LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
Data File : BG022240.D
Acq On : 8 Jun 2016 18:10
Operator : UM/SJ
Sample : H3402-03 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DUP-3

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.984	19	25	43	rVB	482102	777110	65.22%	4.571%
2	5.182	394	399	410	rVB	20806	39151	3.29%	0.230%
3	5.664	473	481	497	rBV	213286	417948	35.08%	2.458%
4	7.274	746	755	770	rBV	251053	516844	43.38%	3.040%
5	7.644	810	818	833	rBV	272660	545250	45.76%	3.207%
6	8.114	891	898	907	rBV	104928	203160	17.05%	1.195%
7	8.443	943	954	963	rBV	213051	424039	35.59%	2.494%
8	9.289	1088	1098	1114	rBV	142867	313570	26.32%	1.844%
9	10.934	1371	1378	1385	rBV	185387	375392	31.51%	2.208%
10	13.366	1783	1792	1803	rBV	604621	1011542	84.90%	5.949%
11	14.066	1904	1911	1916	rBV5	6400	12104	1.02%	0.071%
12	14.189	1927	1932	1938	rBV	28244	44240	3.71%	0.260%
13	14.747	2018	2027	2033	rBV2	324573	562238	47.19%	3.307%
14	14.806	2033	2037	2045	rVB	38145	67715	5.68%	0.398%
15	15.141	2085	2094	2100	rBV2	15678	30323	2.54%	0.178%
16	15.558	2162	2165	2171	rVB3	8620	12815	1.08%	0.075%
17	15.787	2198	2204	2215	rVB	33253	62561	5.25%	0.368%
18	16.081	2249	2254	2262	rVB5	9292	20907	1.75%	0.123%
19	16.234	2274	2280	2290	rBV	333100	584956	49.09%	3.440%
20	16.416	2308	2311	2314	rBV3	9184	13638	1.14%	0.080%
21	16.792	2363	2375	2379	rBV7	11800	32457	2.72%	0.191%
22	17.015	2406	2413	2415	rBV8	7686	16030	1.35%	0.094%
23	17.056	2417	2420	2428	rVB6	8792	17613	1.48%	0.104%
24	17.150	2430	2436	2441	rVB3	11814	19745	1.66%	0.116%
25	17.215	2441	2447	2454	rBV7	17142	38548	3.24%	0.227%
26	17.309	2459	2463	2473	rVB	19706	36753	3.08%	0.216%
27	17.491	2487	2494	2497	rBV2	349677	585456	49.14%	3.443%
28	17.532	2497	2501	2509	rVV	375143	636942	53.46%	3.746%
29	17.620	2511	2516	2522	rVB	81041	135383	11.36%	0.796%
30	17.896	2554	2563	2569	rBV3	28469	72296	6.07%	0.425%
31	17.943	2569	2571	2577	rVB5	9076	14627	1.23%	0.086%
32	17.996	2577	2580	2583	rBV2	11366	15462	1.30%	0.091%
33	18.067	2589	2592	2599	rVB6	10121	15539	1.30%	0.091%
34	18.361	2636	2642	2646	rBV	51132	85449	7.17%	0.503%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
Data File : BG022240.D
Acq On : 8 Jun 2016 18:10
Operator : UM/SJ
Sample : H3402-03 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DUP-3

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	18.413	2646	2651	2656	rVB	65207	111550	9.36%	0.656%
36	18.496	2658	2665	2669	rBV	24579	45162	3.79%	0.266%
37	18.560	2669	2676	2679	rBV2	78784	157604	13.23%	0.927%
38	18.854	2722	2726	2731	rBV	39783	63256	5.31%	0.372%
39	18.901	2731	2734	2739	rVB3	21146	32538	2.73%	0.191%
40	19.095	2761	2767	2772	rBV6	17500	29201	2.45%	0.172%
41	19.148	2772	2776	2780	rBV3	11093	17720	1.49%	0.104%
42	19.189	2780	2783	2788	rVB4	12200	15751	1.32%	0.093%
43	19.265	2790	2796	2802	rBV	40428	77133	6.47%	0.454%
44	19.324	2802	2806	2811	rBV4	12620	24455	2.05%	0.144%
45	19.412	2816	2821	2828	rBV4	33823	72947	6.12%	0.429%
46	19.536	2835	2842	2848	rBV	654340	1038565	87.16%	6.108%
47	19.630	2854	2858	2863	rVB4	18780	27663	2.32%	0.163%
48	19.683	2863	2867	2870	rBV3	9196	16623	1.40%	0.098%
49	19.777	2877	2883	2890	rVB4	20895	49119	4.12%	0.289%
50	19.894	2891	2903	2910	rVV	582914	1089707	91.46%	6.409%
51	19.959	2911	2914	2921	rVB	34572	53489	4.49%	0.315%
52	20.082	2929	2935	2941	rBV	850589	1180439	99.07%	6.943%
53	20.217	2952	2958	2964	rVB3	42556	95845	8.04%	0.564%
54	20.270	2965	2967	2974	rVB5	9495	20513	1.72%	0.121%
55	20.382	2974	2986	2992	rBV2	78038	215246	18.06%	1.266%
56	20.482	2998	3003	3007	rBV	57645	97802	8.21%	0.575%
57	20.540	3007	3013	3017	rVV2	45062	83803	7.03%	0.493%
58	20.681	3033	3037	3040	rBV	33420	43811	3.68%	0.258%
59	20.722	3041	3044	3051	rVB4	32349	52782	4.43%	0.310%
60	21.151	3112	3117	3125	rVB3	42503	83928	7.04%	0.494%
61	21.339	3145	3149	3152	rBV2	25507	44112	3.70%	0.259%
62	21.375	3153	3155	3160	rVB	45212	60613	5.09%	0.356%
63	21.428	3160	3164	3170	rBV3	37228	65202	5.47%	0.383%
64	21.762	3212	3221	3226	rBV2	432874	1191518	100.00%	7.008%
65	21.809	3226	3229	3240	rVB	235000	450925	37.84%	2.652%
66	21.968	3251	3256	3263	rBV	42614	81852	6.87%	0.481%
67	24.019	3596	3605	3623	rVB3	165736	806061	67.65%	4.741%
68	24.271	3644	3648	3659	rVB2	31204	77565	6.51%	0.456%
69	24.753	3723	3730	3744	rVB3	95772	308805	25.92%	1.816%
70	24.906	3747	3756	3770	rVB2	133057	436945	36.67%	2.570%
71	25.064	3772	3783	3792	rBV	171840	587385	49.30%	3.455%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
Data File : BG022240.D
Acq On : 8 Jun 2016 18:10
Operator : UM/SJ
Sample : H3402-03 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DUP-3

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	28.831	4414	4424	4443	rVB4	47104	243234	20.41%	1.431%
73	30.012	4613	4625	4626	rBV3	32854	93658	7.86%	0.551%

Sum of corrected areas: 17002330

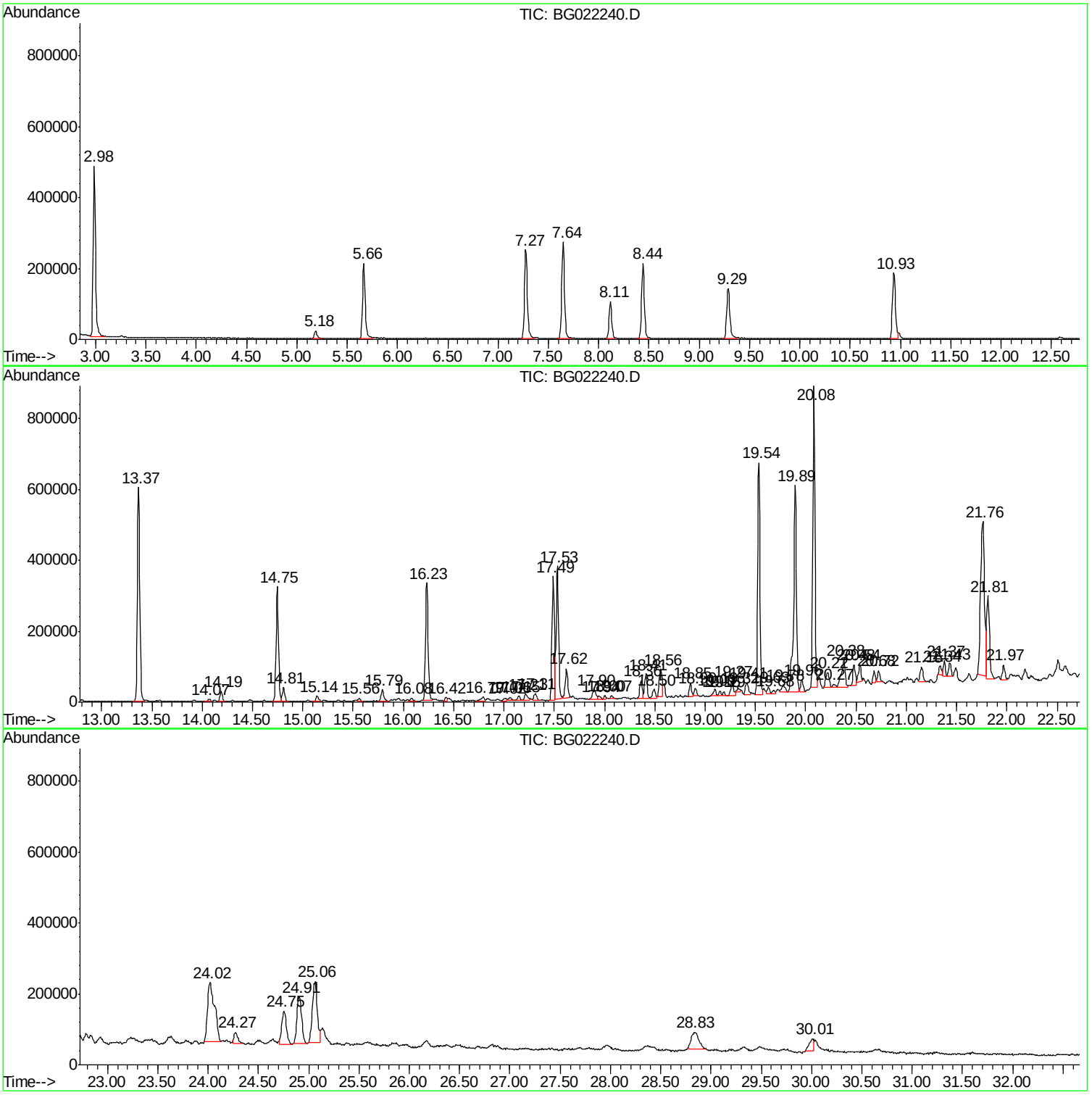
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
Data File : BG022240.D
Acq On : 8 Jun 2016 18:10
Operator : UM/SJ
Sample : H3402-03 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DUP-3

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
Data File : BG022240.D
Acq On : 8 Jun 2016 18:10
Operator : UM/SJ
Sample : H3402-03 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DUP-3

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

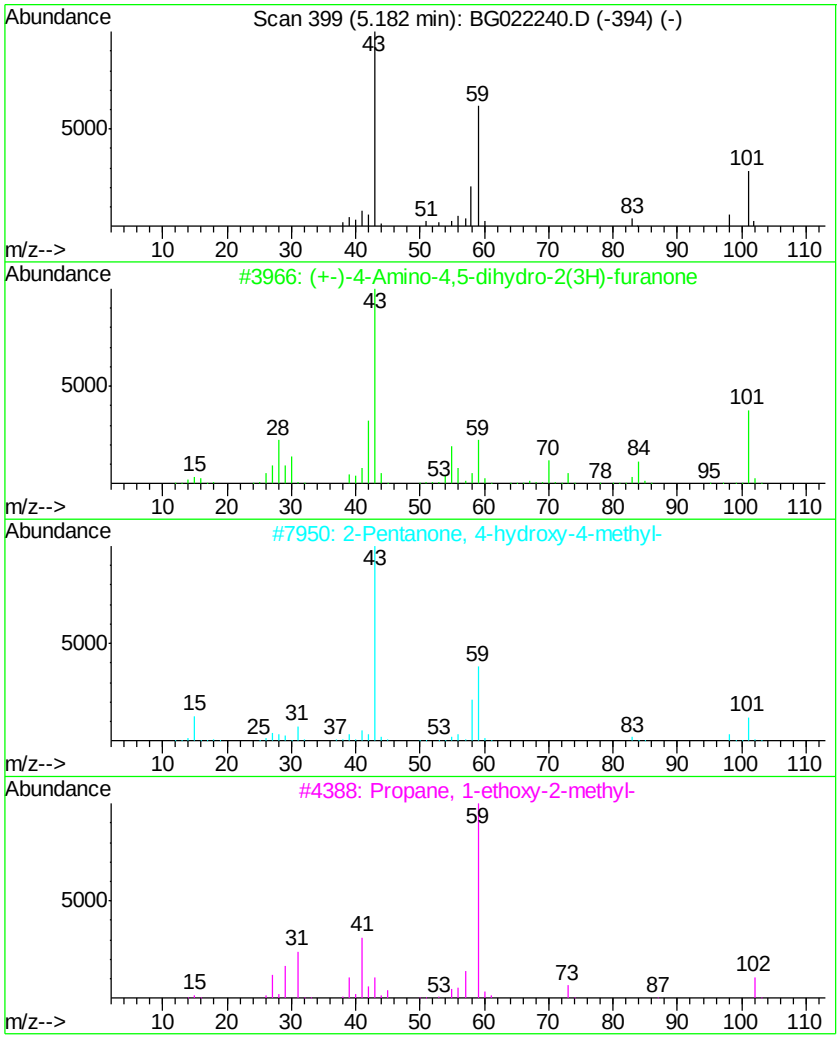
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

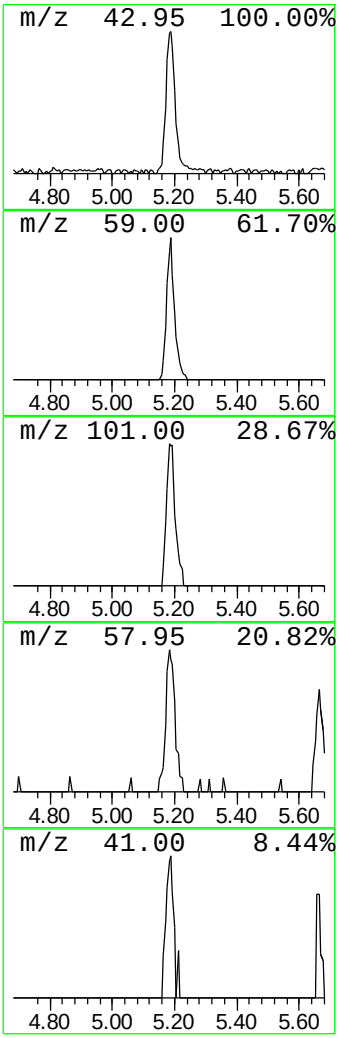
****

Peak Number 2 unknown5.18 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	3.85 ng	39151	1,4-Dichlorobenzene-d4	8.11

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+-)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	43
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
3	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	27
4	2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	23
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	12





Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
Data File : BG022240.D
Acq On : 8 Jun 2016 18:10
Operator : UM/SJ
Sample : H3402-03 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DUP-3

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

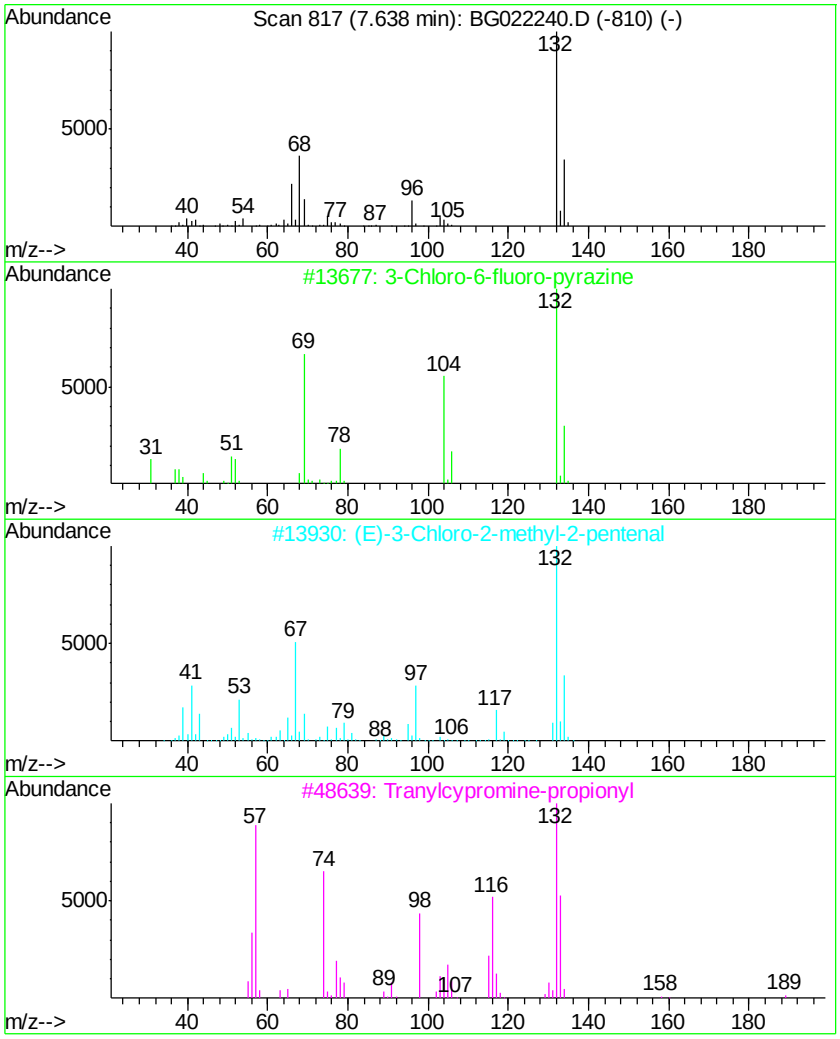
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

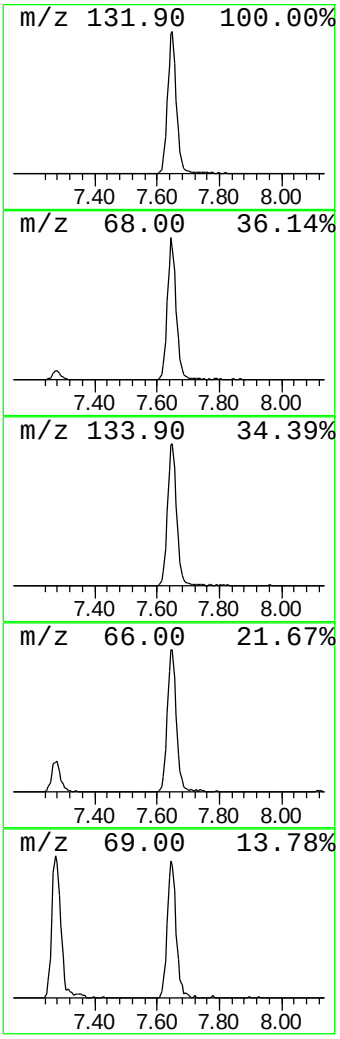
****

Peak Number 3 unknown7.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	53.68 ng	545250	1,4-Dichlorobenzene-d4	8.11

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4	Xenon	132	Xe	007440-63-3	9
5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9





Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
Data File : BG022240.D
Acq On : 8 Jun 2016 18:10
Operator : UM/SJ
Sample : H3402-03 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DUP-3

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

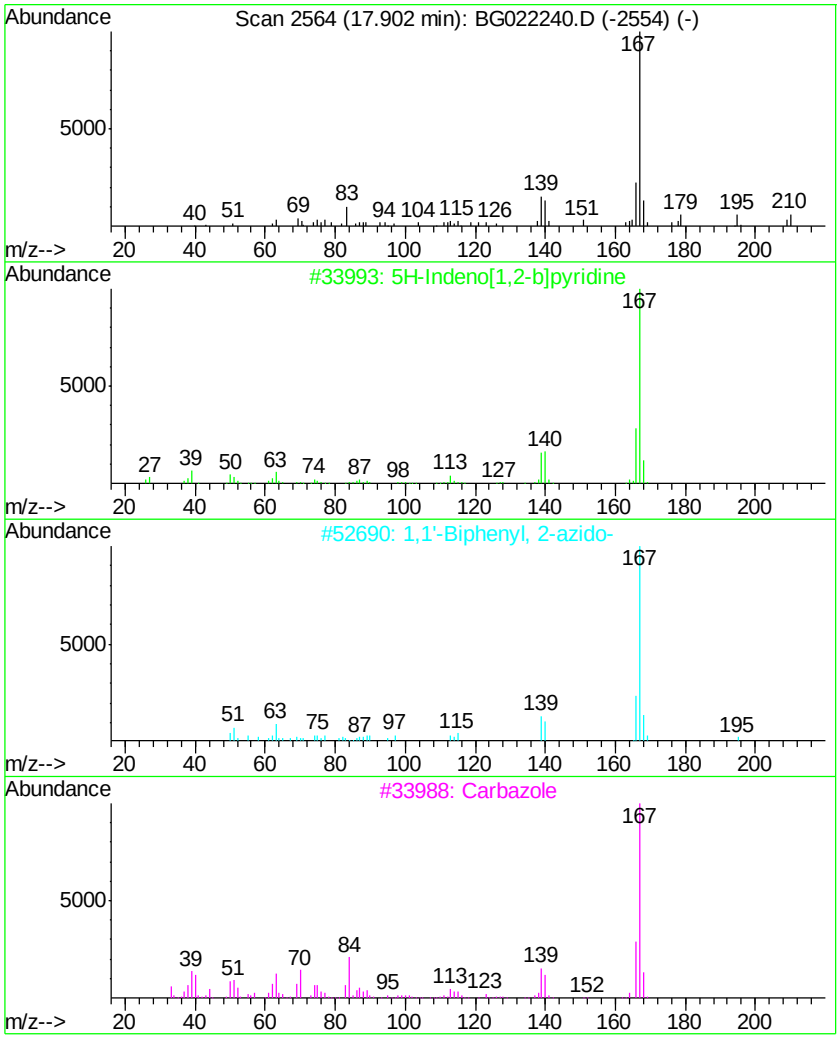
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

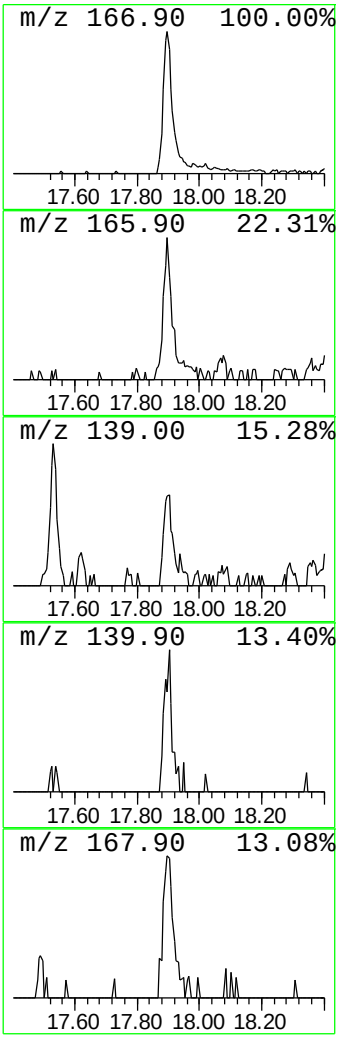
****

Peak Number 6 5H-Indeno[1,2-b]pyridine Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.90	2.47 ng	72296	Phenanthrene-d10	17.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	87
2	1,1'-Biphenyl, 2-azido-	195	C12H9N3	007599-23-7	87
3	Carbazole	167	C12H9N	000086-74-8	86
4	9H-Carbazole-9-methanol	197	C13H11NO	002409-36-1	72
5	2-Propanone, 1,1-diphenyl-	210	C15H14O	000781-35-1	70





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Sample : H3402-03 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DUP-3

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

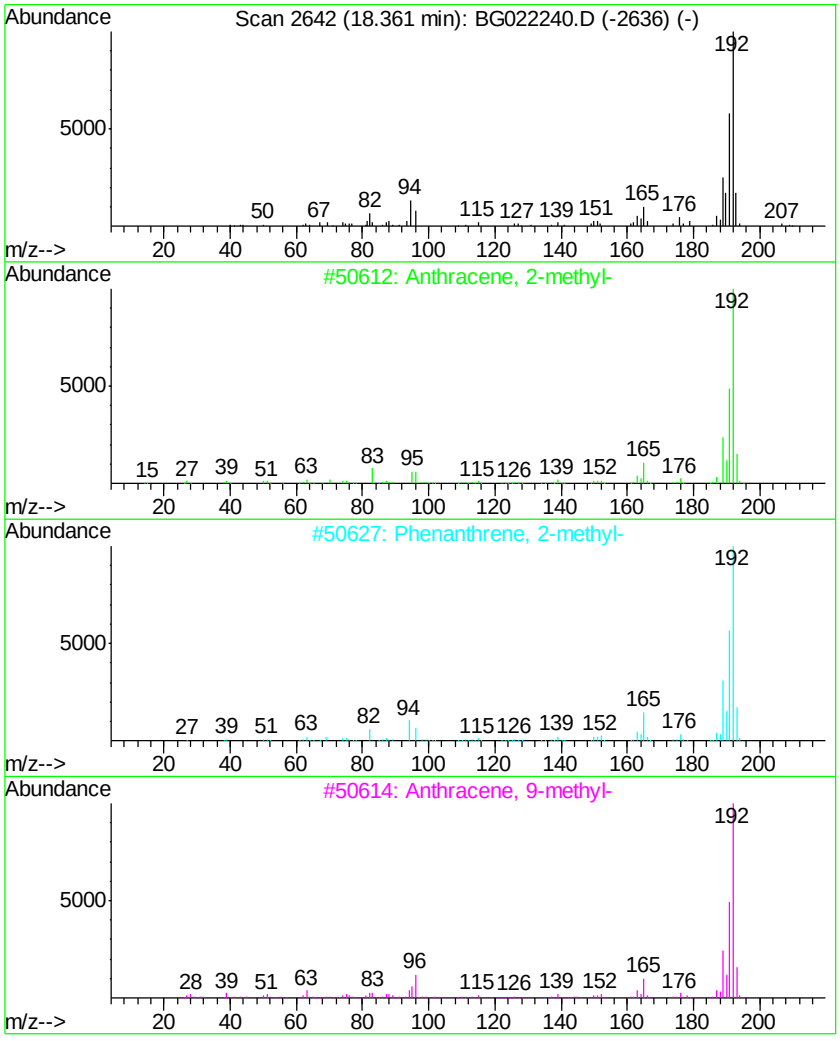
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

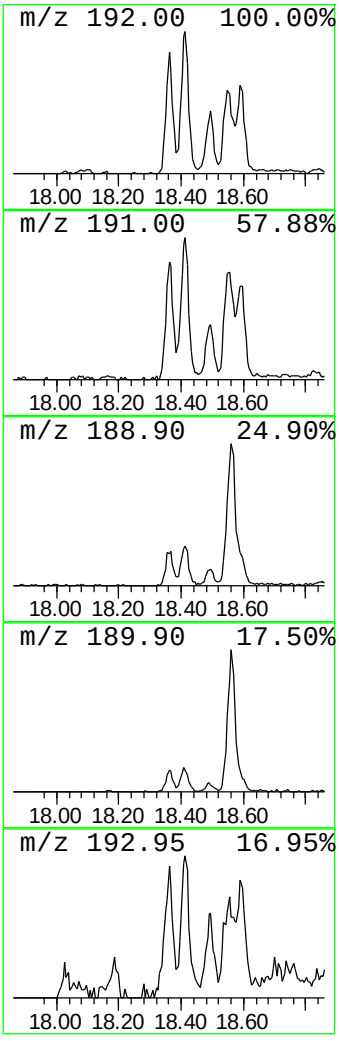
****

Peak Number 7 Anthracene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.36	2.92 ng	85449	Phenanthrene-d10	17.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90





Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
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Acq On : 8 Jun 2016 18:10
Operator : UM/SJ
Sample : H3402-03 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DUP-3

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

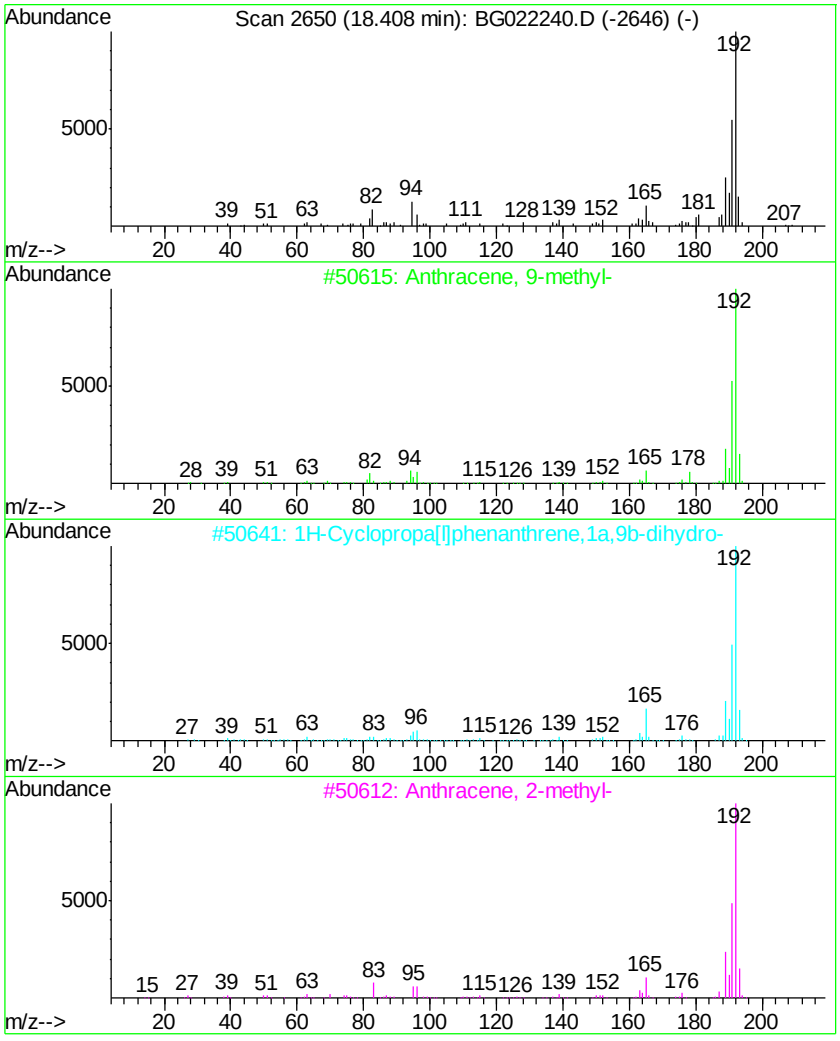
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

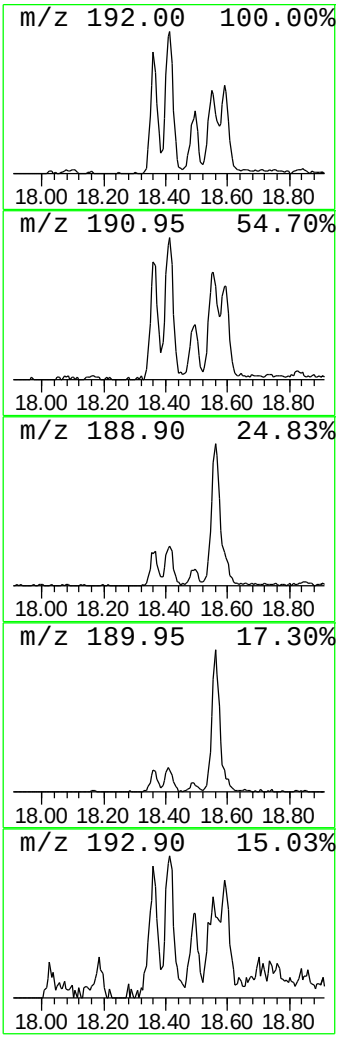
****

Peak Number 8 Anthracene, 9-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.41	3.81 ng	111550	Phenanthrene-d10	17.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 9-methyl-	192	C15H12	000779-02-2	94
2	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	93
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90





Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
Data File : BG022240.D
Acq On : 8 Jun 2016 18:10
Operator : UM/SJ
Sample : H3402-03 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DUP-3

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

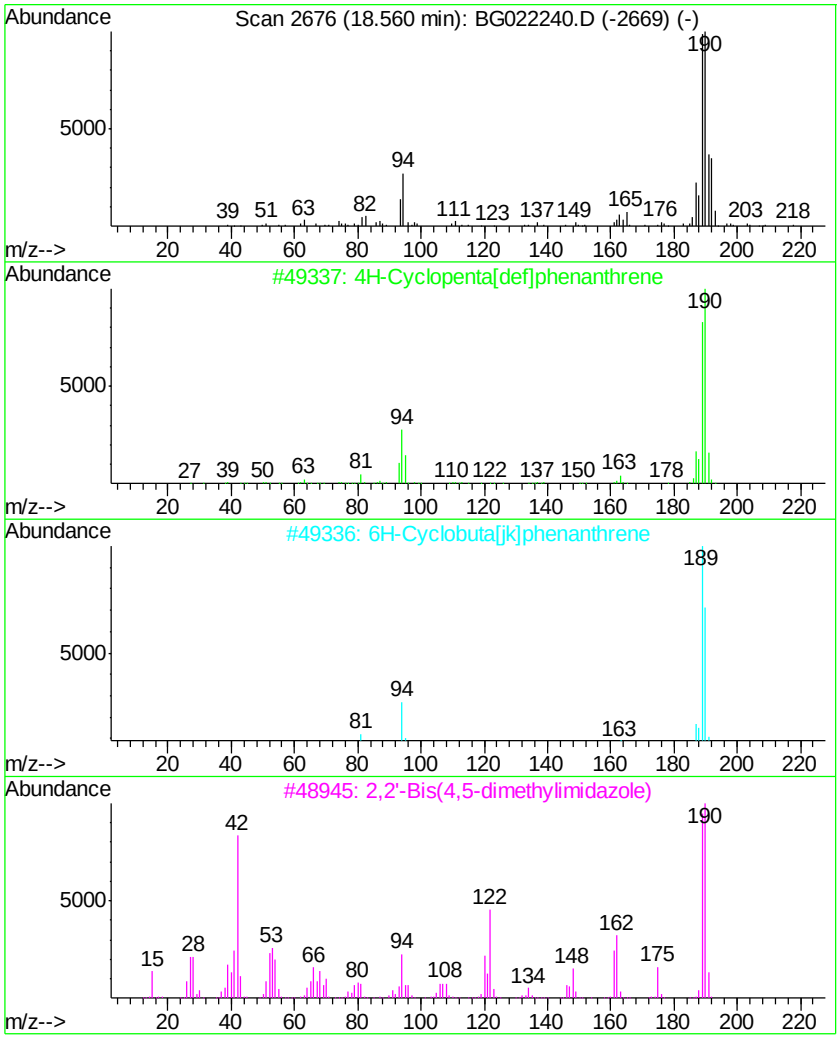
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

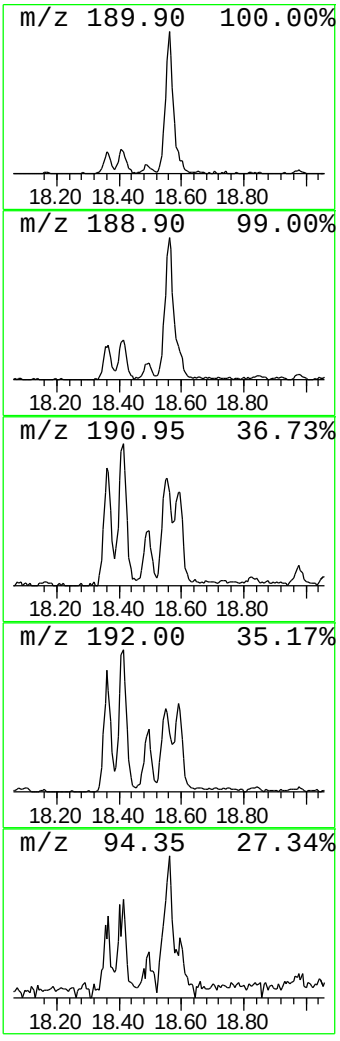
****

Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.56	5.38 ng	157604	Phenanthrene-d10	17.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	64
3	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	38
4	4,4'-Difluorobiphenyl	190	C12H8F2	000398-23-2	23
5	4-Cyano-4-hydroxy-3-methyl-2-phe...	216	C13H16N2O	1000216-11-7	22





Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
Data File : BG022240.D
Acq On : 8 Jun 2016 18:10
Operator : UM/SJ
Sample : H3402-03 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DUP-3

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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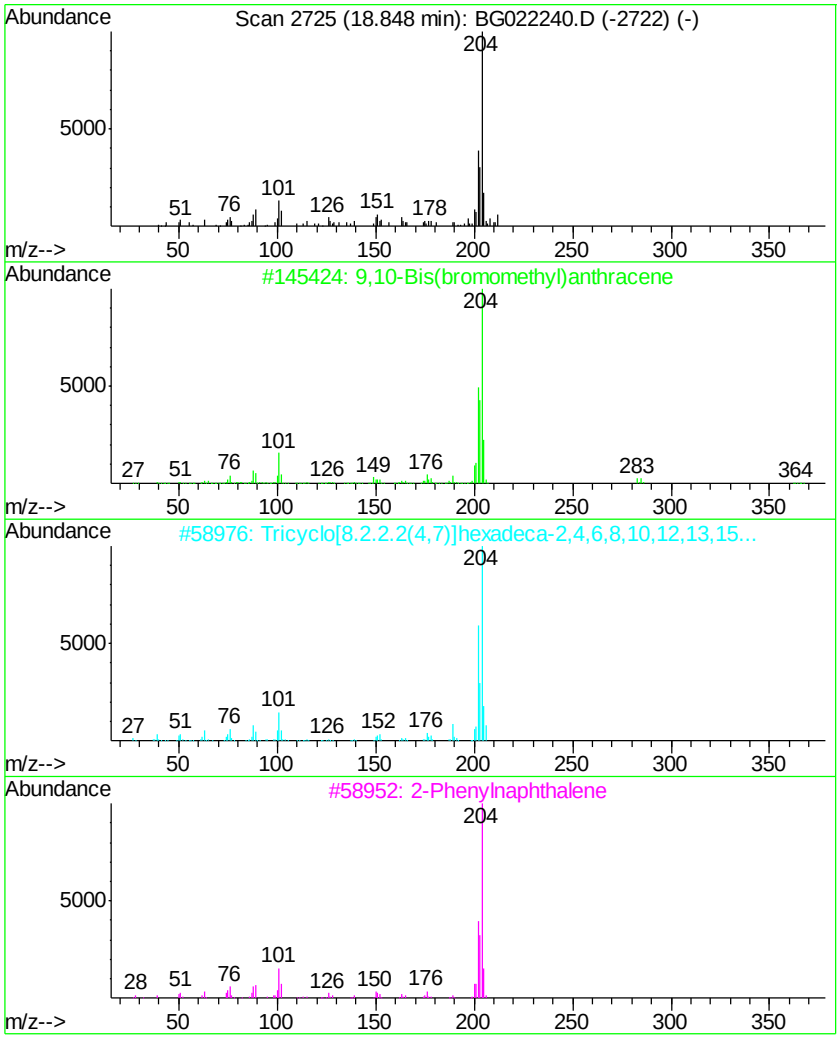
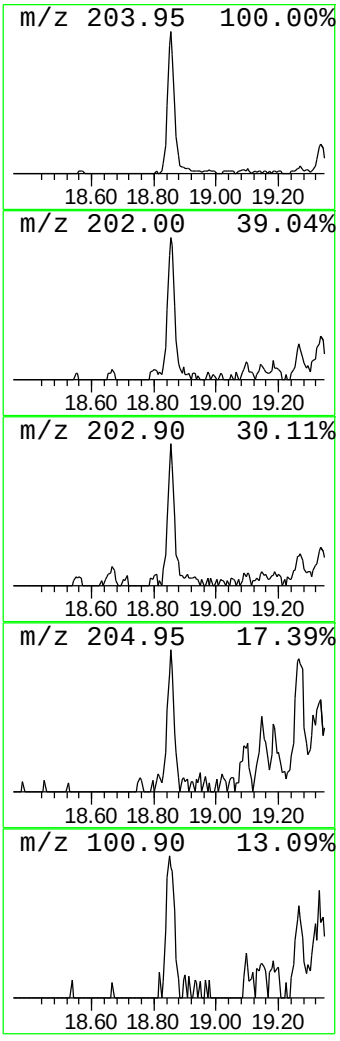
Peak Number 10 9,10-Bis(bromomethyl)anthra... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.85	2.16 ng	63256	Phenanthrene-d10	17.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	86
2	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	86
3	2-Phenylnaphthalene	204	C16H12	035465-71-5	70
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	70
5	[1,1'-Biphenyl]-2-ol, 5-chloro-	204	C12H9ClO	000607-12-5	49





Data Path : Z:\HPCHEM1\BNA G\DATA\BG060816\
Data File : BG022240.D
Acq On : 8 Jun 2016 18:10
Operator : UM/SJ
Sample : H3402-03 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DUP-3

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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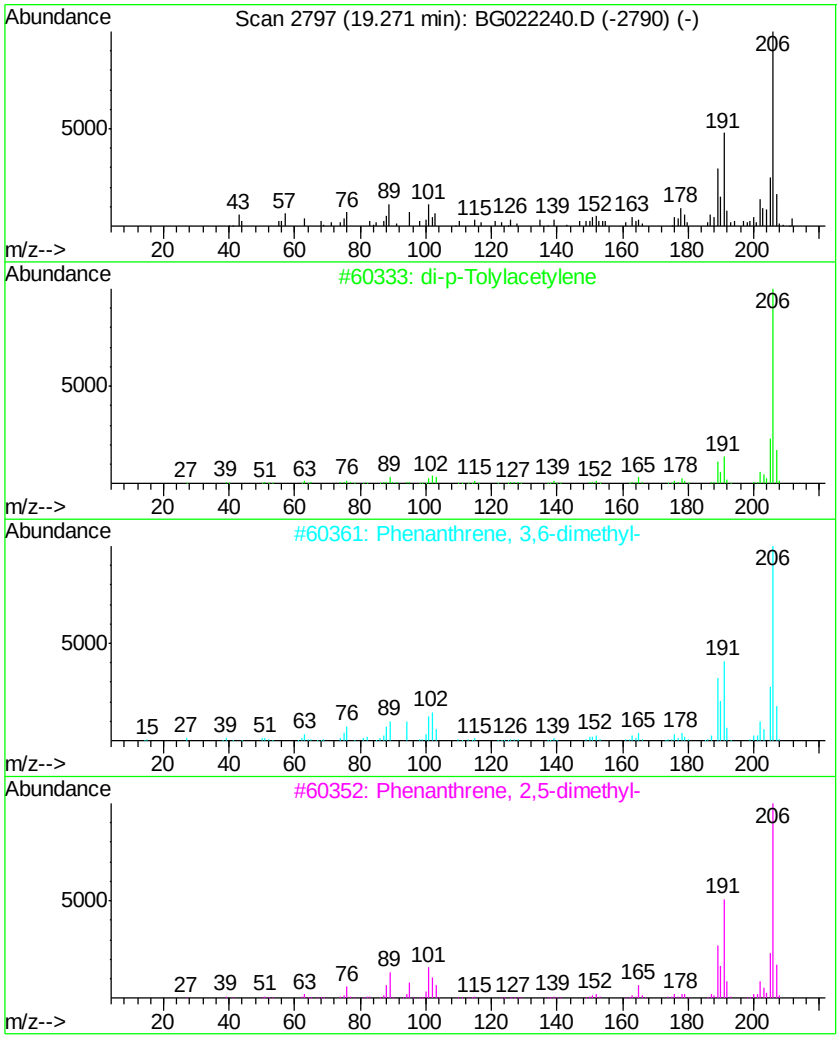
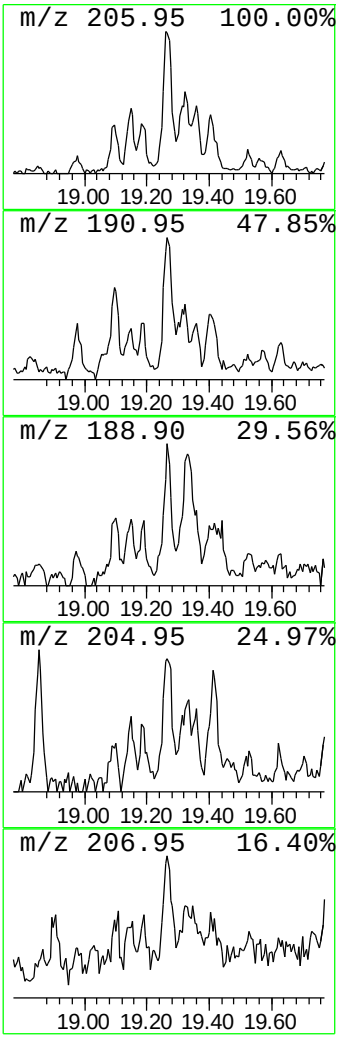
Peak Number 11 di-p-Tolylacetylene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.27	2.63 ng	77133	Phenanthrene-d10	17.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	di-p-Tolylacetylene	206	C16H14	002789-88-0	95
2	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	93
3	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
4	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	93
5	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	91





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Instrument : BNA_G
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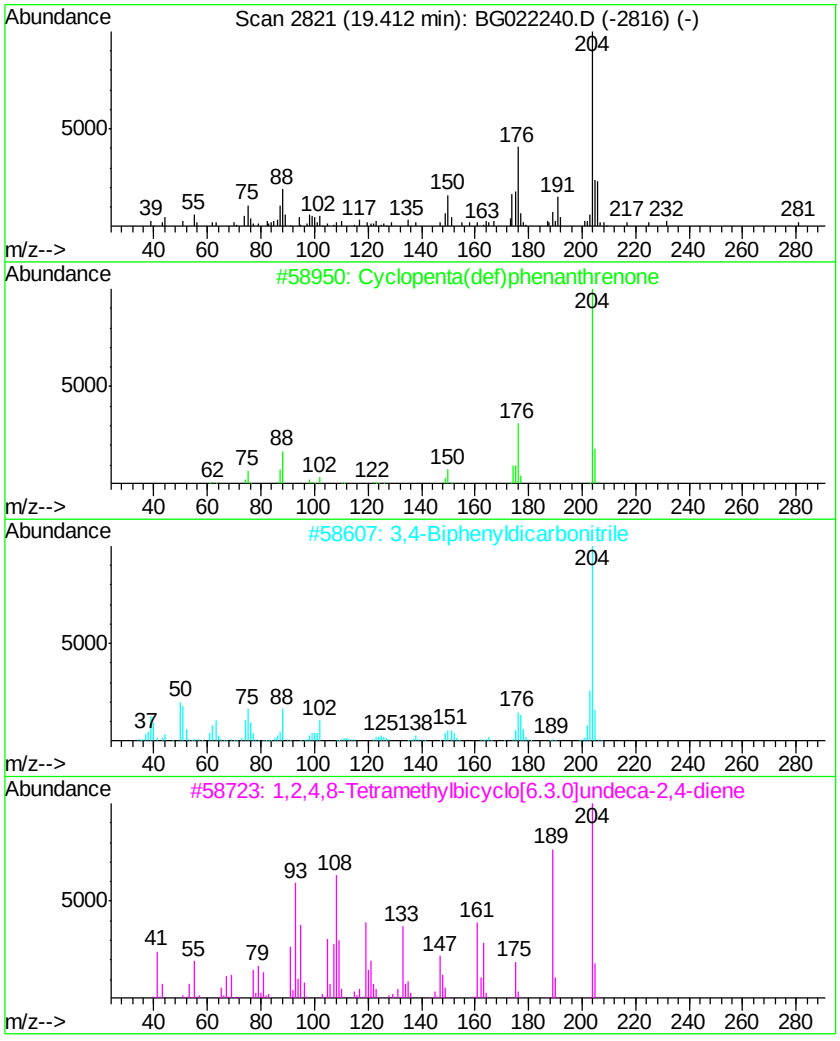
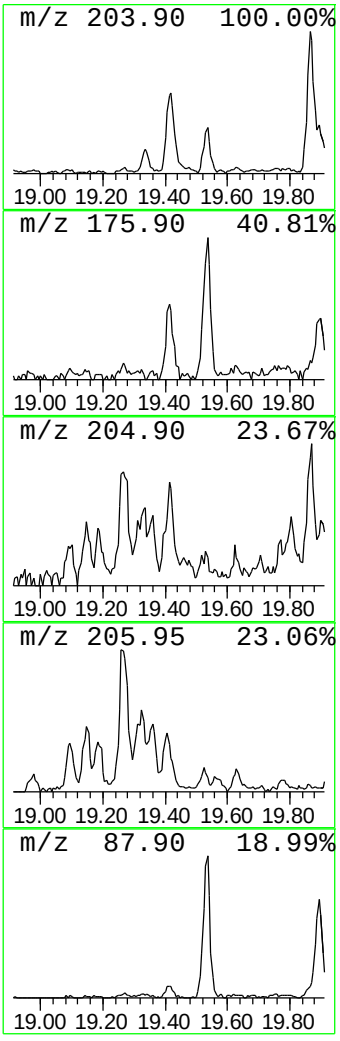
Peak Number 12 Cyclopenta(def)phenanthrenone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.41	2.49 ng	72947	Phenanthrene-d10	17.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	87
2	3,4-Biphenyldicarbonitrile	204	C14H8N2	004128-63-6	42
3	1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	38
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	38
5	[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	38





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Data File : BG022240.D
Acq On : 8 Jun 2016 18:10
Operator : UM/SJ
Sample : H3402-03 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DUP-3

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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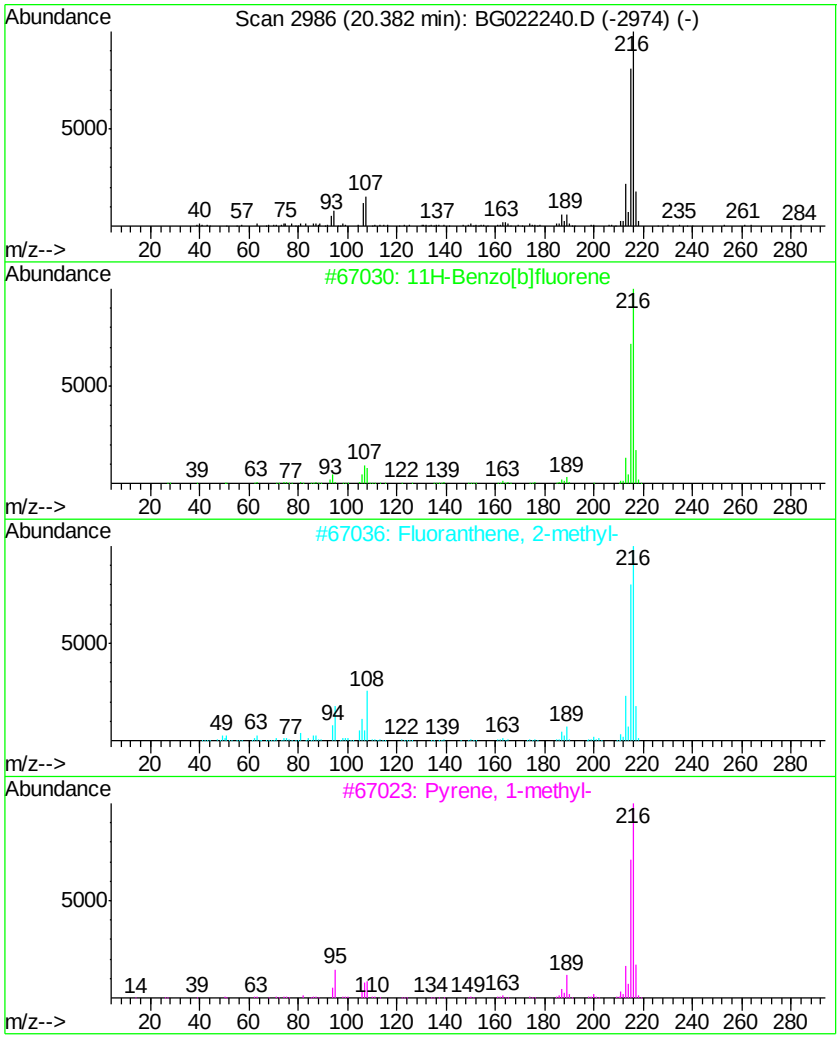
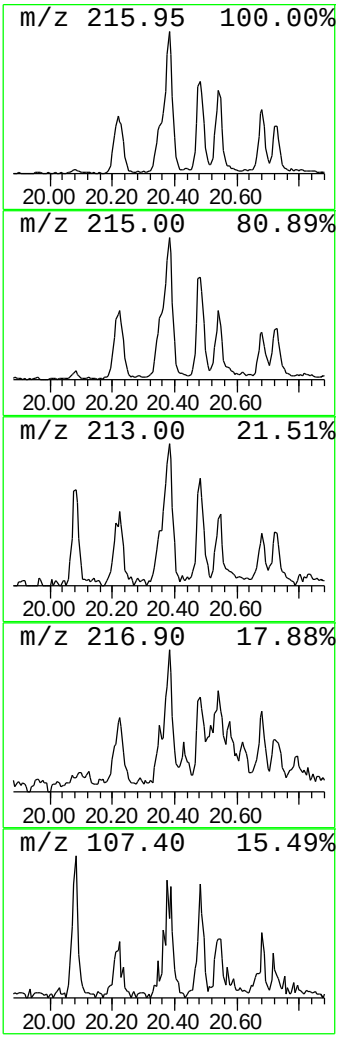
Peak Number 13 11H-Benzo[b]fluorene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.38	3.61 ng	215246	Chrysene-d12	21.77

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	96
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	94
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	86





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Sample : H3402-03 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DUP-3

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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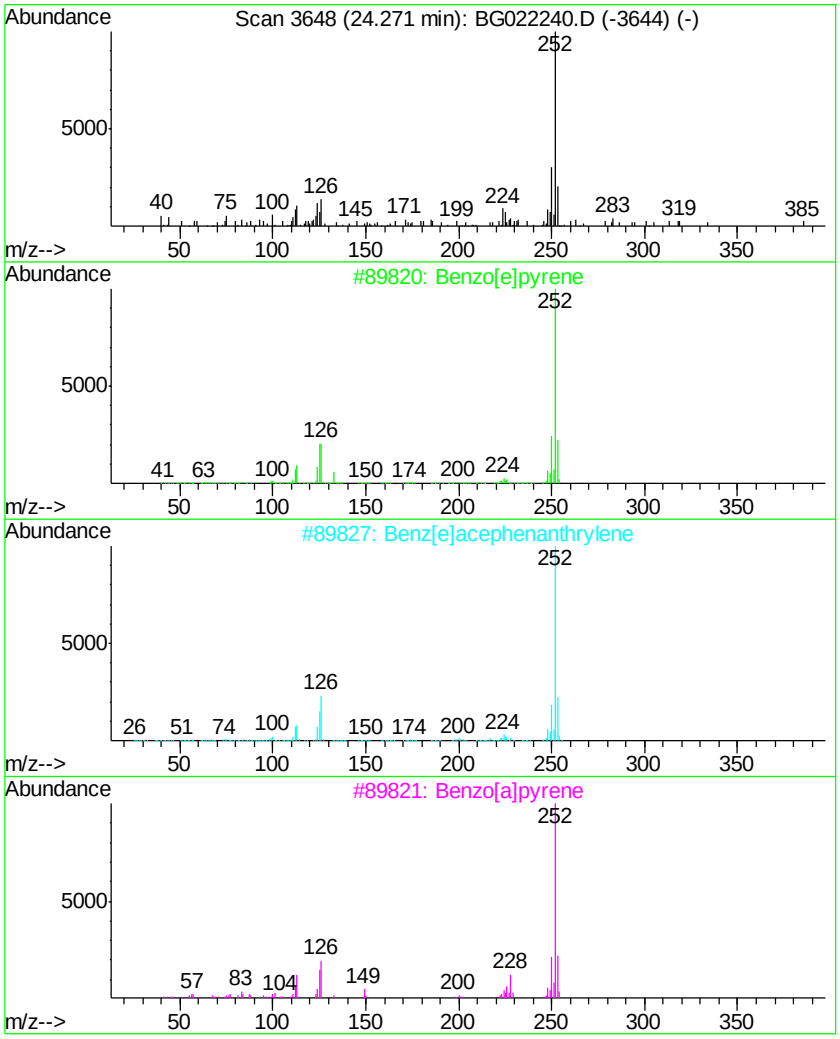
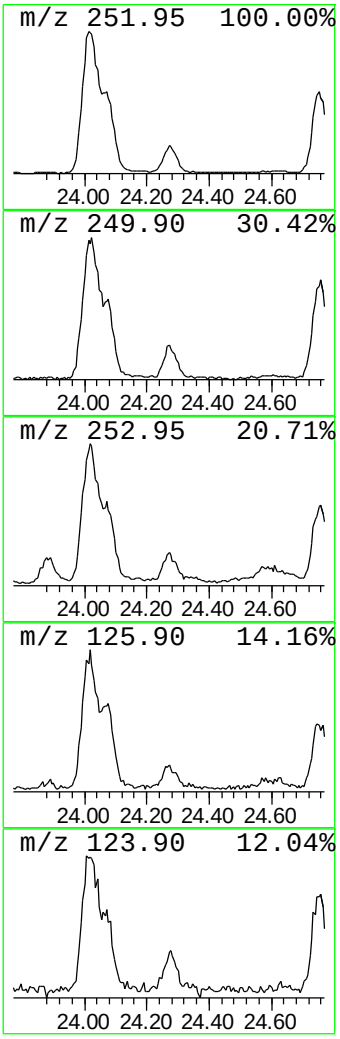
Peak Number 14 Benzo[e]pyrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.27	2.64 ng	77565	Perylene-d12	25.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	81
2	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	76
3	Benzo[a]pyrene	252	C20H12	000050-32-8	68
4	Benzo[k]fluoranthene	252	C20H12	000207-08-9	64
5	Perylene	252	C20H12	000198-55-0	59





Tentatively Identified Compound (LSC) summary

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Data File : BG022240.D
Acq On : 8 Jun 2016 18:10
Operator : UM/SJ
Sample : H3402-03 2X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DUP-3

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
unknown5.18	5.18	3.9	ng	39151	1	8.11	203160	20.0
unknown7.64	7.64	53.7	ng	545250	1	8.11	203160	20.0
5H-Indeno[1,2-b]p...	17.90	2.5	ng	72296	4	17.49	585456	20.0
Anthracene, 2-met...	18.36	2.9	ng	85449	4	17.49	585456	20.0
Anthracene, 9-met...	18.41	3.8	ng	111550	4	17.49	585456	20.0
4H-Cyclopenta[def...	18.56	5.4	ng	157604	4	17.49	585456	20.0
9,10-Bis(bromomet...	18.85	2.2	ng	63256	4	17.49	585456	20.0
di-p-Tolylacetylene	19.27	2.6	ng	77133	4	17.49	585456	20.0
Cyclopenta(def)ph...	19.41	2.5	ng	72947	4	17.49	585456	20.0
11H-Benzo[b]fluorene	20.38	3.6	ng	215246	5	21.77	1191520	20.0
Benzo[e]pyrene	24.27	2.6	ng	77565	6	25.06	587385	20.0