

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122215\
 Data File : BG020462.D
 Acq On : 24 Dec 2015 3:21
 Operator : UM/SJ
 Sample : G4848-04MEDL 20X
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 D9N50MEDL

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-BG121415.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	3.108	41	46	51	rBV3	8448	14846	2.55%	0.334%
2	3.184	56	59	65	rVB4	4563	7951	1.36%	0.179%
3	4.224	231	236	242	rBV	11221	18265	3.13%	0.411%
4	7.409	774	778	783	rBB	1023	1863	0.32%	0.042%
5	7.614	809	813	819	rVB	1374	1816	0.31%	0.041%
6	8.085	887	893	901	rBV	50748	87733	15.05%	1.976%
7	8.625	981	985	990	rBB	2183	3904	0.67%	0.088%
8	9.265	1086	1094	1097	rBB	1196	1940	0.33%	0.044%
9	10.523	1304	1308	1314	rBB	1608	2409	0.41%	0.054%
10	10.899	1365	1372	1376	rBV	75463	135056	23.17%	3.041%
11	10.952	1376	1381	1394	rVB	97545	175812	30.16%	3.959%
12	11.069	1399	1401	1409	rVB	2237	2675	0.46%	0.060%
13	12.550	1647	1653	1661	rBB	53316	86938	14.91%	1.958%
14	12.767	1682	1690	1697	rBB	25339	41320	7.09%	0.930%
15	13.549	1818	1823	1830	rBB	18147	27073	4.64%	0.610%
16	13.748	1852	1857	1864	rBB	5291	8317	1.43%	0.187%
17	13.889	1874	1881	1887	rVB4	7347	18095	3.10%	0.407%
18	14.036	1900	1906	1911	rBV	11403	19542	3.35%	0.440%
19	14.089	1911	1915	1917	rBV2	7214	9784	1.68%	0.220%
20	14.271	1942	1946	1949	rBV	4688	5888	1.01%	0.133%
21	14.307	1949	1952	1955	rVB2	1383	1895	0.33%	0.043%
22	14.412	1963	1970	1972	rBV2	6034	9614	1.65%	0.217%
23	14.436	1972	1974	1980	rVB	5448	7395	1.27%	0.167%
24	14.718	2012	2022	2027	rBV2	137521	228364	39.17%	5.143%
25	14.783	2027	2033	2040	rVV	119181	184609	31.67%	4.157%
26	14.847	2040	2044	2051	rVB2	3706	5199	0.89%	0.117%
27	14.918	2052	2056	2060	rBV2	1638	2529	0.43%	0.057%
28	15.023	2068	2074	2078	rBB	3894	5392	0.92%	0.121%
29	15.112	2083	2089	2097	rBV	79813	121510	20.84%	2.736%
30	15.188	2097	2102	2106	rVB2	2335	3554	0.61%	0.080%
31	15.323	2121	2125	2129	rBV2	1532	2577	0.44%	0.058%
32	15.376	2131	2134	2138	rVB	1430	2007	0.34%	0.045%
33	15.705	2186	2190	2194	rVV2	9536	15068	2.58%	0.339%
34	15.758	2194	2199	2208	rVV	113437	170451	29.24%	3.838%

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35	15.858	2212	2216	2218	rVV2	3961	6361	1.09%	0.143%
36	15.881	2218	2220	2223	rVV	6395	8439	1.45%	0.190%
37	15.922	2223	2227	2232	rVV3	13515	19100	3.28%	0.430%
38	15.969	2232	2235	2237	rVV2	2235	2822	0.48%	0.064%
39	16.011	2238	2242	2245	rVV2	5589	8705	1.49%	0.196%
40	16.052	2245	2249	2256	rVB	12876	19821	3.40%	0.446%
41	16.199	2268	2274	2283	rVB2	15734	29997	5.15%	0.676%
42	16.293	2285	2290	2294	rVB2	2097	3385	0.58%	0.076%
43	16.551	2330	2334	2339	rVB	6893	10171	1.74%	0.229%
44	16.763	2360	2370	2375	rBV2	10169	21735	3.73%	0.489%
45	16.815	2375	2379	2384	rVB2	3690	5639	0.97%	0.127%
46	16.915	2391	2396	2401	rVB5	3962	6200	1.06%	0.140%
47	16.986	2401	2408	2411	rBV4	3517	7287	1.25%	0.164%
48	17.027	2411	2415	2419	rVV3	3580	5081	0.87%	0.114%
49	17.186	2437	2442	2449	rBV3	4104	9339	1.60%	0.210%
50	17.280	2453	2458	2464	rVB2	22567	33081	5.67%	0.745%
51	17.462	2482	2489	2492	rBV	194511	290190	49.78%	6.535%
52	17.503	2492	2496	2502	rVV	388800	582949	100.00%	13.128%
53	17.556	2502	2505	2508	rVV2	13922	21120	3.62%	0.476%
54	17.597	2508	2512	2516	rVB	22199	31554	5.41%	0.711%
55	17.861	2550	2557	2568	rBV2	12720	23370	4.01%	0.526%
56	17.979	2572	2577	2583	rVV3	4863	6819	1.17%	0.154%
57	18.043	2584	2588	2595	rVB3	1788	3093	0.53%	0.070%
58	18.184	2608	2612	2616	rBV	1934	3260	0.56%	0.073%
59	18.337	2632	2638	2641	rBV	22750	32486	5.57%	0.732%
60	18.384	2641	2646	2653	rVB	32189	44832	7.69%	1.010%
61	18.466	2655	2660	2664	rBV	7289	10831	1.86%	0.244%
62	18.531	2664	2671	2676	rBV3	42986	77198	13.24%	1.738%
63	18.772	2709	2712	2717	rVV3	1680	2033	0.35%	0.046%
64	18.825	2717	2721	2726	rVV	16781	22843	3.92%	0.514%
65	18.872	2726	2729	2734	rVB	1963	2905	0.50%	0.065%
66	19.072	2759	2763	2768	rVB2	2563	3973	0.68%	0.089%
67	19.125	2768	2772	2775	rBV	2366	3152	0.54%	0.071%
68	19.160	2775	2778	2782	rVB3	1899	2334	0.40%	0.053%
69	19.242	2786	2792	2798	rBV2	5361	9343	1.60%	0.210%
70	19.289	2798	2800	2801	rVV2	1744	1778	0.31%	0.040%
71	19.307	2801	2803	2811	rVB3	3577	6814	1.17%	0.153%

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72	19.377	2811	2815	2818	rBV3	1555	2592	0.44%	0.058%
73	19.407	2818	2820	2825	rVB	1636	1932	0.33%	0.044%
74	19.506	2831	2837	2845	rVV	234863	328905	56.42%	7.407%
75	19.565	2845	2847	2850	rVV2	1910	2237	0.38%	0.050%
76	19.600	2850	2853	2858	rVB2	2183	2719	0.47%	0.061%
77	19.753	2874	2879	2884	rVV2	4543	7175	1.23%	0.162%
78	19.841	2888	2894	2895	rBV2	49375	67061	11.50%	1.510%
79	19.871	2895	2899	2906	rVV	151451	218830	37.54%	4.928%
80	19.935	2907	2910	2915	rVB3	6402	9216	1.58%	0.208%
81	20.047	2924	2929	2934	rVB	7342	10389	1.78%	0.234%
82	20.112	2936	2940	2944	rVB2	1726	2541	0.44%	0.057%
83	20.188	2947	2953	2960	rBV6	6094	15317	2.63%	0.345%
84	20.247	2960	2963	2968	rVB2	3058	3739	0.64%	0.084%
85	20.358	2973	2982	2989	rVB	23458	41325	7.09%	0.931%
86	20.458	2994	2999	3003	rBV	28345	36071	6.19%	0.812%
87	20.517	3003	3009	3013	rVV5	5428	11108	1.91%	0.250%
88	20.552	3013	3015	3019	rVB3	3471	3701	0.63%	0.083%
89	20.652	3029	3032	3037	rBV3	3049	5599	0.96%	0.126%
90	20.699	3038	3040	3047	rVB4	3559	5072	0.87%	0.114%
91	20.881	3068	3071	3076	rVB2	3047	4260	0.73%	0.096%
92	21.081	3102	3105	3108	rBV2	1980	2817	0.48%	0.063%
93	21.310	3141	3144	3148	rBV3	5207	6009	1.03%	0.135%
94	21.357	3148	3152	3156	rVB3	5579	6762	1.16%	0.152%
95	21.404	3156	3160	3164	rBV3	4391	7788	1.34%	0.175%
96	21.733	3207	3216	3221	rBV	250693	442876	75.97%	9.973%
97	21.939	3247	3251	3258	rVB2	6409	10713	1.84%	0.241%
98	23.966	3589	3596	3601	rBV4	7023	16280	2.79%	0.367%
99	24.777	3727	3734	3739	rBV4	5362	12827	2.20%	0.289%
100	25.000	3762	3772	3784	rBV2	117565	343347	58.90%	7.732%

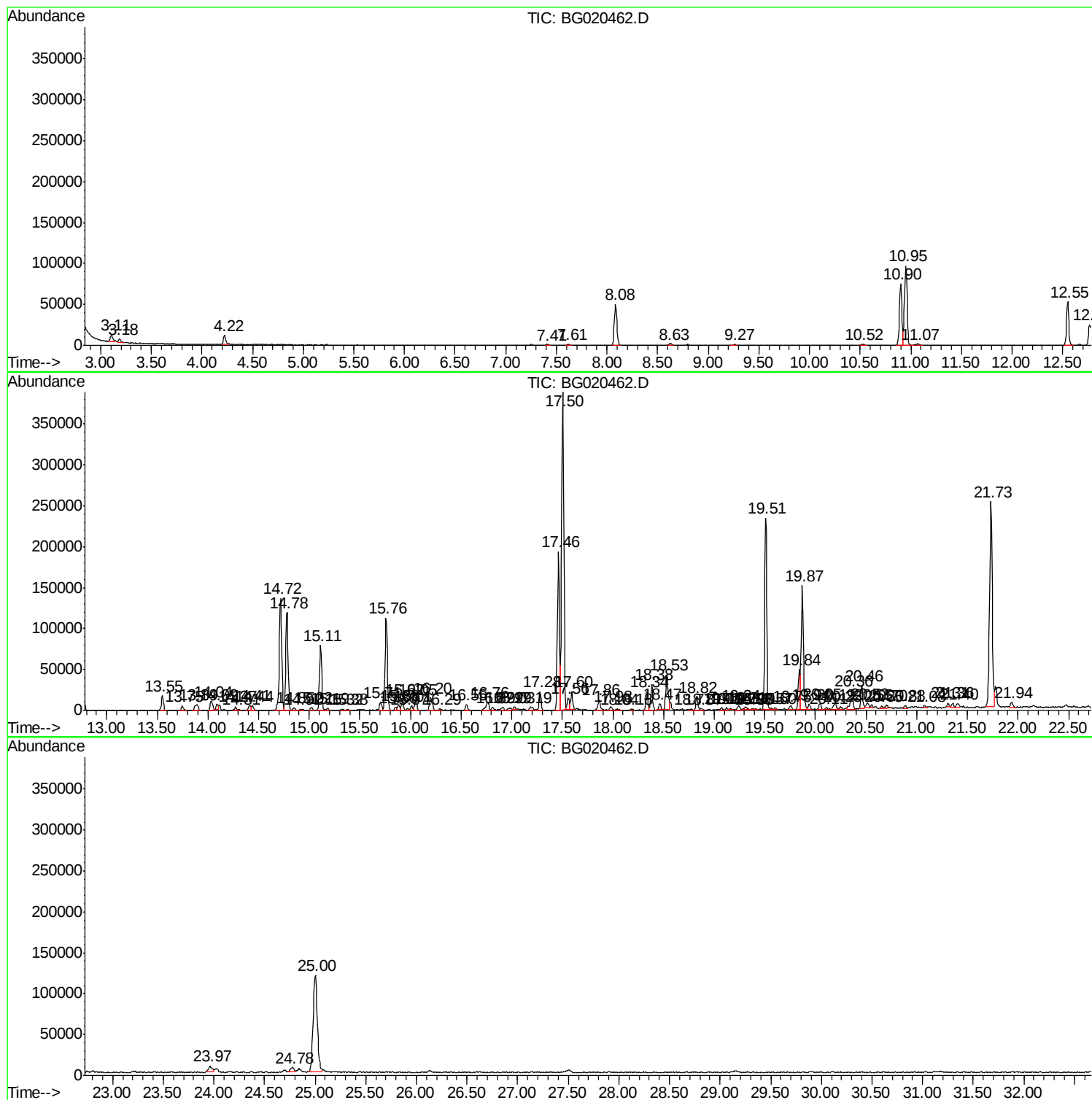
Sum of corrected areas: 4440639

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

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



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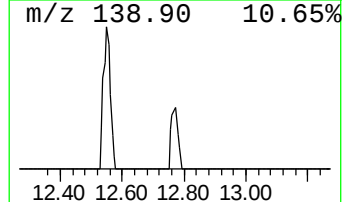
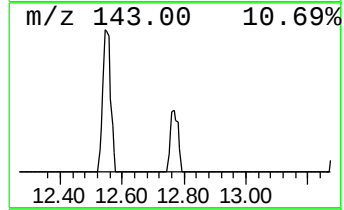
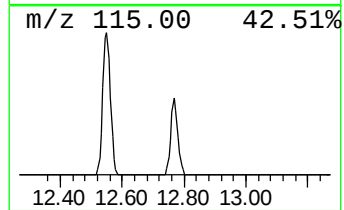
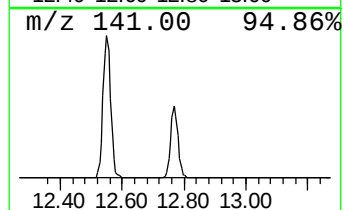
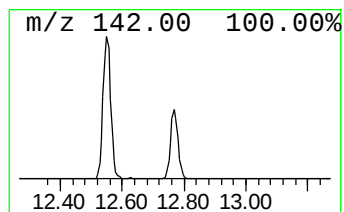
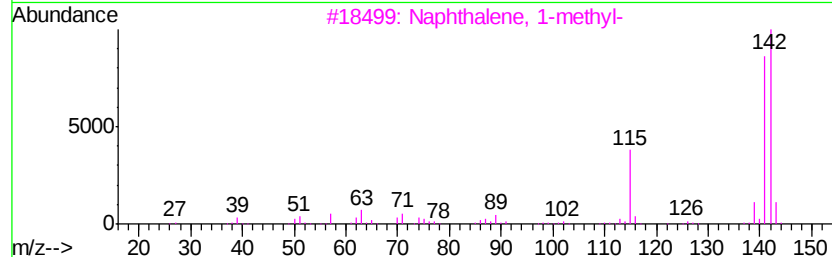
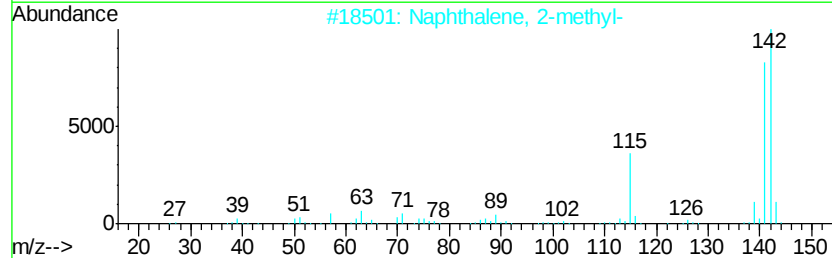
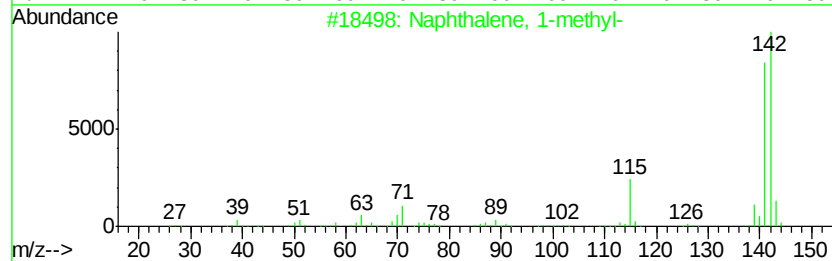
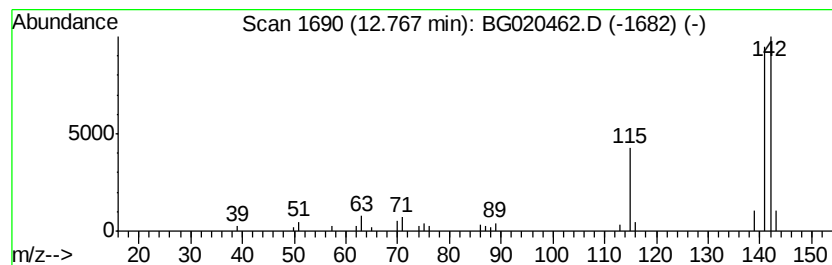
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	6.12 ng/ul	41320	Naphthalene-d8	10.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	91
3		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5		Benzocycloheptatriene	142	C11H10	000264-09-5	90



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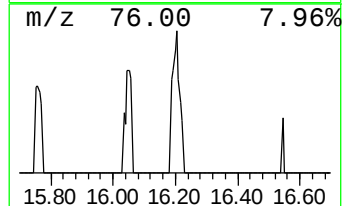
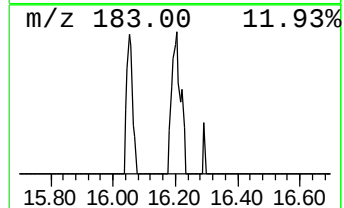
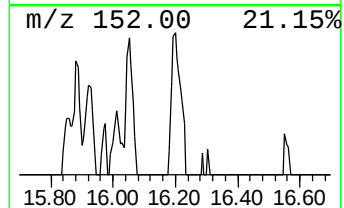
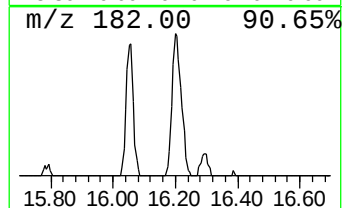
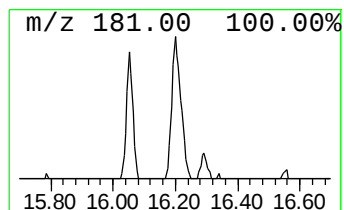
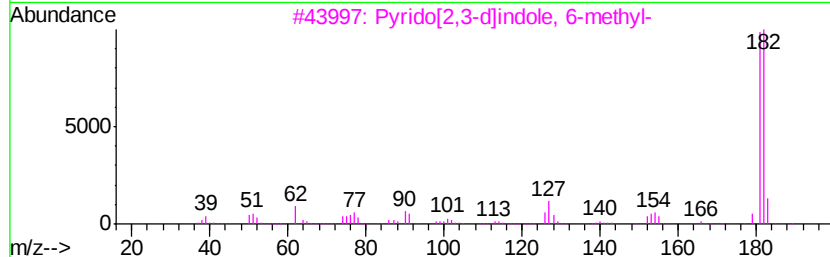
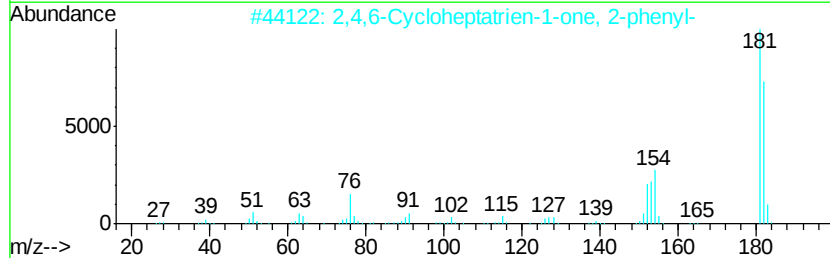
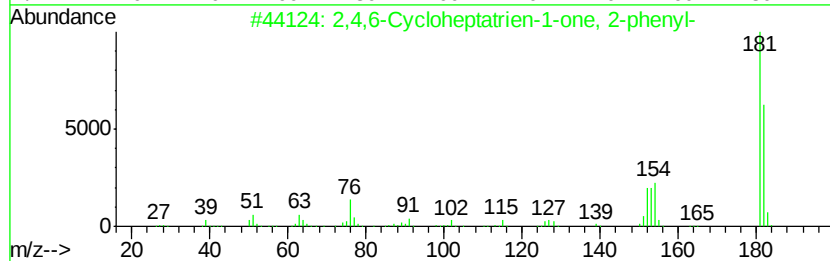
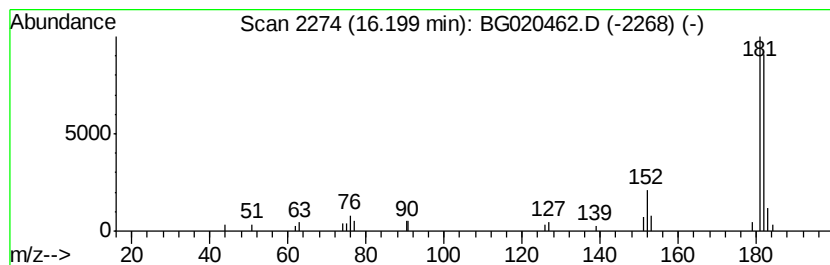
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2,4,6-Cycloheptatrien-1-one... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.20	2.07 ng/ul	29997	Phenanthrene-d10	17.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	78
2		2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	78
3		Pyrido[2,3-d]indole, 6-methyl-	182	C12H10N2	108349-67-3	72
4		9H-Fluoren-9-ol	182	C13H10O	001689-64-1	72
5		9H-Fluoren-9-ol	182	C13H10O	001689-64-1	64



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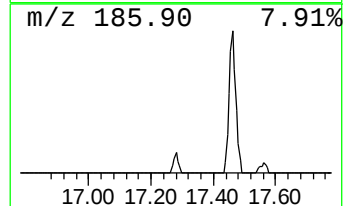
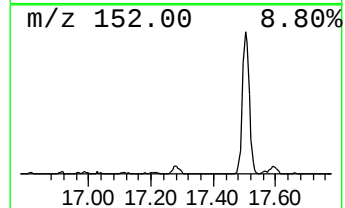
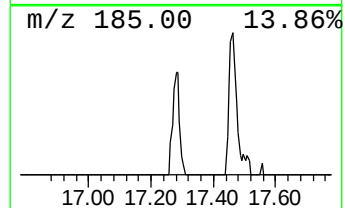
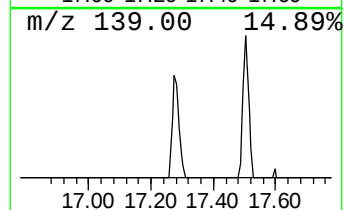
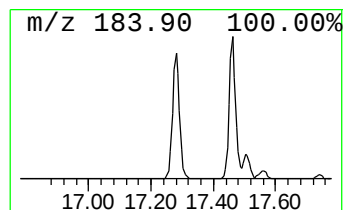
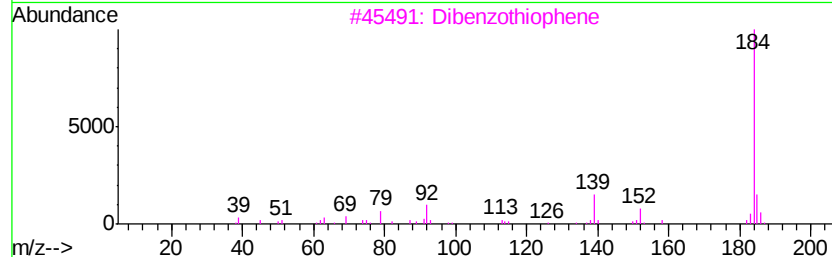
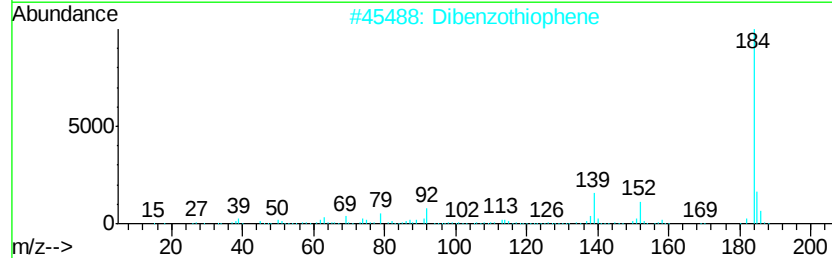
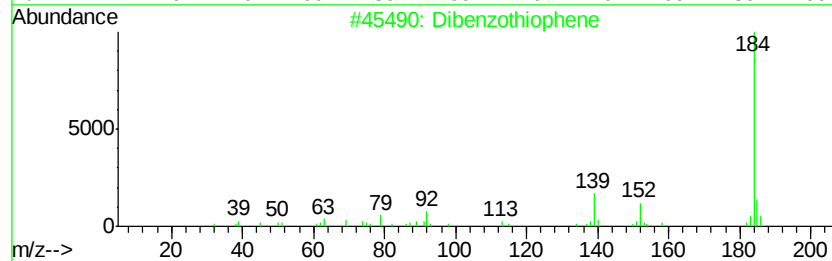
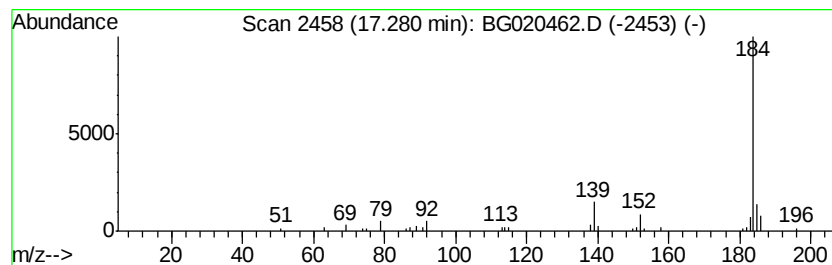
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TIC Integration Parameters: LSCINT.P

Peak Number 5 Dibenzothiophene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.28	2.28 ng/ul	33081	Phenanthrene-d10	17.46

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzothiophene		184	C12H8S	000132-65-0	95
2	Dibenzothiophene		184	C12H8S	000132-65-0	95
3	Dibenzothiophene		184	C12H8S	000132-65-0	95
4	Dibenzothiophene		184	C12H8S	000132-65-0	94
5	Azuleno(2,1-b)thiophene		184	C12H8S	000248-13-5	91



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122215\
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Acq On : 24 Dec 2015 3:21
Operator : UM/SJ
Sample : G4848-04MEDL 20X
Misc :
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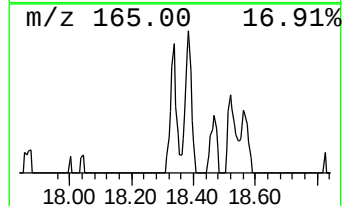
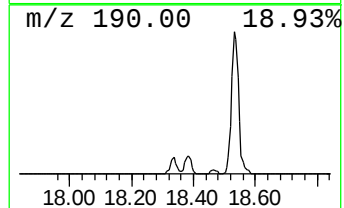
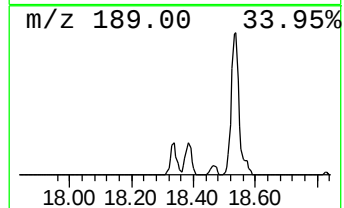
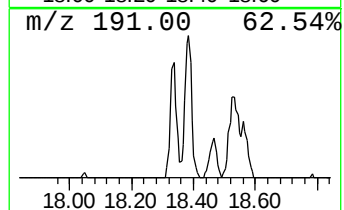
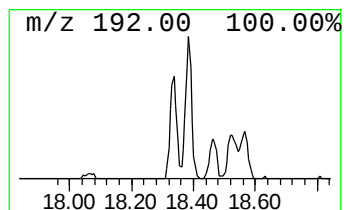
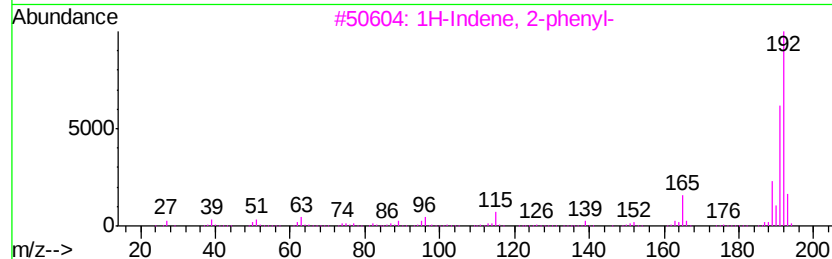
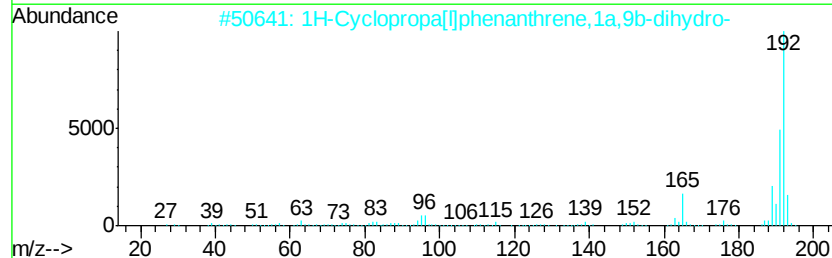
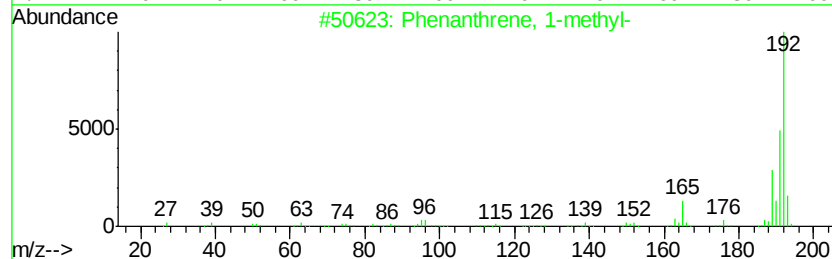
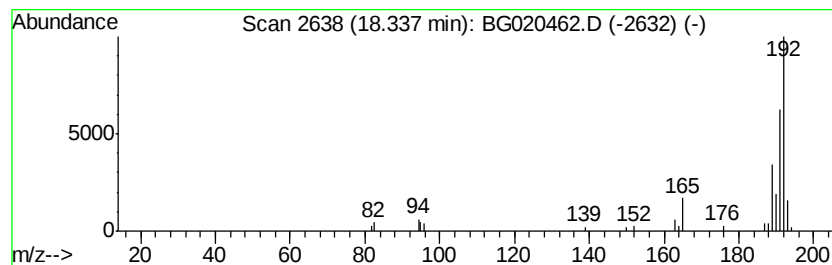
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Phenanthrene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.34	2.24 ng/ul	32486	Phenanthrene-d10	17.46	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
2	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	93
3	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	93
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	90
5	Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	90



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122215\
Data File : BG020462.D
Acq On : 24 Dec 2015 3:21
Operator : UM/SJ
Sample : G4848-04MEDL 20X
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
D9N50MEDL

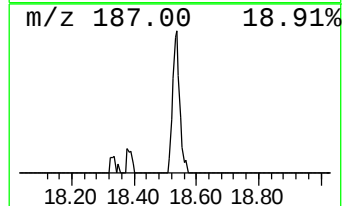
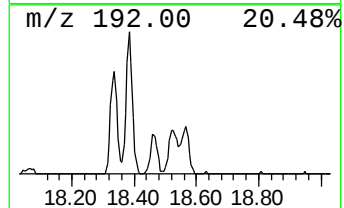
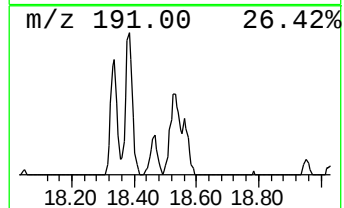
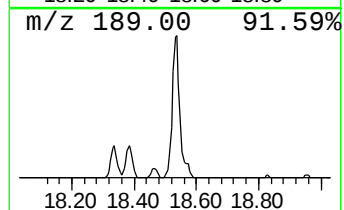
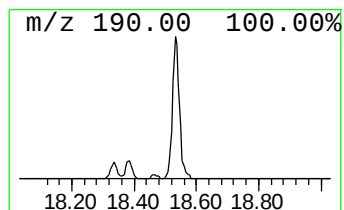
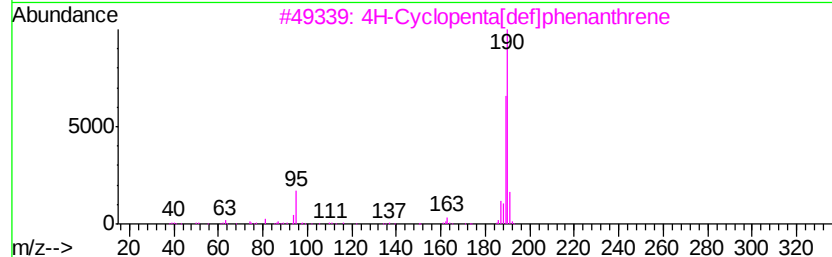
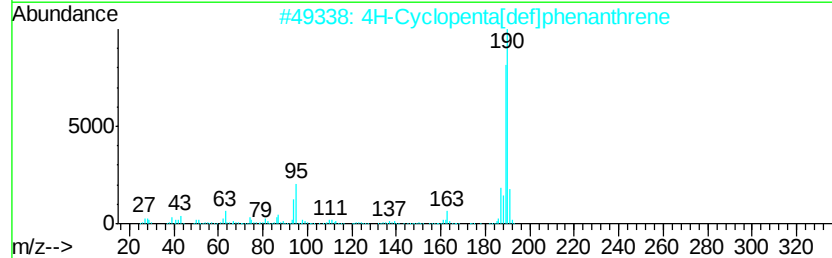
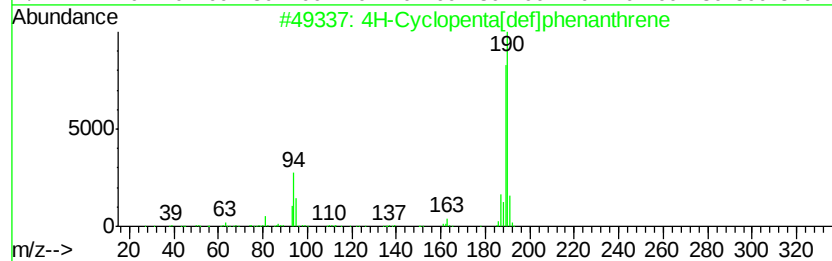
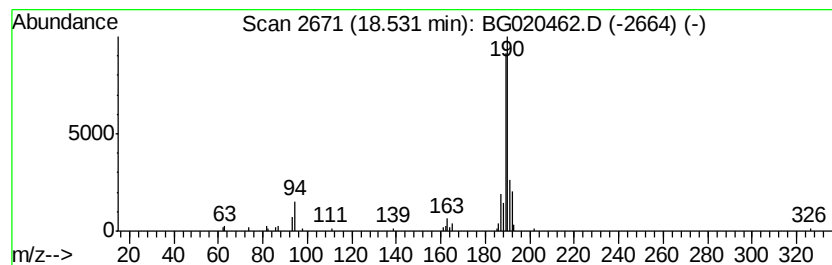
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG121415.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.53	5.32 ng/ul	77198	Phenanthrene-d10	17.46

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	83
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
3		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	74
4		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	42
5		Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	42



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG122215\
Data File : BG020462.D
Acq On : 24 Dec 2015 3:21
Operator : UM/SJ
Sample : G4848-04MEDL 20X
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
D9N50MEDL

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG121415.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
Naphthalene, 1-me...	12.77	6.1	ng/ul	41320	2	10.90	135056	20.0
2,4,6-Cycloheptat...	16.20	2.1	ng/ul	29997	4	17.46	290190	20.0
Dibenzothiophene	17.28	2.3	ng/ul	33081	4	17.46	290190	20.0
Phenanthrene, 1-m...	18.34	2.2	ng/ul	32486	4	17.46	290190	20.0
4H-Cyclopenta[def...	18.53	5.3	ng/ul	77198	4	17.46	290190	20.0