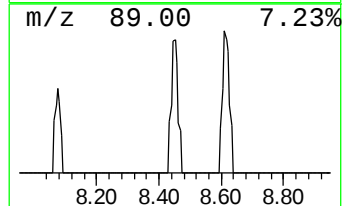
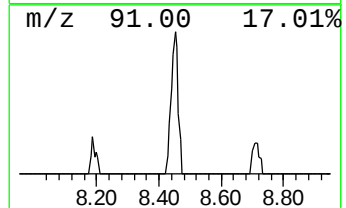
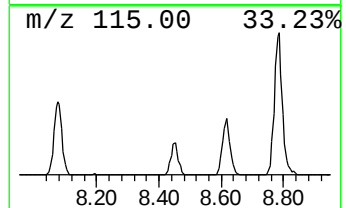
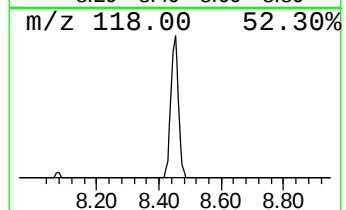
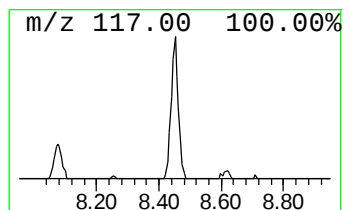
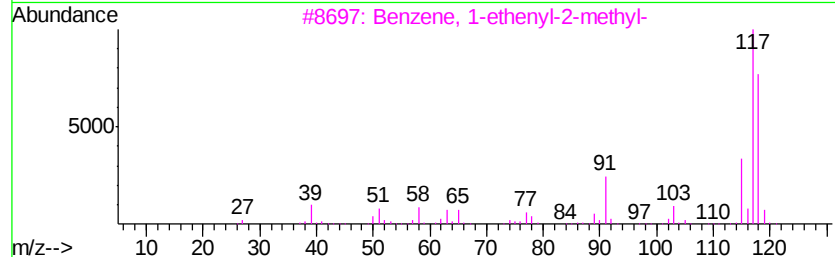
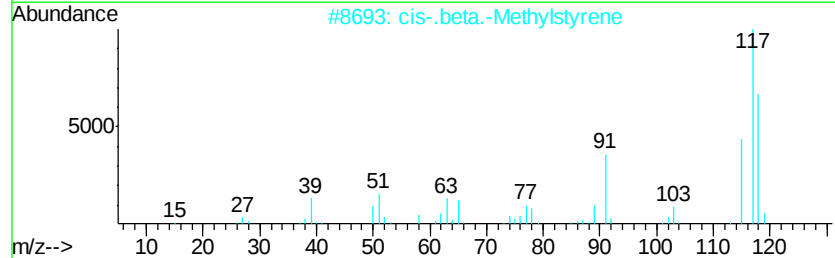
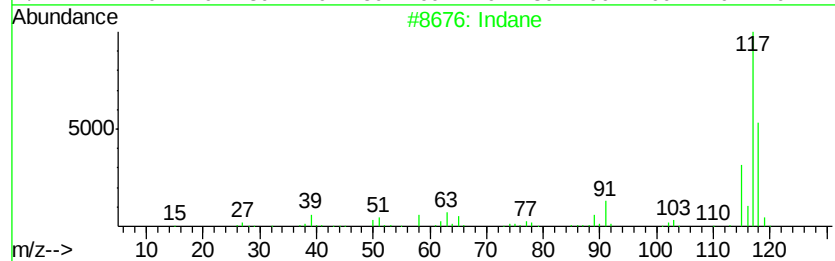
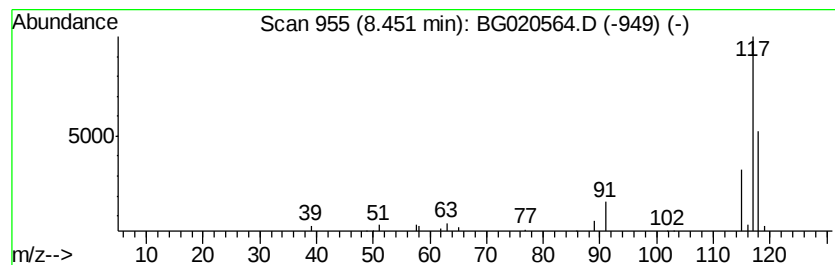
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TIC Integration Parameters: LSCINT.P

Peak Number 4 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.45	6.58 ng/ul	36844	1,4-Dichlorobenzene-d4	8.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	80
2			cis-.beta.-Methylstyrene	118	C9H10	000766-90-5	72
3			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	64
4			Benzeneethanol, .beta.-ethenyl-	148	C10H12O	006052-63-7	64
5			Deltacyclene	118	C9H10	007785-10-6	59



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG122615\
Data File : BG020564.D
Acq On : 27 Dec 2015 21:13
Operator : UM/SJ
Sample : G4846-11
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
D9N57

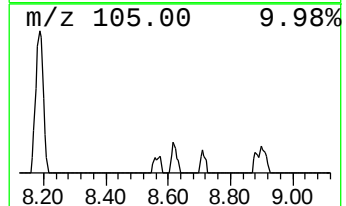
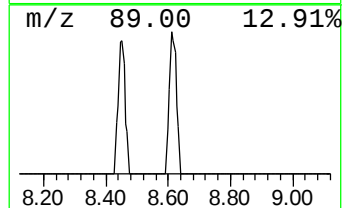
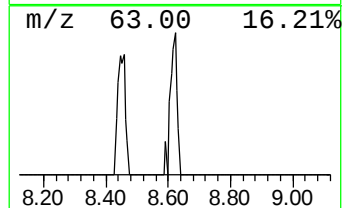
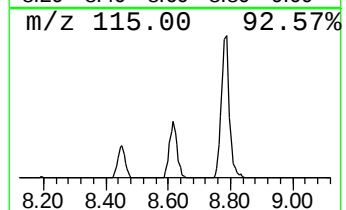
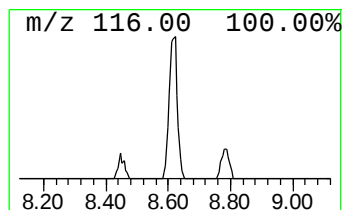
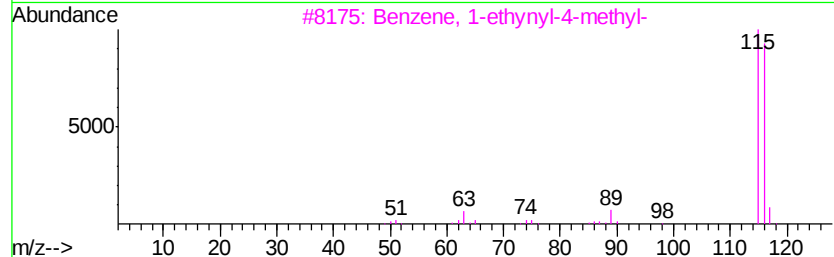
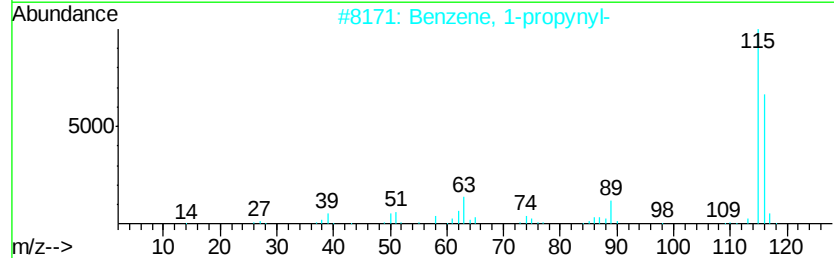
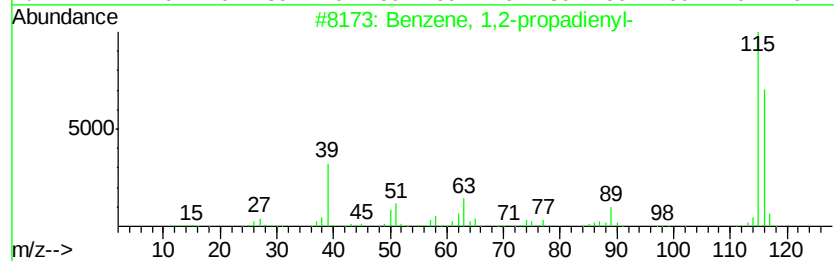
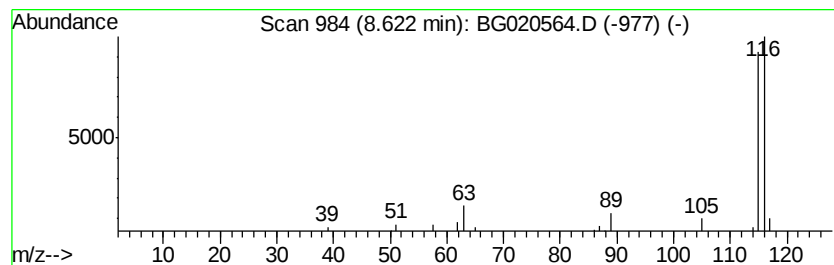
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1,2-propadienyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.62	4.15 ng/ul	23267	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	86
2		Benzene, 1-propynyl-	116	C9H8	000673-32-5	86
3		Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	64
4		Benzene, 1-propynyl-	116	C9H8	000673-32-5	64
5		Benzene, 1-propynyl-	116	C9H8	000673-32-5	64



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG122615\
Data File : BG020564.D
Acq On : 27 Dec 2015 21:13
Operator : UM/SJ
Sample : G4846-11
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
D9N57

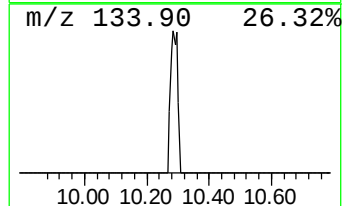
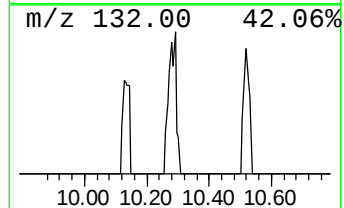
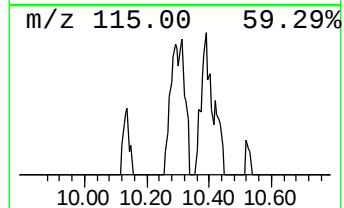
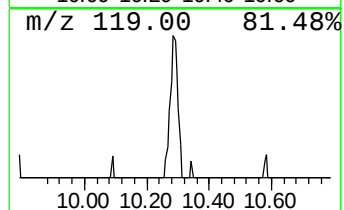
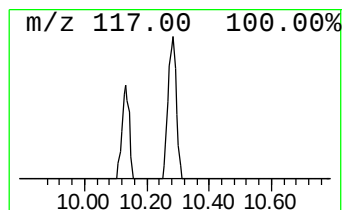
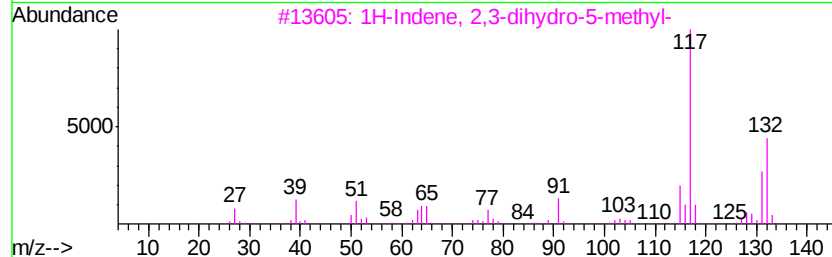
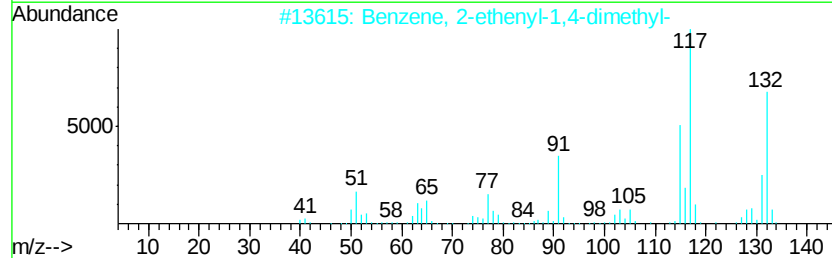
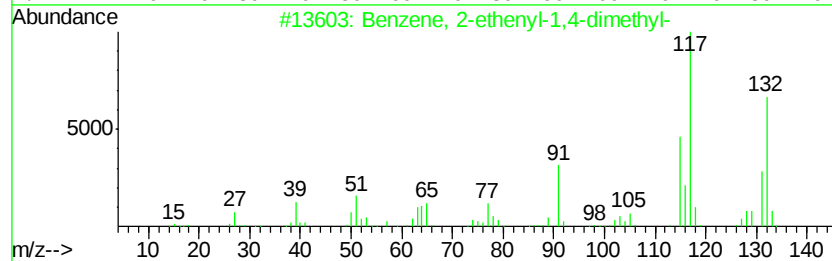
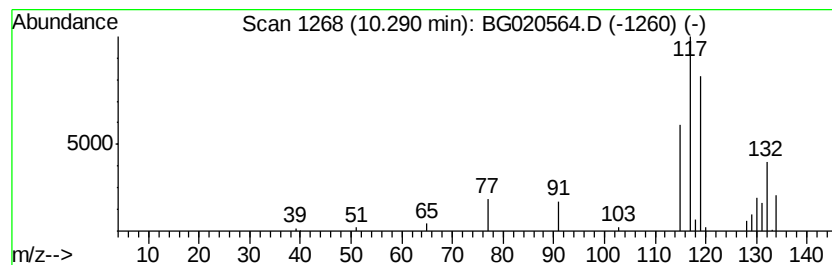
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.29	2.17 ng/ul	19663	Naphthalene-d8	10.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	49
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	46
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	43
4		2,4-Dimethylstyrene	132	C10H12	002234-20-0	43
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	43



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG122615\
Data File : BG020564.D
Acq On : 27 Dec 2015 21:13
Operator : UM/SJ
Sample : G4846-11
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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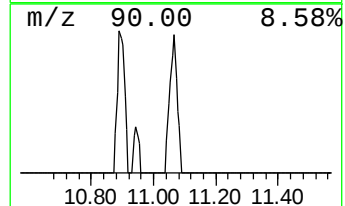
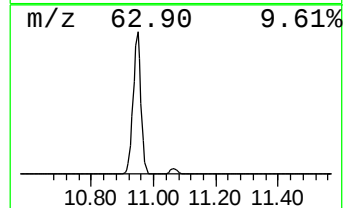
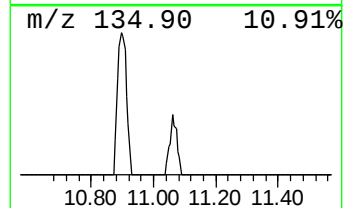
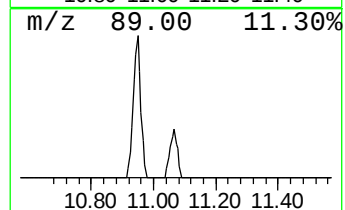
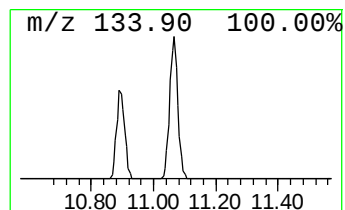
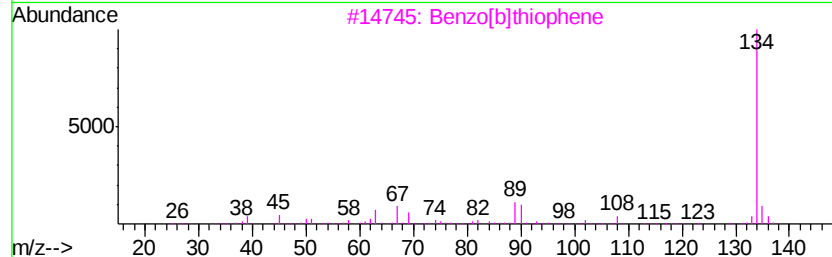
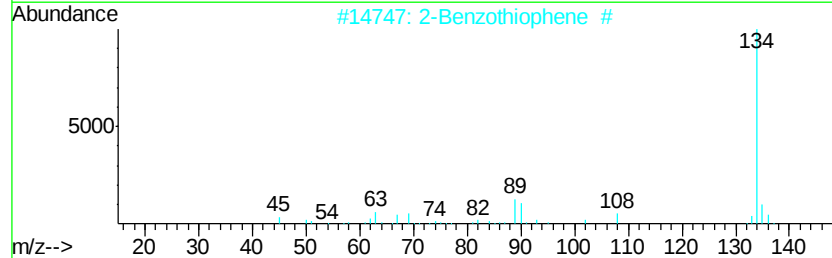
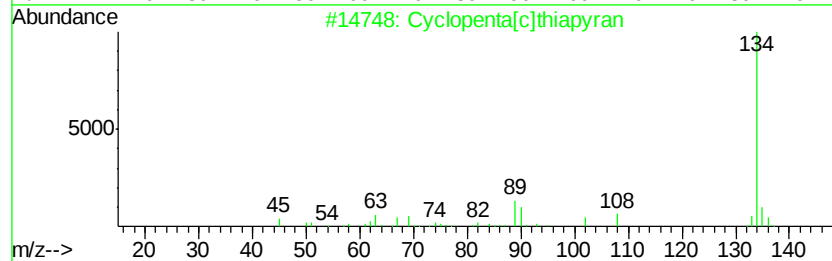
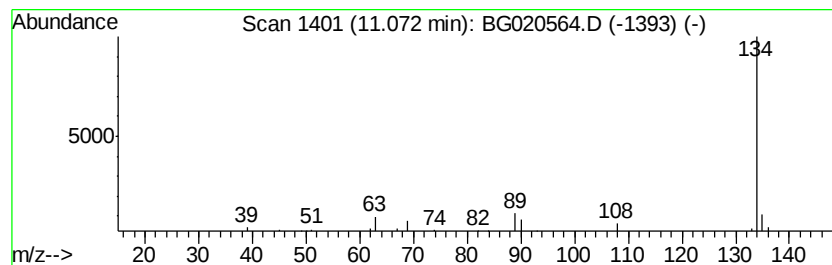
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Cyclopenta[c]thiapyran Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.07	3.24 ng/ul	29338	Napthalene-d8	10.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	94
2		2-Benzothiophene #	134	C8H6S	000270-82-6	94
3		Benzo[b]thiophene	134	C8H6S	000095-15-8	91
4		Cyclopenta[b]thiapyran	134	C8H6S	000271-17-0	91
5		Benzo[b]thiophene	134	C8H6S	000095-15-8	91



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG122615\
Data File : BG020564.D
Acq On : 27 Dec 2015 21:13
Operator : UM/SJ
Sample : G4846-11
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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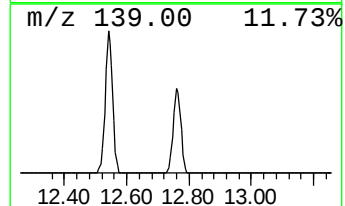
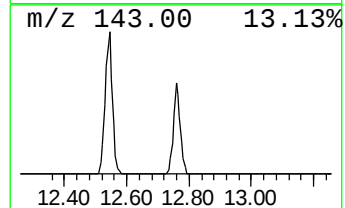
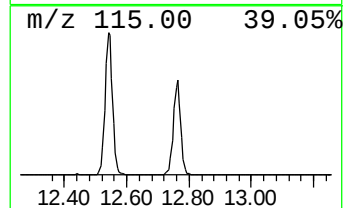
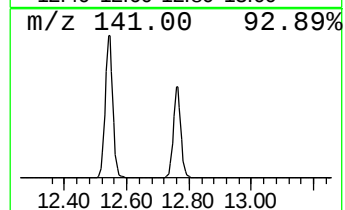
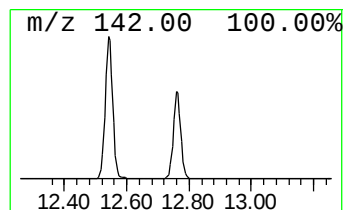
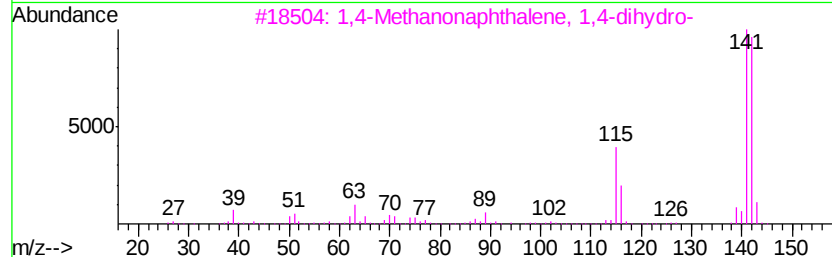
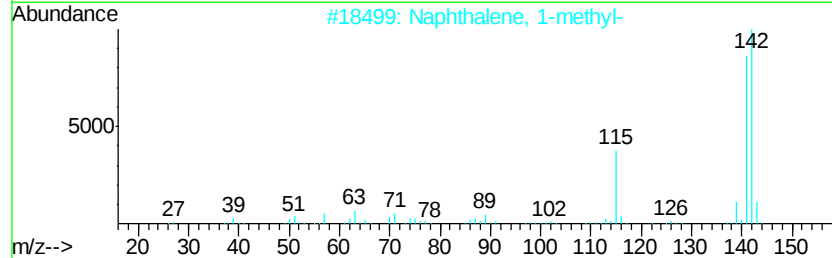
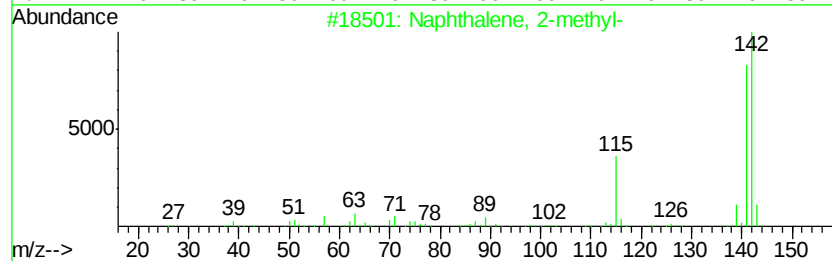
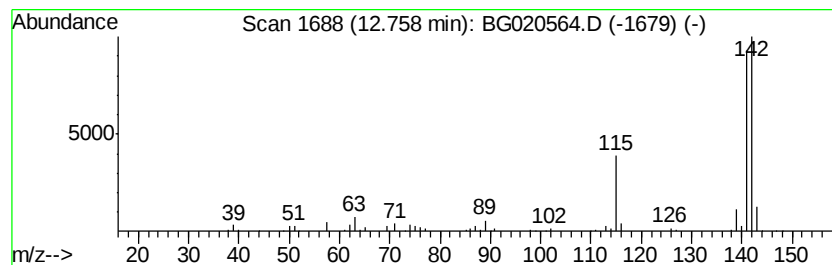
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	22.83 ng/ul	207052	Naphthalene-d8	10.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
4		Benzocycloheptatriene	142	C11H10	000264-09-5	94
5		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG122615\
Data File : BG020564.D
Acq On : 27 Dec 2015 21:13
Operator : UM/SJ
Sample : G4846-11
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
D9N57

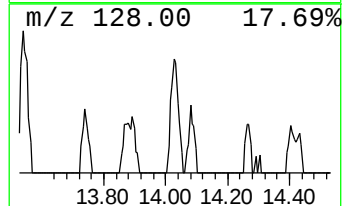
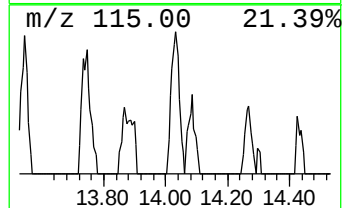
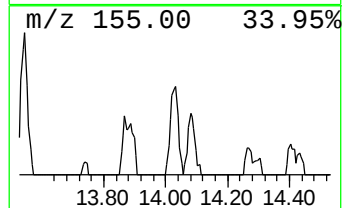
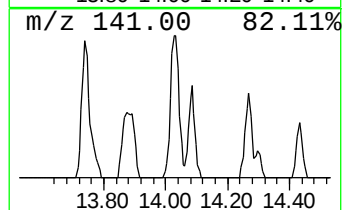
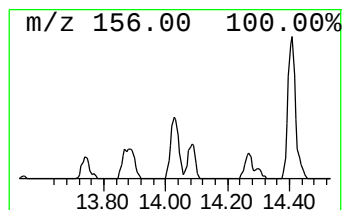
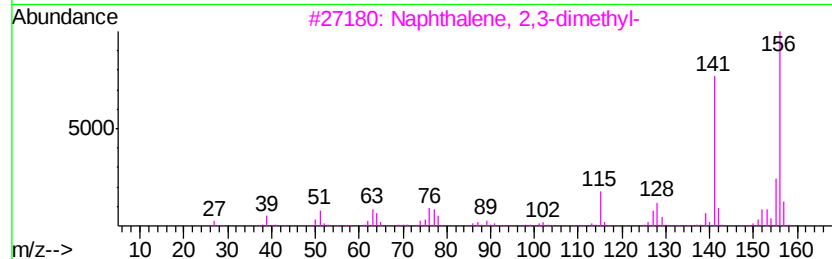
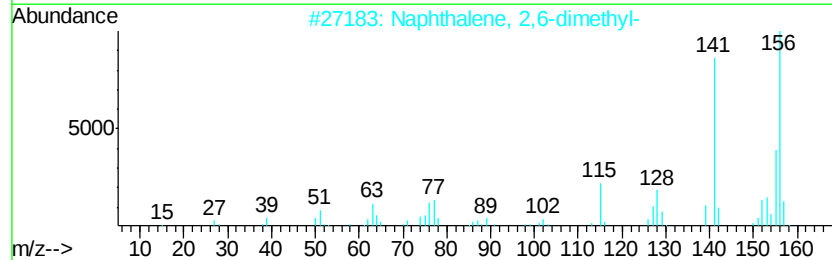
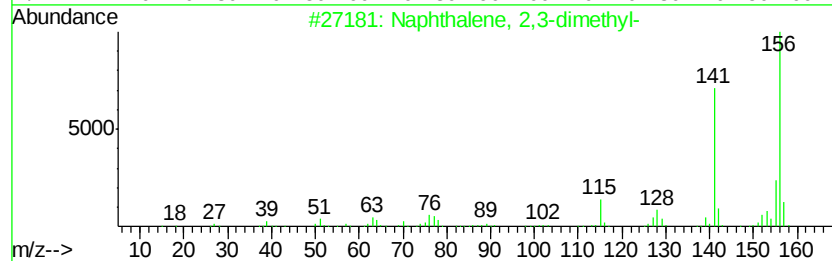
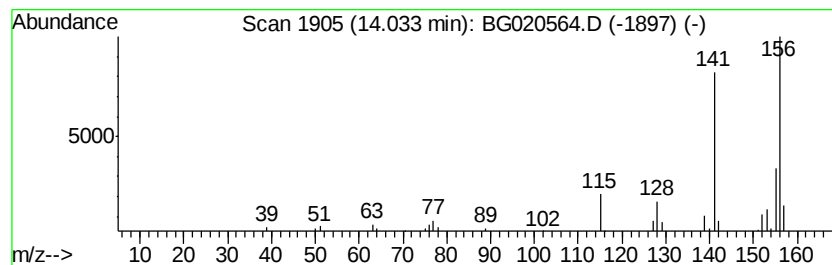
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Naphthalene, 2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.03	2.50 ng/ul	39567	Acenaphthene-d10	14.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
5		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG122615\
Data File : BG020564.D
Acq On : 27 Dec 2015 21:13
Operator : UM/SJ
Sample : G4846-11
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
D9N57

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG122415.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...	3.18	6.8	ng/ul	38383	1	8.08	112053	20.0
Benzene, 1-ethyl-...	7.72	3.2	ng/ul	17734	1	8.08	112053	20.0
Indane	8.45	6.6	ng/ul	36844	1	8.08	112053	20.0
Benzene, 1,2-prop...	8.62	4.2	ng/ul	23267	1	8.08	112053	20.0
unknown-01	10.29	2.2	ng/ul	19663	2	10.90	181372	20.0
Cyclopenta[c]thia...	11.07	3.2	ng/ul	29338	2	10.90	181372	20.0
Naphthalene, 1-me...	12.76	22.8	ng/ul	207052	2	10.90	181372	20.0
Naphthalene, 2,3-...	14.03	2.5	ng/ul	39567	3	14.71	316443	20.0