LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060615\
Data File : BG017531.D
Acq On : 7 Jun 2015 12:50
Operator : TP/IZ
Sample : G2457-07 2X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-4

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-BG060315.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	2.764	9	13	19	rVB	31419	42283	5.99%	0.458%
2	4.726	342	347	351	rBV	12494	15454	2.19%	0.168%
3	5.220	425	431	440	rBV	232315	342981	48.62%	3.718%
4	6.836	700	706	715	rBV	249677	390009	55.28%	4.228%
5	7.165	756	762	770	rBV	263243	406310	57.59%	4.405%
6	7.617	833	839	847	rVB	110028	174014	24.67%	1.886%
7	7.940	887	894	902	rBV	193952	296519	42.03%	3.215%
8	8.786	1030	1038	1049	rBV	149976	253004	35.86%	2.743%
9	10.414	1309	1315	1321	rBV	141195	240417	34.08%	2.606%
10	10.461	1321	1323	1329	rVB3	6658	9430	1.34%	0.102%
11	12.905	1733	1739	1748	rBV	345570	499183	70.76%	5.412%
12	13.604	1853	1858	1863	rBV4	5066	8703	1.23%	0.094%
13	13.751	1878	1883	1891	rBV2	77198	110399	15.65%	1.197%
14	14.010	1923	1927	1931	rVB3	5702	8525	1.21%	0.092%
15	14.292	1968	1975	1981	rBV2	211416	310117	43.96%	3.362%
16	14.356	1981	1986	1991	rVB2	20995	30660	4.35%	0.332%
17	14.697	2037	2044	2051	rBV3	10850	19300	2.74%	0.209%
18	15.091	2104	2111	2117	rBV5	4332	9699	1.37%	0.105%
19	15.149	2117	2121	2126	rVB3	7317	8210	1.16%	0.089%
20	15.343	2149	2154	2160	rBV	22863	32808	4.65%	0.356%
21	15.637	2202	2204	2209	rBV4	7127	9043	1.28%	0.098%
22	15.790	2224	2230	2238	rBV2	296028	415139	58.85%	4.500%
23	16.013	2265	2268	2272	rBV5	12207	20863	2.96%	0.226%
24	16.195	2296	2299	2305	rVB4	4920	9190	1.30%	0.100%
25	16.260	2305	2310	2312	rBV5	6492	8519	1.21%	0.092%
26	16.354	2320	2326	2329	rBV5	7832	13140	1.86%	0.142%
27	16.401	2331	2334	2338	rVB5	5275	8190	1.16%	0.089%
28	16.460	2340	2344	2348	rBV5	5969	10158	1.44%	0.110%
29	16.589	2358	2366	2367	rBV6	7078	10610	1.50%	0.115%
30	16.701	2381	2385	2390	rVB4	9294	14619	2.07%	0.158%
31	16.806	2399	2403	2407	rVB7	10557	15562	2.21%	0.169%
32	16.859	2407	2412	2416	rBV2	14019	19276	2.73%	0.209%
33	17.041	2436	2443	2447	rBV2	241602	350968	49.75%	3.805%
34	17.082	2447	2450	2457	rVV	239390	317645	45.03%	3.444%

LSC Area Percent Report

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Data File : BG017531.D
Acq On : 7 Jun 2015 12:50
Operator : TP/IZ
Sample : G2457-07 2X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-4

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-BG060315.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	17.176	2457	2466	2471	rVB2	60262	93978	13.32%	1.019%
36	17.241	2471	2477	2481	rBV7	8109	15149	2.15%	0.164%
37	17.458	2510	2514	2520	rVV	12239	16691	2.37%	0.181%
38	17.547	2524	2529	2530	rBV4	6652	8204	1.16%	0.089%
39	17.564	2530	2532	2535	rVB2	6417	7182	1.02%	0.078%
40	17.923	2588	2593	2595	rBV2	29234	38239	5.42%	0.415%
41	17.970	2595	2601	2605	rVV2	49392	87442	12.39%	0.948%
42	18.046	2609	2614	2621	rBV	21902	36771	5.21%	0.399%
43	18.116	2621	2626	2630	rVV2	50715	92554	13.12%	1.003%
44	18.152	2630	2632	2638	rVB2	23184	31952	4.53%	0.346%
45	18.240	2642	2647	2650	rVB6	8563	13976	1.98%	0.152%
46	18.281	2650	2654	2658	rBV6	6399	10445	1.48%	0.113%
47	18.428	2673	2679	2682	rBV2	26367	36131	5.12%	0.392%
48	18.469	2682	2686	2690	rVB4	15430	19264	2.73%	0.209%
49	18.669	2715	2720	2724	rVB8	10563	18644	2.64%	0.202%
50	18.722	2726	2729	2733	rVV5	7902	10615	1.50%	0.115%
51	18.757	2733	2735	2739	rVB4	6983	8113	1.15%	0.088%
52	18.839	2746	2749	2753	rBV3	18426	29187	4.14%	0.316%
53	18.910	2758	2761	2762	rVV3	10813	12273	1.74%	0.133%
54	18.939	2764	2766	2769	rVV3	7922	7630	1.08%	0.083%
55	18.986	2771	2774	2782	rVB6	14764	26950	3.82%	0.292%
56	19.109	2789	2795	2802	rBV	401671	514921	72.99%	5.582%
57	19.174	2802	2806	2809	rVV6	11387	15169	2.15%	0.164%
58	19.209	2809	2812	2815	rVV5	11917	14358	2.04%	0.156%
59	19.256	2818	2820	2824	rVB3	11166	14474	2.05%	0.157%
60	19.356	2833	2837	2840	rBV	16519	18740	2.66%	0.203%
61	19.444	2848	2852	2853	rBV	66614	73505	10.42%	0.797%
62	19.474	2853	2857	2866	rVV	328942	457668	64.87%	4.962%
63	19.550	2866	2870	2878	rVB4	21167	39547	5.61%	0.429%
64	19.679	2884	2892	2897	rBV	539812	687488	97.45%	7.453%
65	19.797	2908	2912	2914	rBV3	23219	32493	4.61%	0.352%
66	19.973	2939	2942	2948	rVB	55957	73333	10.39%	0.795%
67	20.073	2956	2959	2963	rVB	44617	56128	7.96%	0.608%
68	20.132	2963	2969	2973	rBV3	38602	67853	9.62%	0.736%
69	20.214	2981	2983	2988	rVB6	10218	10557	1.50%	0.114%
70	20.273	2990	2993	2996	rVB2	18800	18312	2.60%	0.199%
71	20.320	2998	3001	3004	rVV3	22037	26028	3.69%	0.282%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060615\
Data File : BG017531.D
Acq On : 7 Jun 2015 12:50
Operator : TP/IZ
Sample : G2457-07 2X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-4

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-BG060315.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	20.725	3067	3070	3074	rVB3	22037	26688	3.78%	0.289%
73	20.884	3095	3097	3101	rVB2	18526	20155	2.86%	0.218%
74	20.925	3101	3104	3106	rBV	26668	32156	4.56%	0.349%
75	20.954	3106	3109	3116	rVB3	48876	87086	12.34%	0.944%
76	21.236	3151	3157	3161	rBV2	398489	705466	100.00%	7.648%
77	21.278	3161	3164	3172	rVB	174669	215877	30.60%	2.340%
78	21.389	3181	3183	3188	rVB	26314	27915	3.96%	0.303%
79	22.805	3419	3424	3428	rBV	139780	274858	38.96%	2.980%
80	23.287	3502	3506	3511	rVB	76848	129853	18.41%	1.408%
81	23.381	3517	3522	3527	rVB	139559	227968	32.31%	2.471%
82	23.481	3534	3539	3544	rVB	185505	303391	43.01%	3.289%

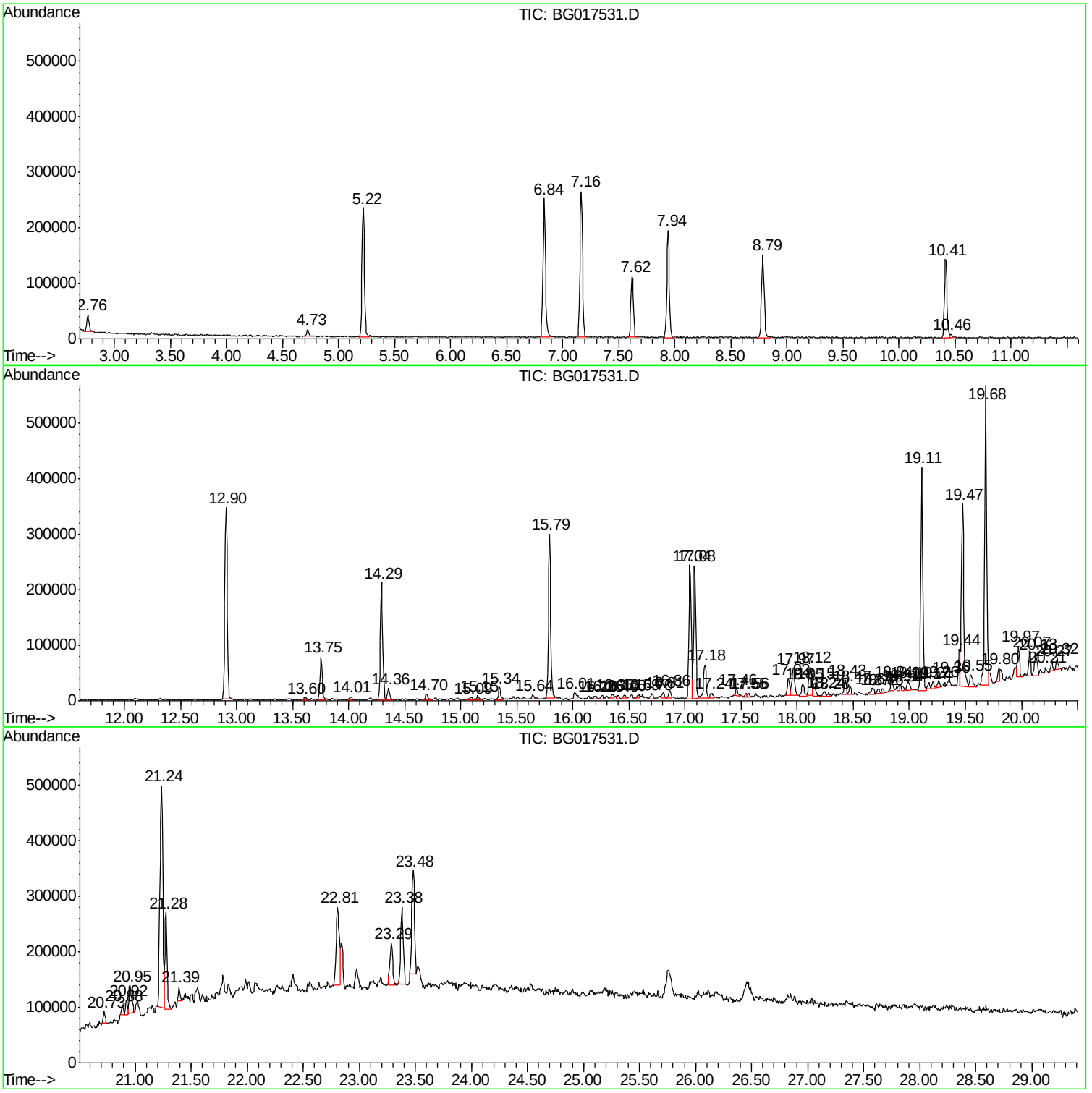
Sum of corrected areas: 9224324

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060615\
Data File : BG017531.D
Acq On : 7 Jun 2015 12:50
Operator : TP/IZ
Sample : G2457-07 2X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SP-4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.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\BG060615\
Data File : BG017531.D
Acq On : 7 Jun 2015 12:50
Operator : TP/IZ
Sample : G2457-07 2X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

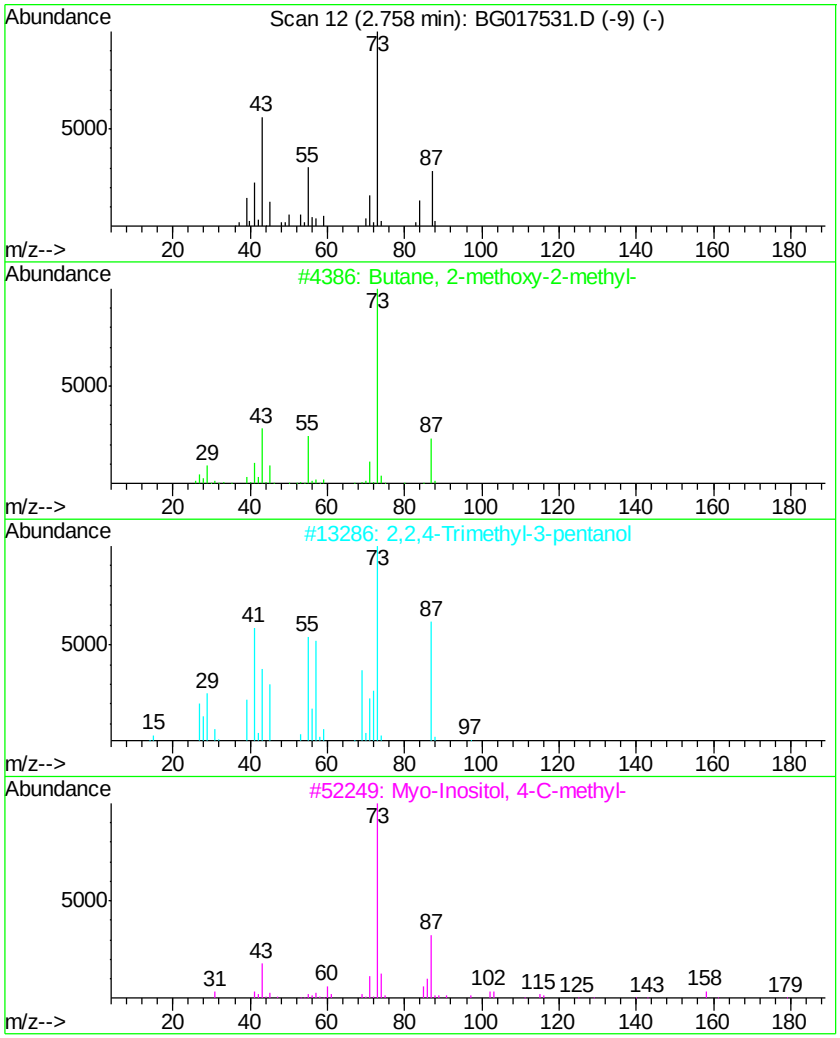
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

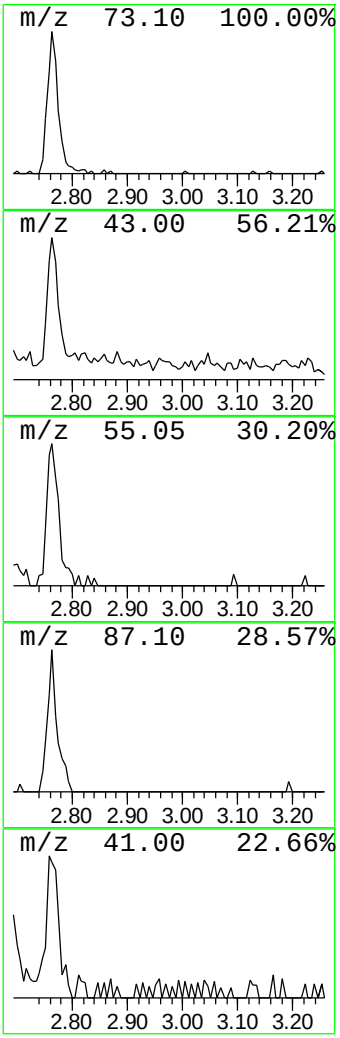
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.76	4.86 ng	42283	1,4-Dichlorobenzene-d4	7.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	36
3	Myo-Inositol, 4-C-methyl-	194	C7H14O6	000472-95-7	36
4	Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	36
5	Myo-Inositol, 2-C-methyl-	194	C7H14O6	000472-96-8	36





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060615\
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Acq On : 7 Jun 2015 12:50
Operator : TP/IZ
Sample : G2457-07 2X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

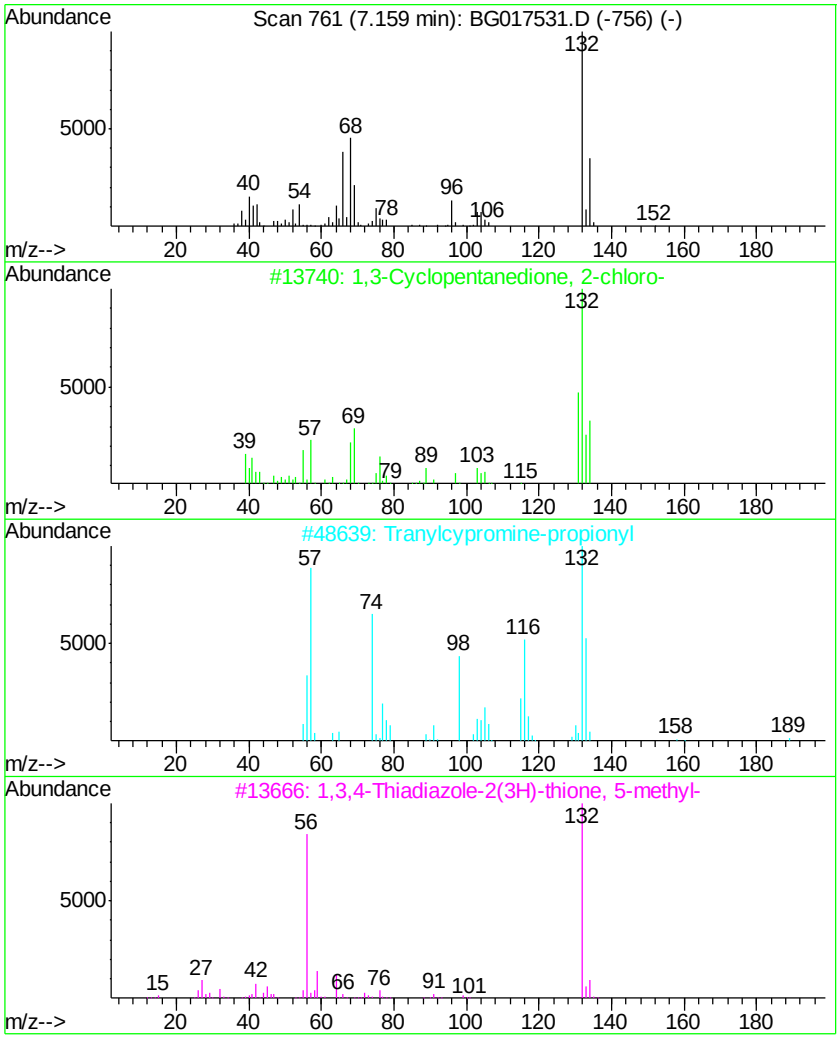
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

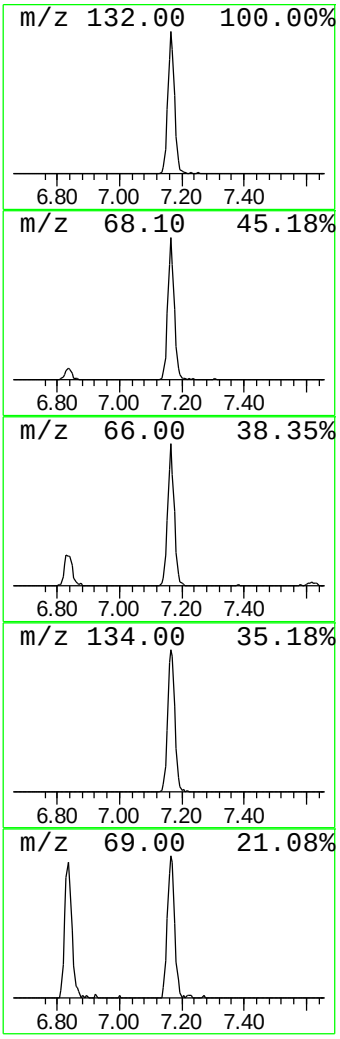
****

Peak Number 2 unknown7.16 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.16	46.70 ng	406310	1,4-Dichlorobenzene-d4	7.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	32
2	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
3	1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	16
4	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12
5	Benzene, 1-methyl-4-(1-methyleth...	132	C10H12	001195-32-0	12





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060615\
Data File : BG017531.D
Acq On : 7 Jun 2015 12:50
Operator : TP/IZ
Sample : G2457-07 2X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

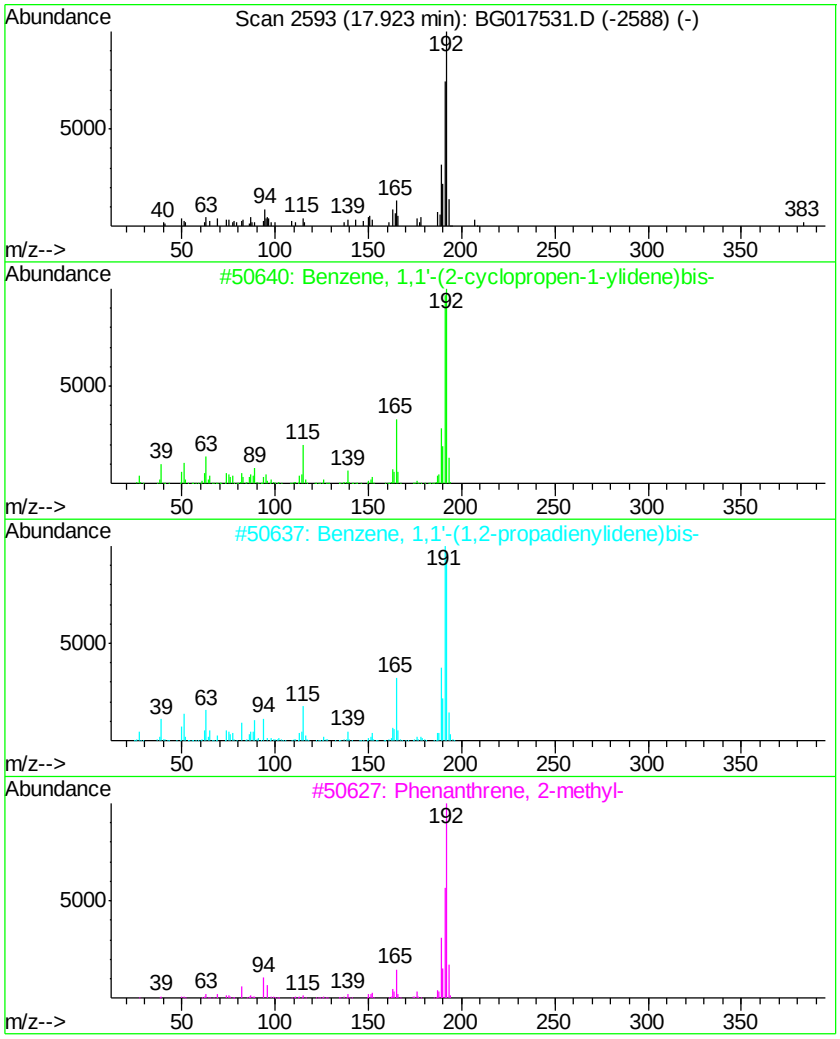
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

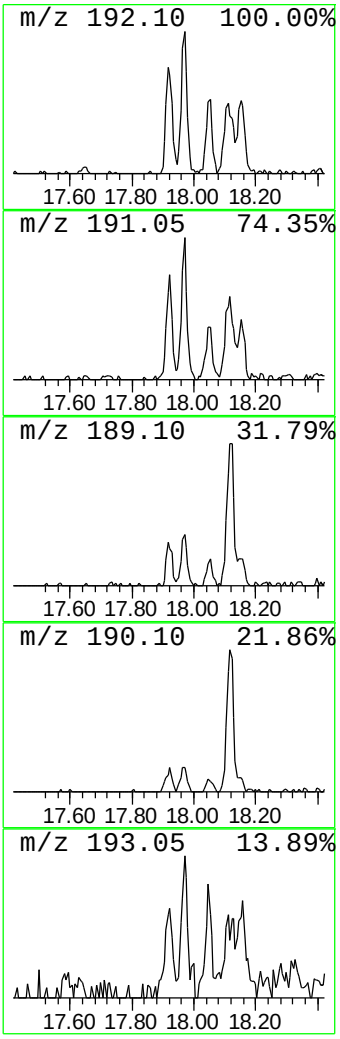
****

Peak Number 5 Benzene, 1,1'-(2-cycloprope... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.92	2.18 ng	38239	Phenanthrene-d10	17.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(2-cyclopropen-1-y...	192	C15H12	022825-21-4	81
2	Benzene, 1,1'-(1,2-propadienylid...	192	C15H12	014251-57-1	76
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	74
4	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	74
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	70





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060615\
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Operator : TP/IZ
Sample : G2457-07 2X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.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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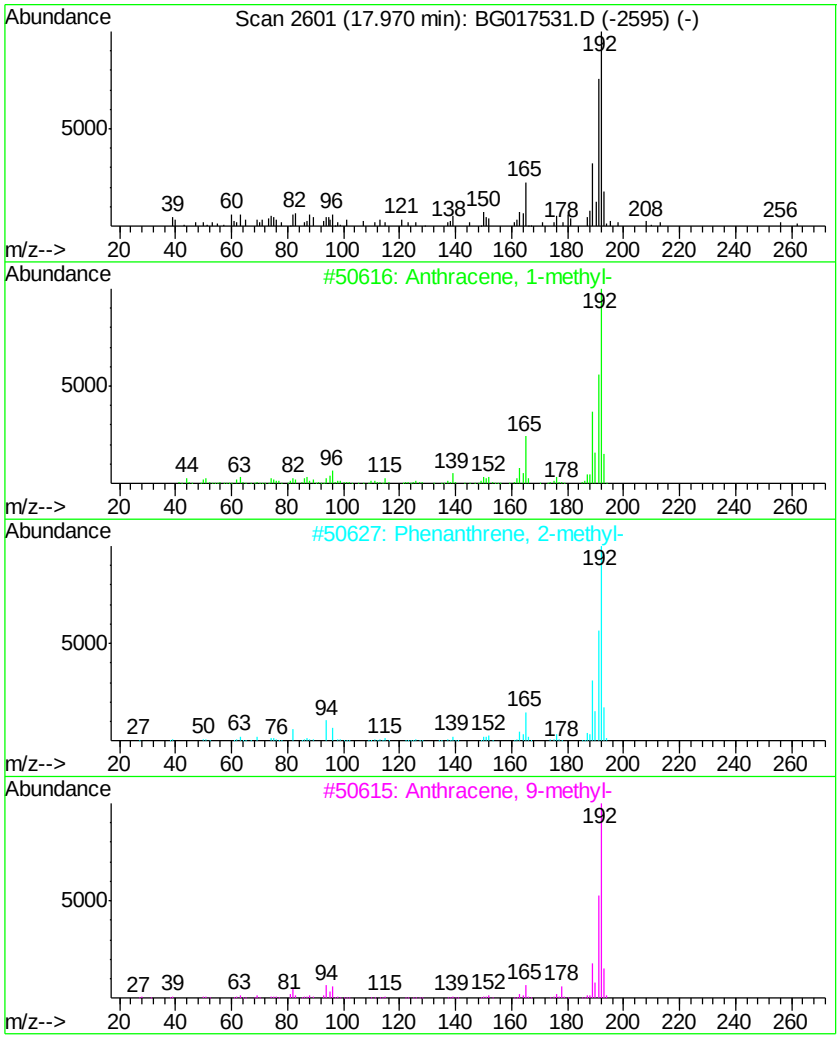
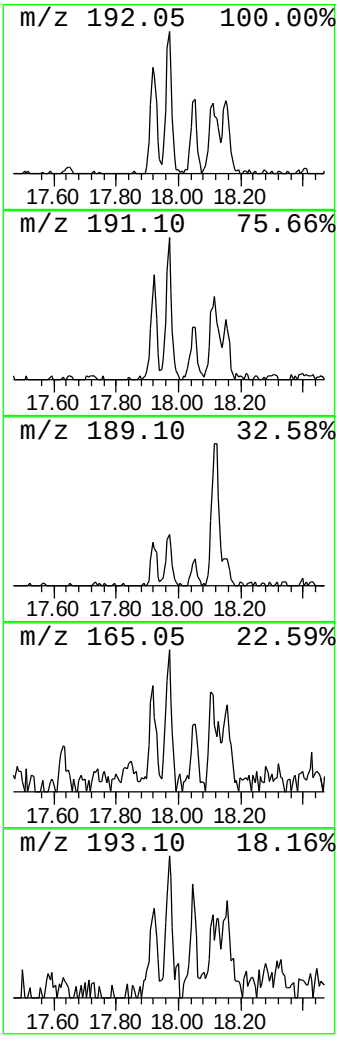
Peak Number 6 Anthracene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.97	4.98 ng	87442	Phenanthrene-d10	17.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	92
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
4	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	90
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060615\
Data File : BG017531.D
Acq On : 7 Jun 2015 12:50
Operator : TP/IZ
Sample : G2457-07 2X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.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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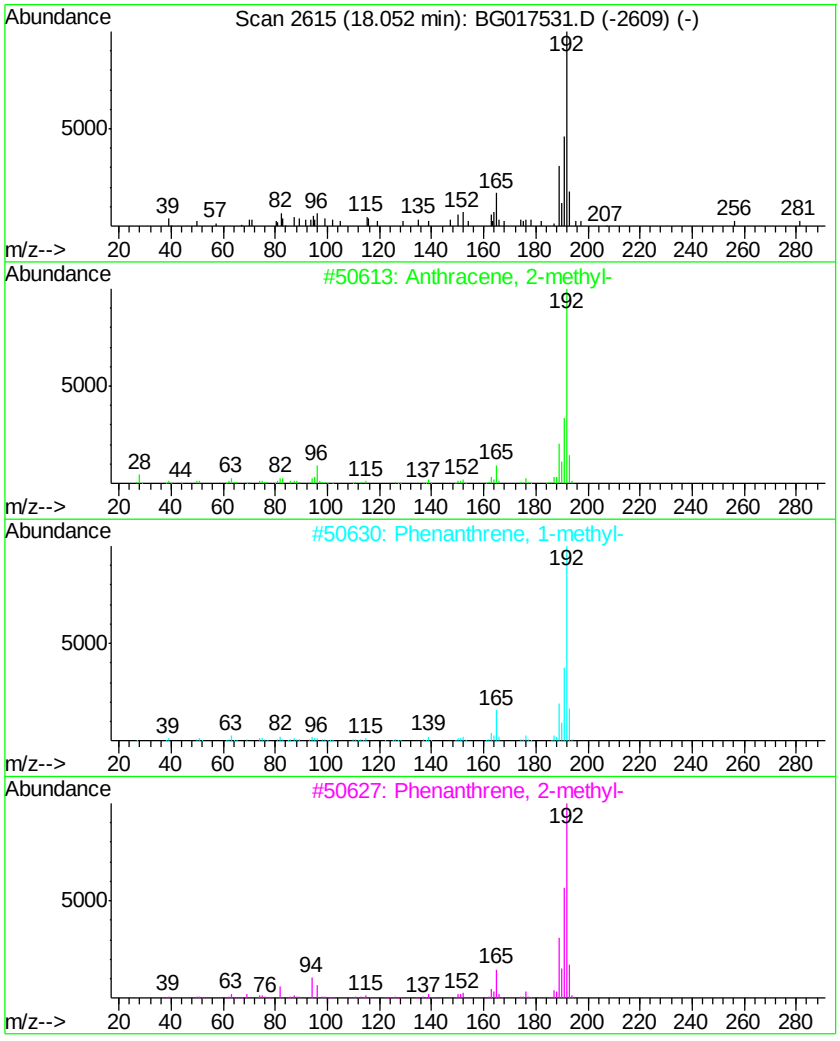
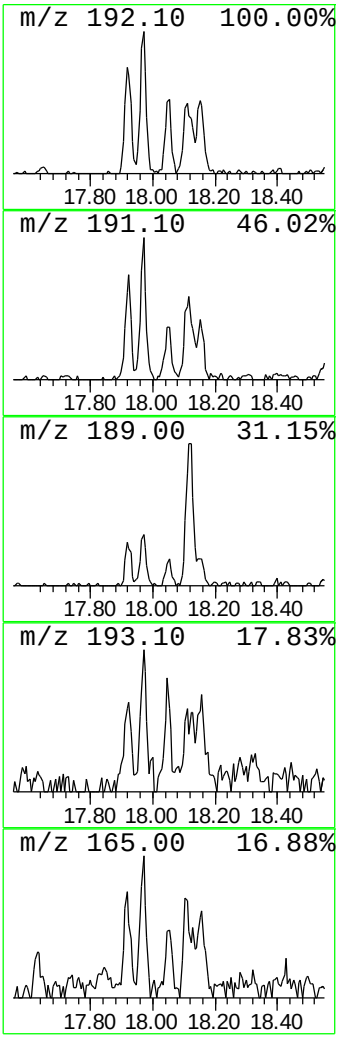
Peak Number 7 Anthracene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.05	2.10 ng	36771	Phenanthrene-d10	17.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	74
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	64
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	64
4	1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	64
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	64





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060615\
Data File : BG017531.D
Acq On : 7 Jun 2015 12:50
Operator : TP/IZ
Sample : G2457-07 2X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

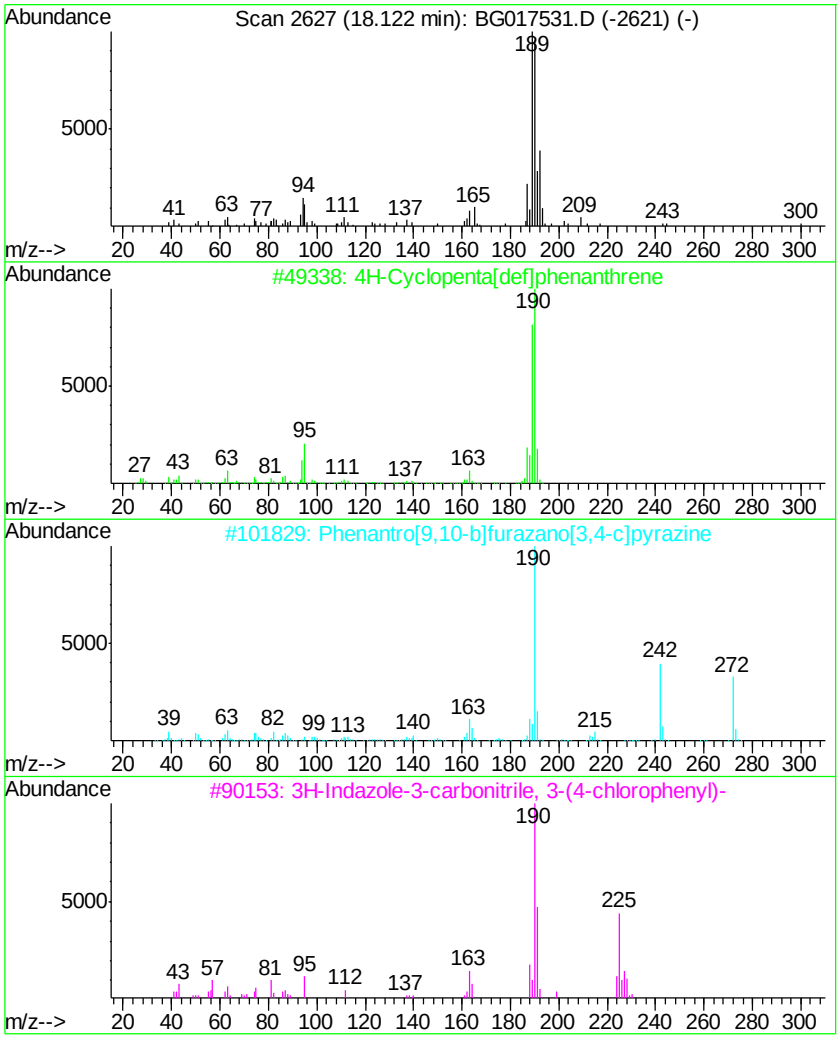
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

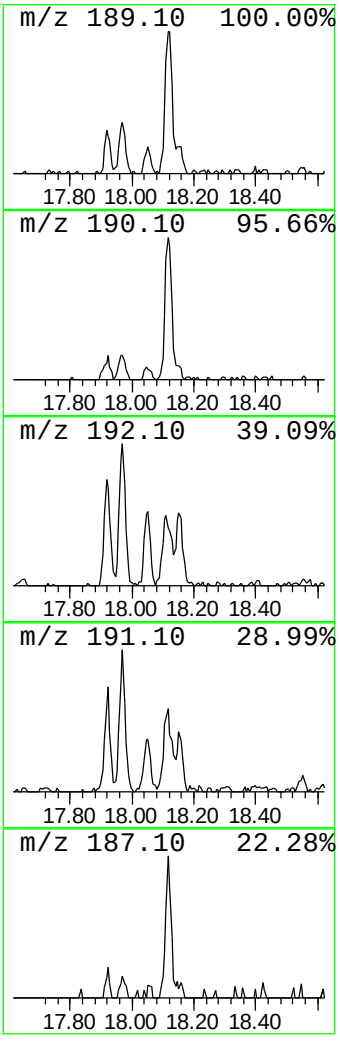
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.12	5.27 ng	92554	Phenanthrene-d10	17.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
2	Phenantro[9,10-b]furazano[3,4-c]...	272	C16H8N4O	024294-86-8	14
3	3H-Indazole-3-carbonitrile, 3-(4...	253	C14H8ClN3	056630-97-8	13
4	Propanoic acid, 2-(2,4-dichlorop...	378	C15H10Cl4O3	284488-02-4	13
5	Pyrimidine, 2-chloro-5-phenyl-	190	C10H7ClN2	022536-62-5	12





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060615\
Data File : BG017531.D
Acq On : 7 Jun 2015 12:50
Operator : TP/IZ
Sample : G2457-07 2X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.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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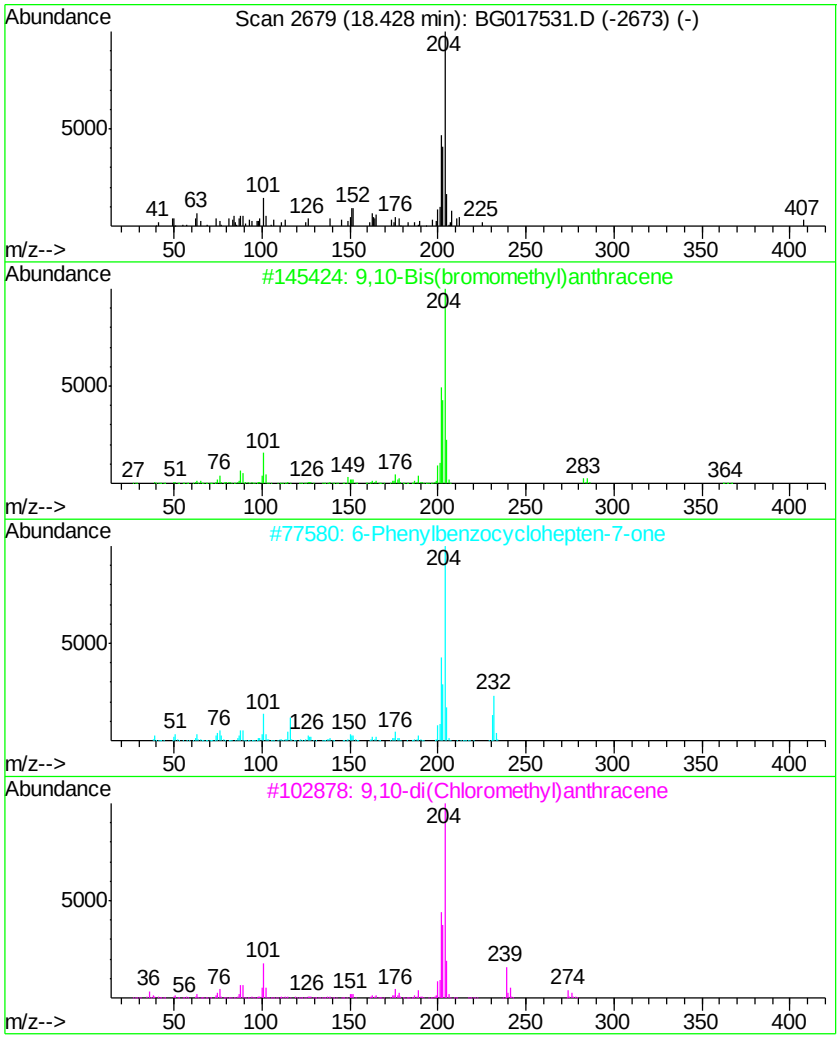
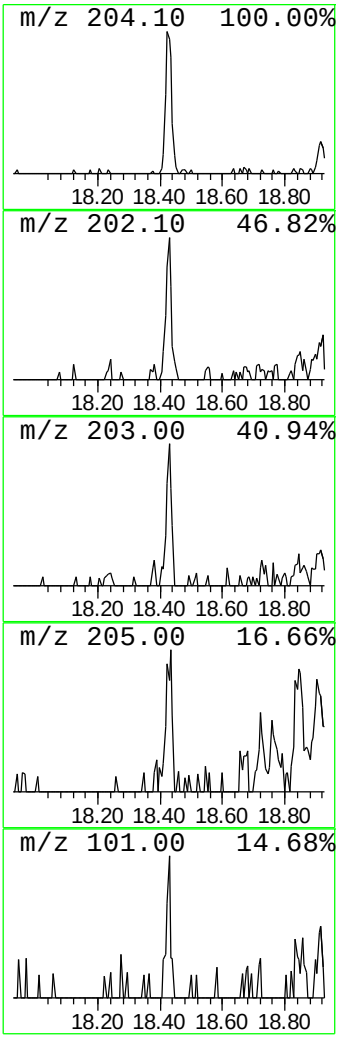
Peak Number 9 9,10-Bis(bromomethyl)anthra... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.43	2.06 ng	36131	Phenanthrene-d10	17.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	80
2	6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	72
3	9,10-di(Chloromethyl)anthracene	274	C16H12Cl2	010387-13-0	72
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	60
5	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	59





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Data File : BG017531.D
Acq On : 7 Jun 2015 12:50
Operator : TP/IZ
Sample : G2457-07 2X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.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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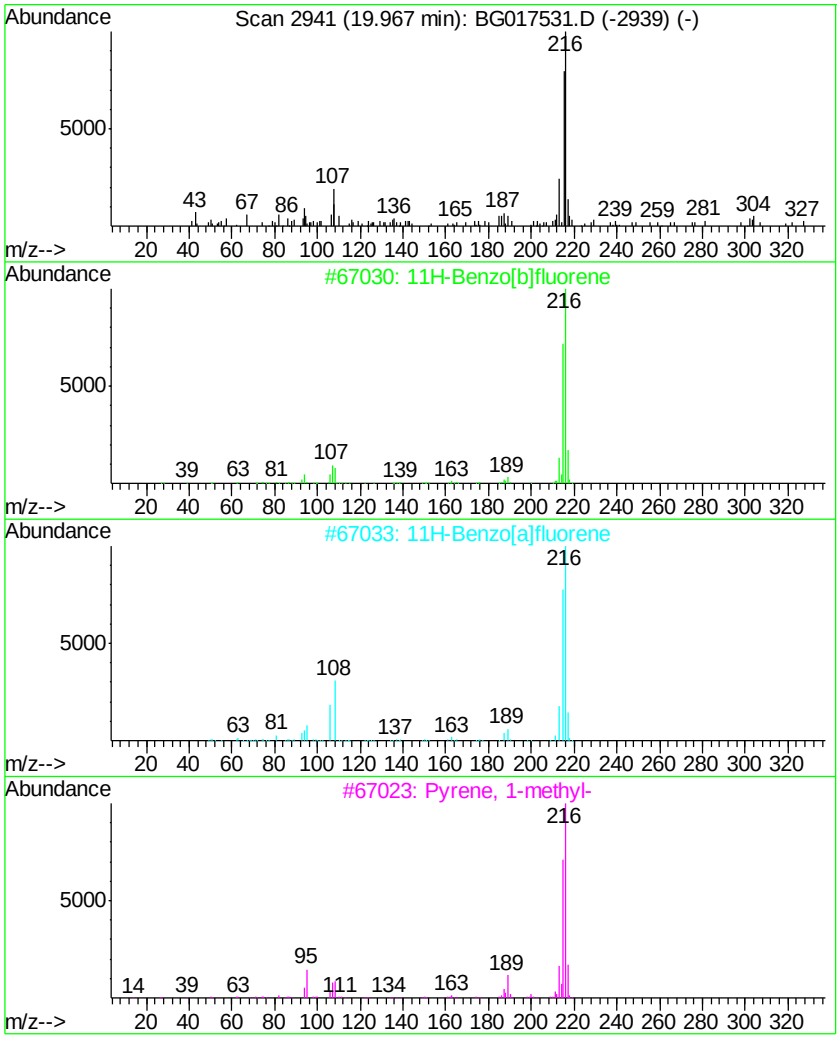
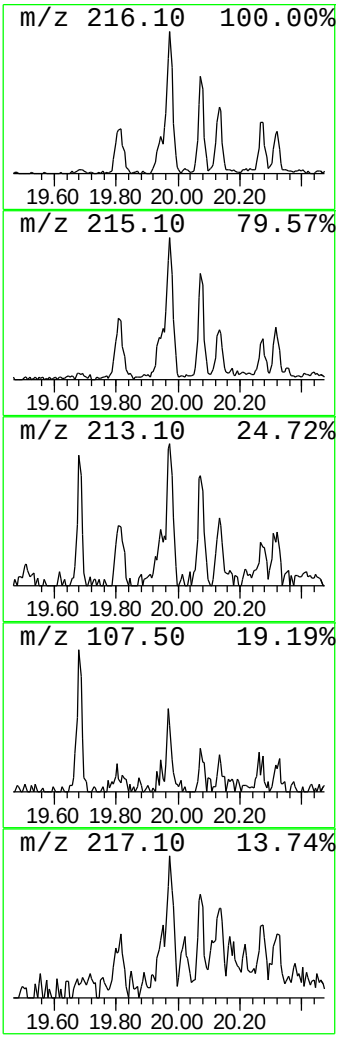
Peak Number 10 11H-Benzo[b]fluorene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.97	2.08 ng	73333	Chrysene-d12	21.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	80
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	80





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Data File : BG017531.D
Acq On : 7 Jun 2015 12:50
Operator : TP/IZ
Sample : G2457-07 2X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.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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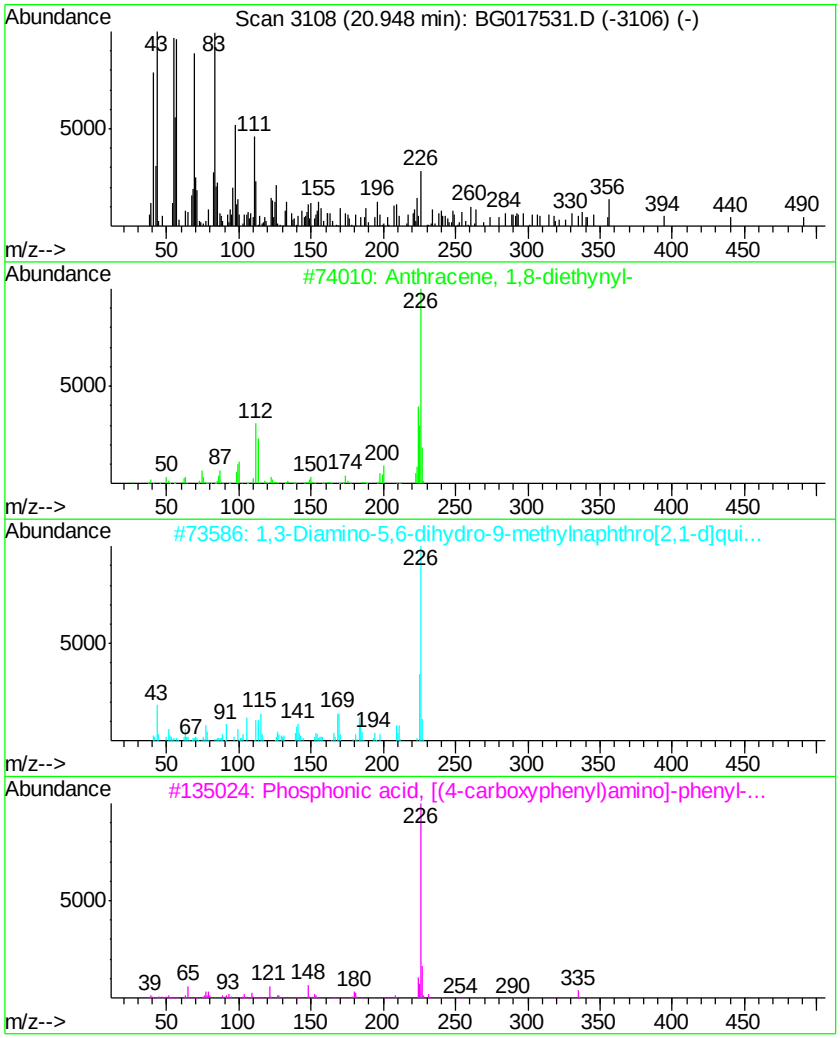
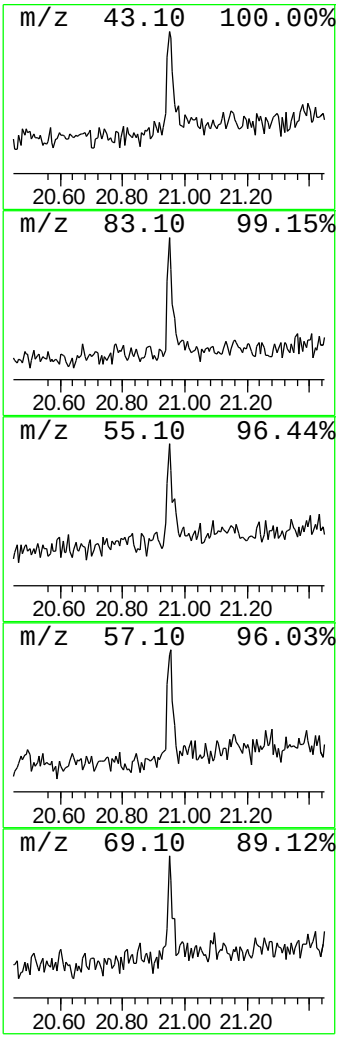
Peak Number 11 unknown20.95 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.95	2.47 ng	87086	Chrysene-d12	21.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 1,8-diethynyl-	226	C18H10	078053-58-4	38
2	1,3-Diamino-5,6-dihydro-9-methyl...	226	C13H14N4	037436-39-8	27
3	Phosphonic acid, [(4-carboxyphen...	335	C16H18NO5P	037753-62-1	17
4	Malononitrile, 1-(4-t-butyl-1-cy...	282	C18H22N2O	1000164-49-6	16
5	Benzene, [4-(3-ethynylphenyl)-1,...	226	C18H10	1000115-87-3	16





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Data File : BG017531.D
Acq On : 7 Jun 2015 12:50
Operator : TP/IZ
Sample : G2457-07 2X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.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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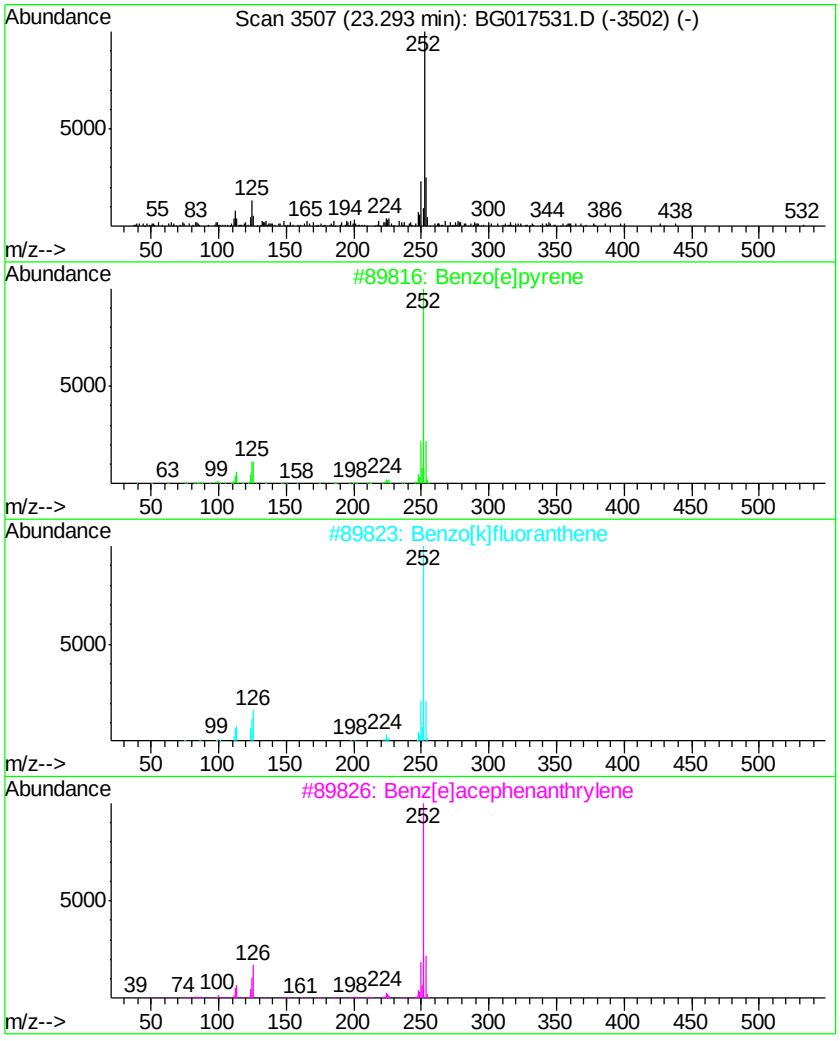
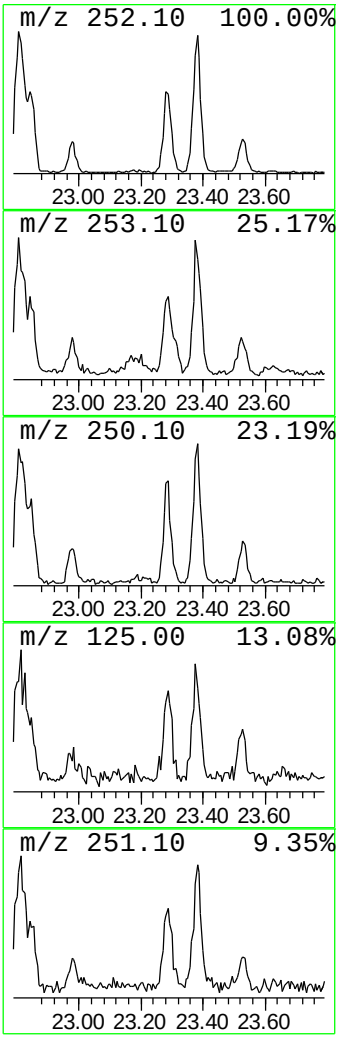
Peak Number 12 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.29	8.56 ng	129853	Perylene-d12	23.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	83
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	81
3	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	81
4	Benzo[a]pyrene	252	C20H12	000050-32-8	74
5	Perylene	252	C20H12	000198-55-0	72





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060615\
Data File : BG017531.D
Acq On : 7 Jun 2015 12:50
Operator : TP/IZ
Sample : G2457-07 2X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.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
Butane, 2-methoxy...	2.76	4.9	ng	42283	1	7.62	174014	20.0
unknown7.16	7.16	46.7	ng	406310	1	7.62	174014	20.0
Benzene, 1,1'-(2-...	17.92	2.2	ng	38239	4	17.04	350968	20.0
Anthracene, 1-met...	17.97	5.0	ng	87442	4	17.04	350968	20.0
Anthracene, 2-met...	18.05	2.1	ng	36771	4	17.04	350968	20.0
4H-Cyclopenta[def...	18.12	5.3	ng	92554	4	17.04	350968	20.0
9,10-Bis(bromomet...	18.43	2.1	ng	36131	4	17.04	350968	20.0
11H-Benzo[b]fluorene	19.97	2.1	ng	73333	5	21.24	705466	20.0
unknown20.95	20.95	2.5	ng	87086	5	21.24	705466	20.0
Benzo[e]pyrene	23.29	8.6	ng	129853	6	23.48	303391	20.0