

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060217\
 Data File : BF095645.D
 Acq On : 2 Jun 2017 21:09
 Operator : SJ/MA
 Sample : I3441-11 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-RO-17273-060117

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % 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_F\METHODS\8270-BF050217.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	4.787	524	527	532	rBV2	5065	10637	1.69%	0.113%
2	5.487	643	646	670	rBV2	63559	233255	37.13%	2.476%
3	7.128	922	925	936	rBV	66821	193193	30.76%	2.051%
4	7.216	936	940	957	rVV	159451	395736	63.00%	4.200%
5	7.340	957	961	966	rVB3	16995	22023	3.51%	0.234%
6	7.540	991	995	1006	rBV	319615	444807	70.81%	4.721%
7	7.775	1031	1035	1050	rBV	163409	248639	39.58%	2.639%
8	8.469	1150	1153	1167	rBV	48457	131026	20.86%	1.391%
9	8.569	1167	1170	1176	rVB3	14206	22634	3.60%	0.240%
10	9.569	1336	1340	1343	rBV	435171	525251	83.62%	5.575%
11	9.604	1343	1346	1362	rVB	110068	198088	31.54%	2.103%
12	10.733	1534	1538	1549	rBV	89245	127176	20.25%	1.350%
13	10.886	1557	1564	1573	rBV	76422	106673	16.98%	1.132%
14	11.345	1638	1642	1651	rBV	262066	323674	51.53%	3.436%
15	11.510	1664	1670	1677	rBV	16925	24279	3.87%	0.258%
16	11.651	1690	1694	1699	rBV2	7114	12204	1.94%	0.130%
17	11.757	1708	1712	1713	rBV2	22037	26826	4.27%	0.285%
18	11.875	1726	1732	1736	rBV	46056	68320	10.88%	0.725%
19	11.916	1736	1739	1749	rVB2	25598	35365	5.63%	0.375%
20	12.051	1758	1762	1772	rVB4	25458	55958	8.91%	0.594%
21	12.175	1778	1783	1792	rBV2	28429	51269	8.16%	0.544%
22	12.392	1816	1820	1826	rBV2	401453	526416	83.81%	5.588%
23	12.445	1826	1829	1838	rVV	72449	105991	16.87%	1.125%
24	12.610	1852	1857	1862	rBV3	7655	14419	2.30%	0.153%
25	12.745	1874	1880	1889	rBV2	43746	77487	12.34%	0.822%
26	12.833	1889	1895	1900	rVB	11930	15275	2.43%	0.162%
27	12.945	1908	1914	1920	rBV4	13361	23102	3.68%	0.245%
28	13.086	1933	1938	1941	rBV4	6509	10838	1.73%	0.115%
29	13.245	1958	1965	1968	rBV4	13634	23354	3.72%	0.248%
30	13.292	1968	1973	1983	rVV	110479	150908	24.02%	1.602%
31	13.386	1983	1989	1991	rVV5	6822	13318	2.12%	0.141%
32	13.433	1994	1997	2001	rVV2	10423	14840	2.36%	0.158%
33	13.569	2017	2020	2029	rBV4	12635	28063	4.47%	0.298%
34	13.704	2039	2043	2051	rBV2	73325	110015	17.51%	1.168%

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35	14.198	2120	2127	2131	rBV4	17068	33675	5.36%	0.357%
36	14.239	2131	2134	2139	rVV3	12870	19898	3.17%	0.211%
37	14.321	2143	2148	2153	rVB4	10291	13777	2.19%	0.146%
38	14.533	2178	2184	2189	rVB5	7034	11978	1.91%	0.127%
39	14.633	2197	2201	2209	rVB	31987	41426	6.60%	0.440%
40	14.798	2224	2229	2232	rBV	362205	441124	70.23%	4.682%
41	14.833	2232	2235	2244	rVB	396733	499855	79.58%	5.306%
42	14.927	2247	2251	2261	rVB	76008	110663	17.62%	1.175%
43	15.033	2263	2269	2272	rBV4	6973	12198	1.94%	0.129%
44	15.227	2299	2302	2310	rVV	22675	33438	5.32%	0.355%
45	15.327	2316	2319	2324	rVB	23497	26218	4.17%	0.278%
46	15.468	2338	2343	2347	rVV3	13496	22147	3.53%	0.235%
47	15.592	2360	2364	2368	rVB	76782	81841	13.03%	0.869%
48	15.633	2368	2371	2378	rVB	80177	88966	14.16%	0.944%
49	15.710	2378	2384	2386	rBV2	24428	27965	4.45%	0.297%
50	15.745	2386	2390	2394	rVV2	86116	129857	20.67%	1.378%
51	15.786	2394	2397	2406	rVB	54069	73357	11.68%	0.779%
52	16.039	2432	2440	2445	rBV	31625	52572	8.37%	0.558%
53	16.121	2450	2454	2461	rBV6	6860	14702	2.34%	0.156%
54	16.210	2466	2469	2472	rVB4	9548	11786	1.88%	0.125%
55	16.251	2472	2476	2481	rBV	13099	19682	3.13%	0.209%
56	16.298	2481	2484	2488	rVV2	19005	22089	3.52%	0.234%
57	16.327	2488	2489	2493	rVB	10723	10839	1.73%	0.115%
58	16.398	2497	2501	2504	rBV2	59706	75554	12.03%	0.802%
59	16.439	2504	2508	2512	rVV3	33520	48822	7.77%	0.518%
60	16.474	2512	2514	2517	rVV2	16181	14382	2.29%	0.153%
61	16.515	2517	2521	2526	rVV4	24075	35231	5.61%	0.374%
62	16.598	2531	2535	2541	rBV	264092	286262	45.57%	3.038%
63	16.698	2549	2552	2555	rBV4	17052	19431	3.09%	0.206%
64	16.739	2555	2559	2564	rVB4	15624	21382	3.40%	0.227%
65	16.798	2564	2569	2573	rBV3	13379	17973	2.86%	0.191%
66	16.904	2582	2587	2592	rVV	281660	332366	52.91%	3.528%
67	17.027	2605	2608	2612	rVB	16812	18838	3.00%	0.200%
68	17.098	2616	2620	2624	rVV4	8696	12540	2.00%	0.133%
69	17.139	2624	2627	2634	rVB	244577	287510	45.77%	3.052%
70	17.227	2637	2642	2647	rBV4	25654	47826	7.61%	0.508%
71	17.339	2656	2661	2664	rBV5	27053	50568	8.05%	0.537%

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72	17.368	2664	2666	2671	rVB	51096	61903	9.86%	0.657%
73	17.421	2671	2675	2677	rBV4	11514	14308	2.28%	0.152%
74	17.457	2678	2681	2685	rVB	41954	45191	7.19%	0.480%
75	17.504	2685	2689	2693	rBV2	44564	61441	9.78%	0.652%
76	17.539	2693	2695	2700	rVB5	12784	13905	2.21%	0.148%
77	17.621	2705	2709	2713	rVV	42397	59301	9.44%	0.629%
78	17.662	2713	2716	2721	rVV	31839	48141	7.66%	0.511%
79	17.727	2725	2727	2731	rVB5	9080	11598	1.85%	0.123%
80	17.804	2736	2740	2742	rBV4	8053	11255	1.79%	0.119%
81	17.862	2747	2750	2753	rBV5	11430	17534	2.79%	0.186%
82	18.015	2773	2776	2781	rBV7	11749	23913	3.81%	0.254%
83	18.115	2791	2793	2797	rBV3	10701	10925	1.74%	0.116%
84	18.174	2797	2803	2806	rBV2	29085	55470	8.83%	0.589%
85	18.204	2806	2808	2811	rVV2	28110	28353	4.51%	0.301%
86	18.239	2811	2814	2817	rVV2	11774	15727	2.50%	0.167%
87	18.274	2818	2820	2823	rVB4	13996	15891	2.53%	0.169%
88	18.345	2829	2832	2836	rVB3	22040	26117	4.16%	0.277%
89	18.398	2838	2841	2844	rBV5	13362	21501	3.42%	0.228%
90	18.492	2852	2857	2861	rBV2	477301	628137	100.00%	6.667%
91	18.533	2861	2864	2870	rVB2	79603	131790	20.98%	1.399%
92	18.621	2876	2879	2884	rVB6	21462	27747	4.42%	0.295%
93	18.721	2894	2896	2901	rBV5	18307	25165	4.01%	0.267%
94	19.003	2942	2944	2948	rVB2	33332	34728	5.53%	0.369%
95	19.045	2949	2951	2954	rVB	25444	26101	4.16%	0.277%
96	19.156	2968	2970	2972	rBV3	19658	22780	3.63%	0.242%
97	19.768	3072	3074	3078	rBV	54255	68350	10.88%	0.725%
98	20.062	3121	3124	3127	rBV	47827	52231	8.32%	0.554%
99	20.121	3131	3134	3136	rVV	57066	58619	9.33%	0.622%
100	20.168	3138	3142	3148	rVV	298957	363314	57.84%	3.856%

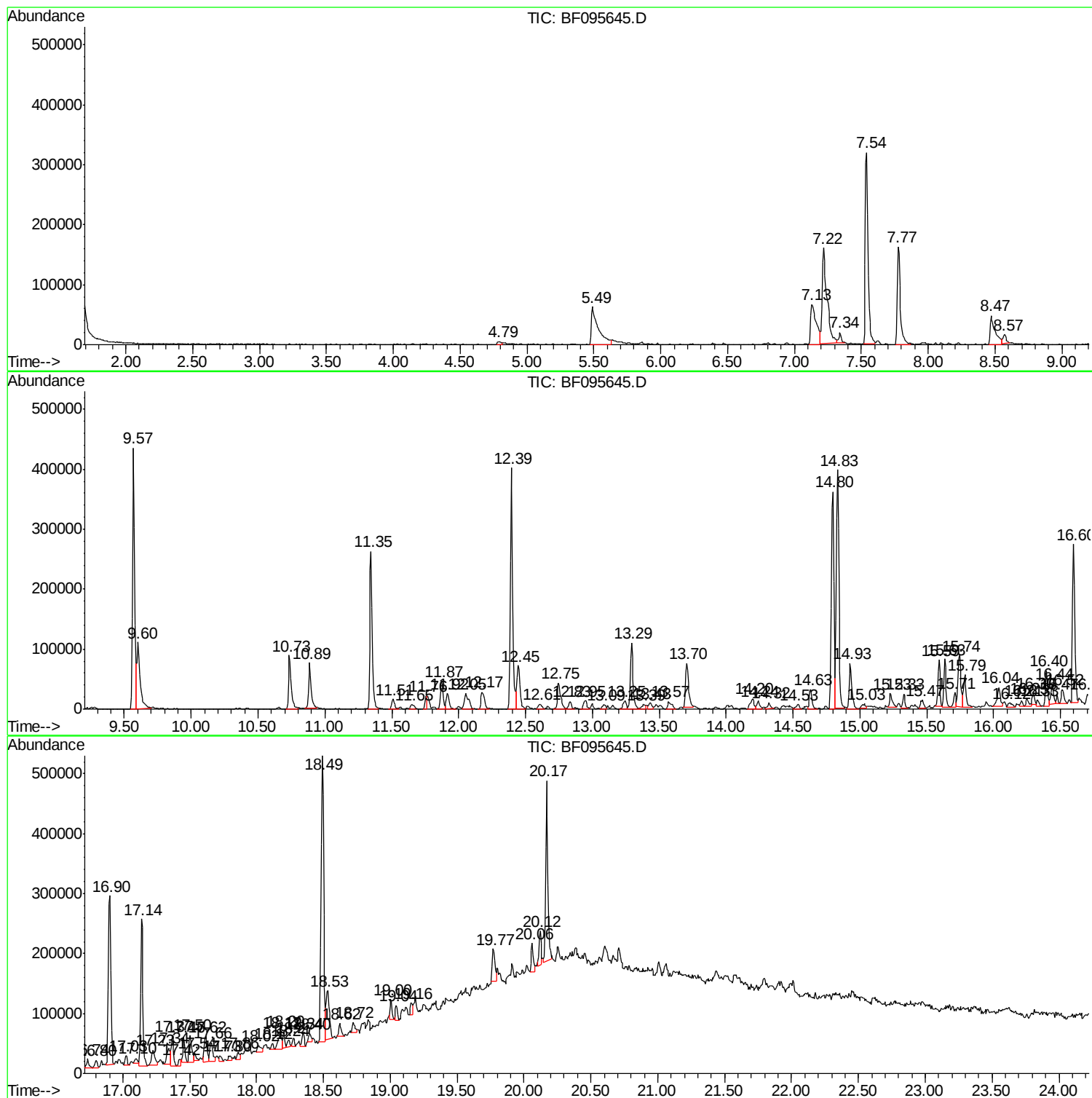
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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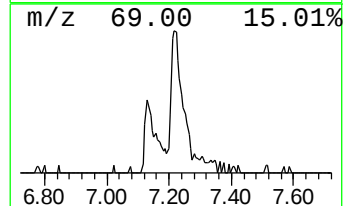
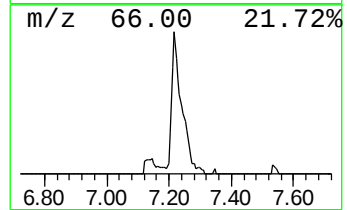
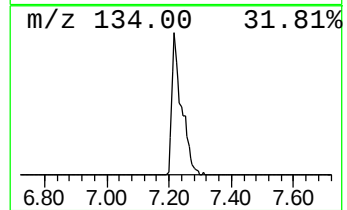
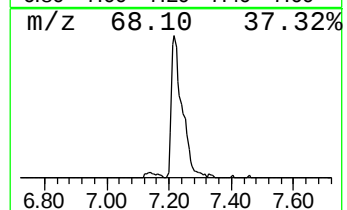
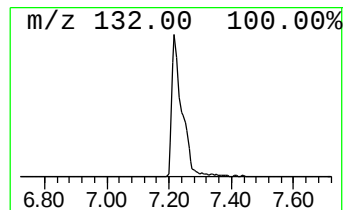
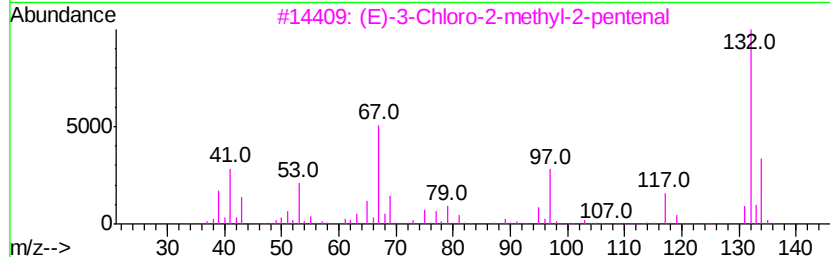
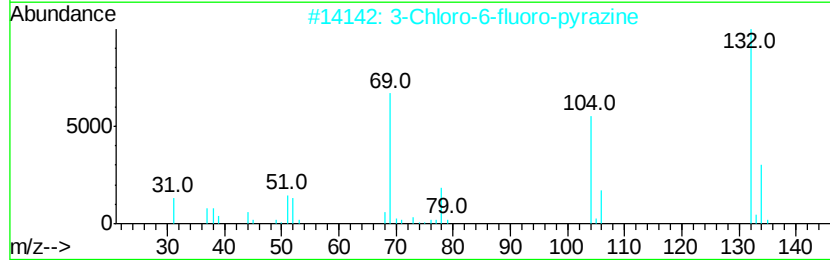
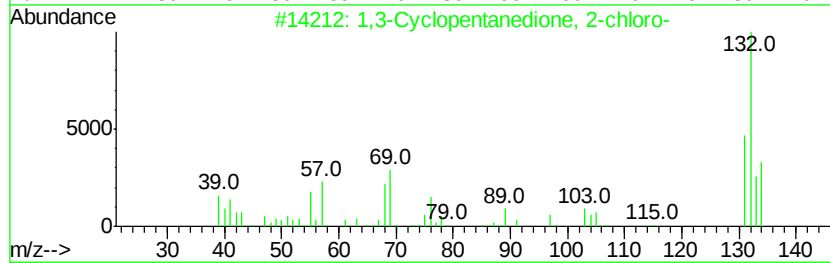
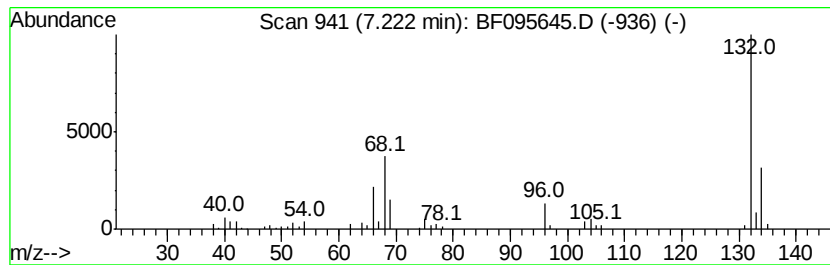
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown7.22 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.22	17.79 ng	395736	1,4-Dichlorobenzene-d4	7.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	32
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
4		Benzeneethanamine, .alpha.-methy...	223	C16H17N	002980-02-1	12
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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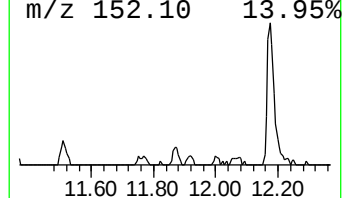
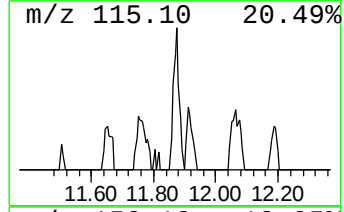
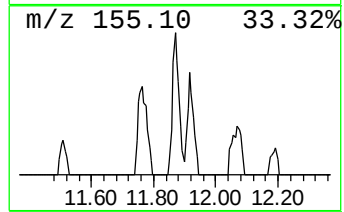
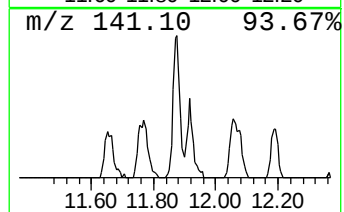
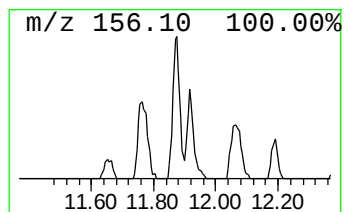
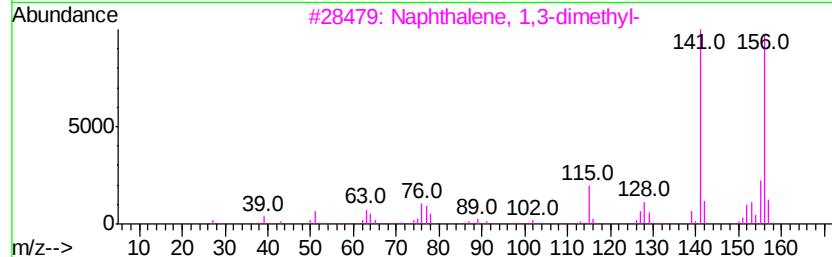
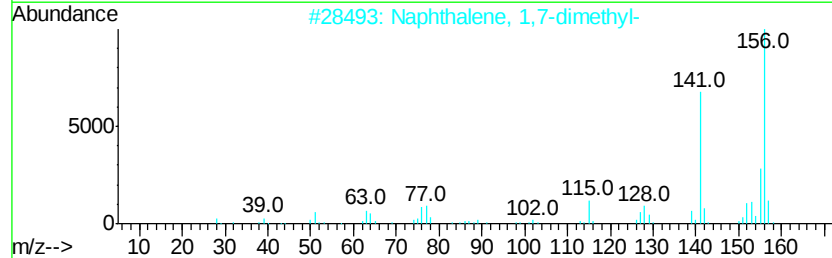
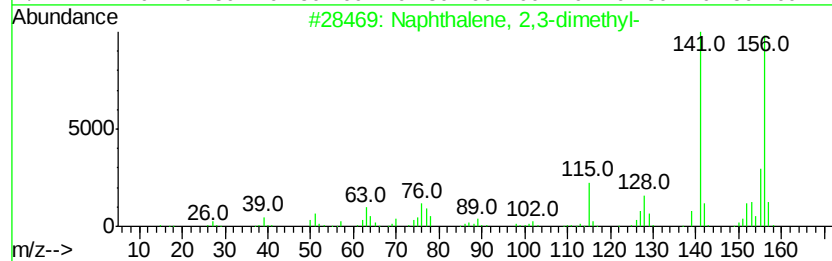
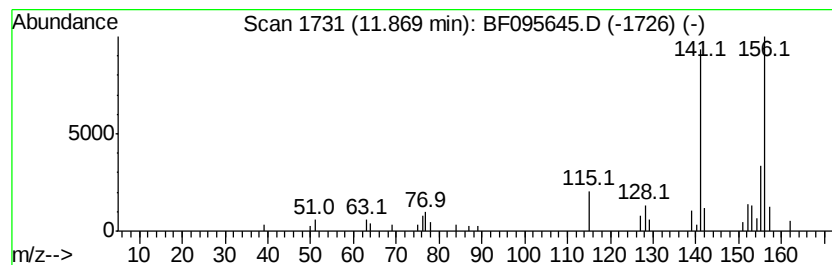
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Naphthalene, 2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.87	2.60 ng	68320	Acenaphthene-d10	12.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
3		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
4		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
5		Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	93



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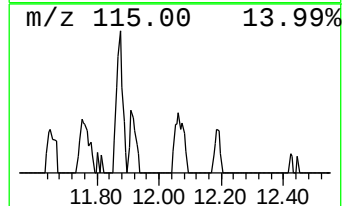
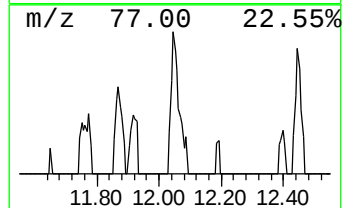
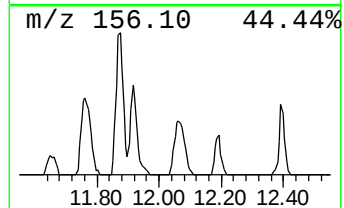
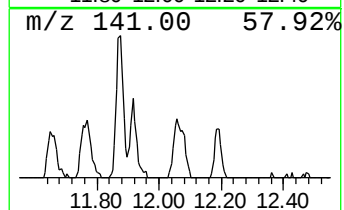
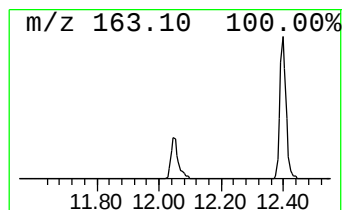
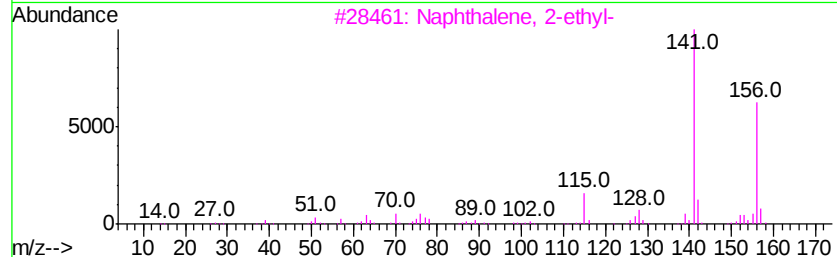
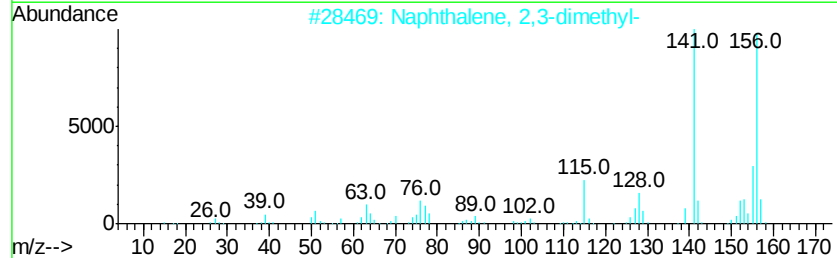
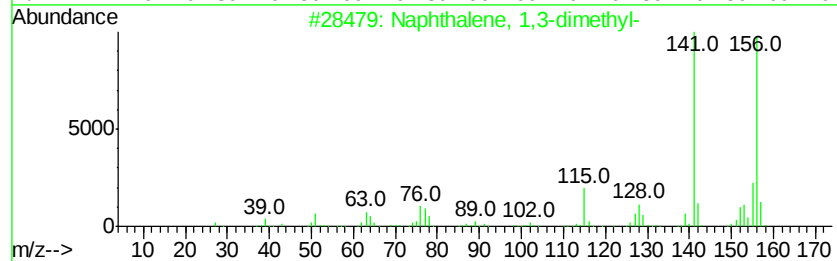
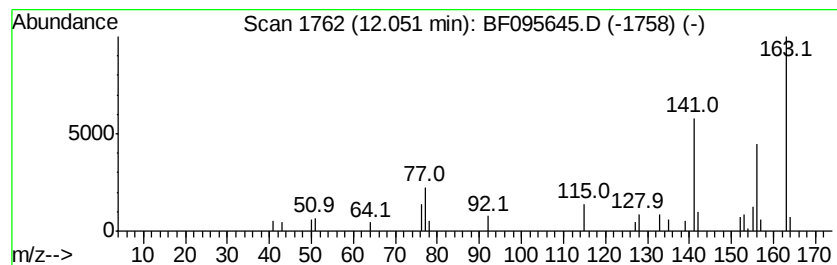
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Peak Number 5 Naphthalene, 1,3-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.05	2.13 ng	55958	Acenaphthene-d10	12.39
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,3-dimethyl-	156 C12H12	000575-41-7 55
2		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 50
3		Naphthalene, 2-ethyl-	156 C12H12	000939-27-5 46
4		Naphthalene, 1-ethyl-	156 C12H12	001127-76-0 46
5		Benzonitrile, 3,4-dimethoxy-	163 C9H9NO2	002024-83-1 38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060217\
Data File : BF095645.D
Acq On : 2 Jun 2017 21:09
Operator : SJ/MA
Sample : I3441-11 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OR-RO-17273-060117

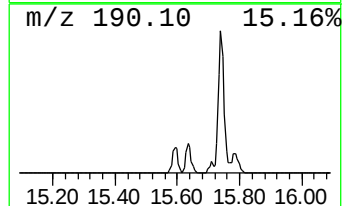
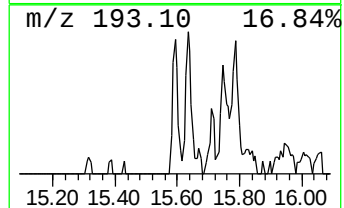
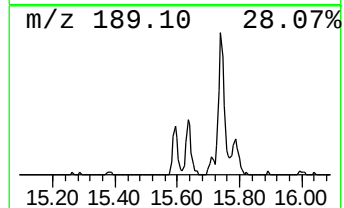
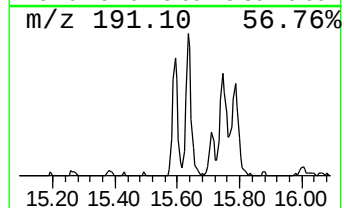
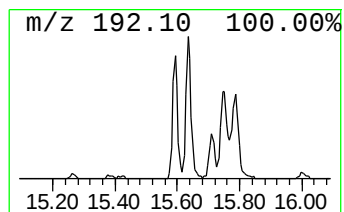
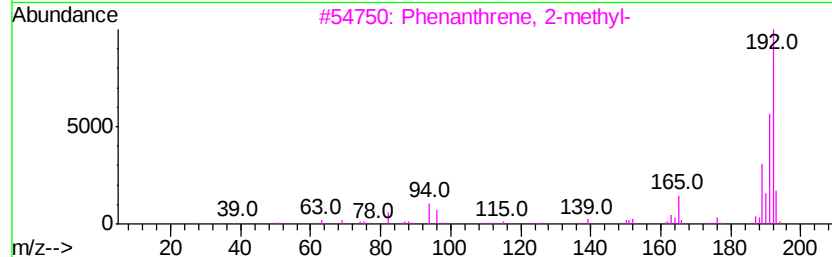
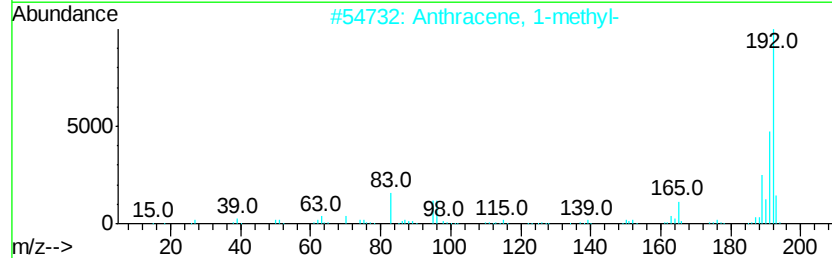
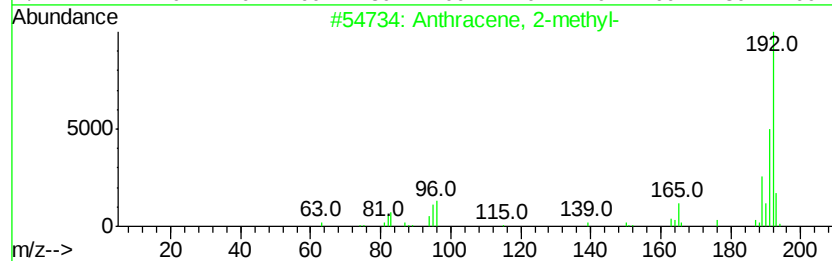
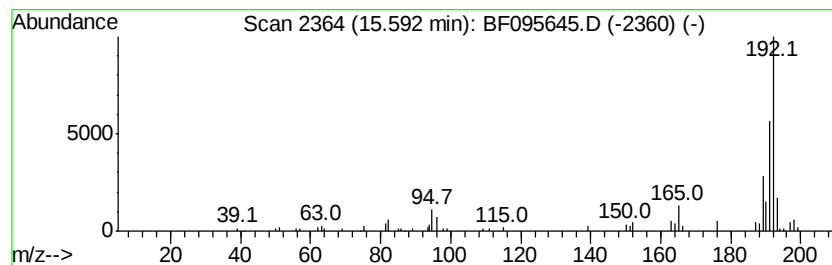
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Anthracene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.59	3.71 ng	81841	Phenanthrene-d10	14.80

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



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060217\
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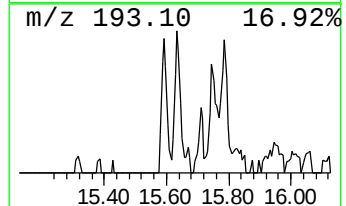
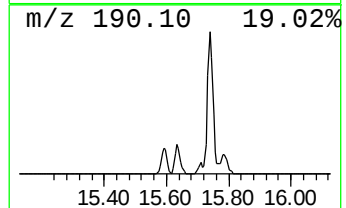
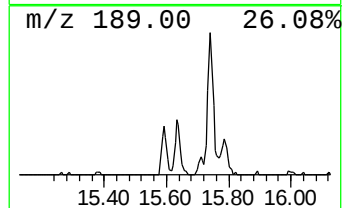
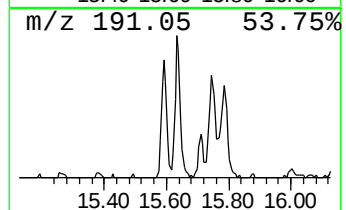
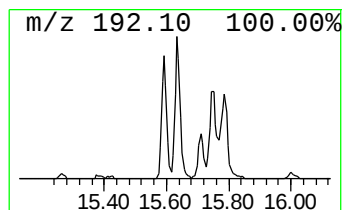
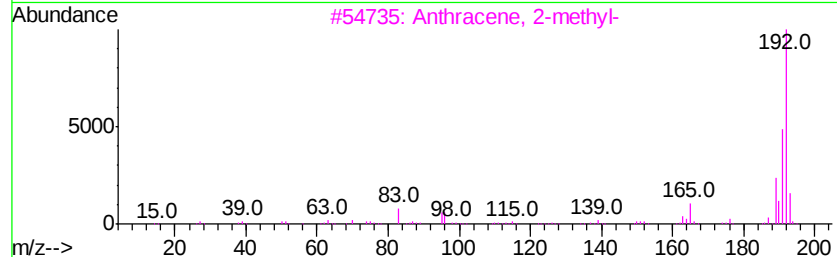
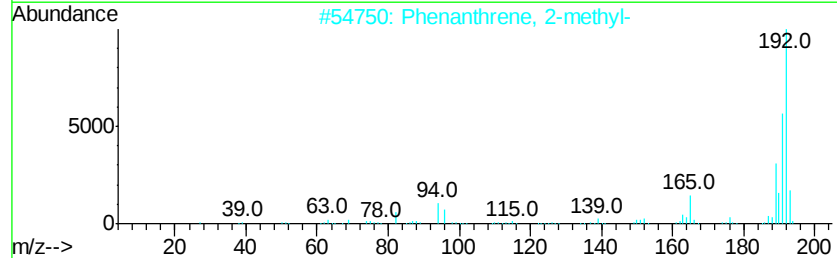
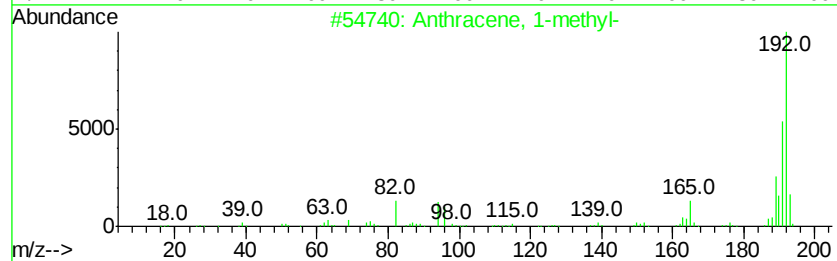
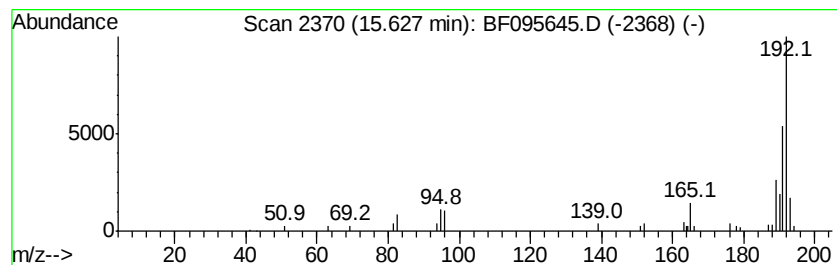
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Anthracene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.63	4.03 ng	88966	Phenanthrene-d10	14.80	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 1-methyl-	192	C15H12	000610-48-0	98
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
5	1H-Cyclopropa[1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060217\
Data File : BF095645.D
Acq On : 2 Jun 2017 21:09
Operator : SJ/MA
Sample : I3441-11 5X
Misc :
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Instrument :
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ClientSampleId :
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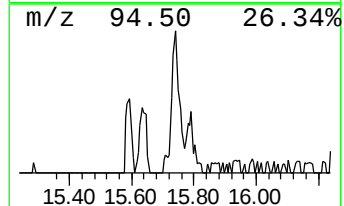
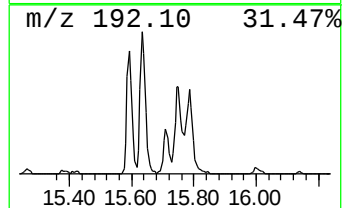
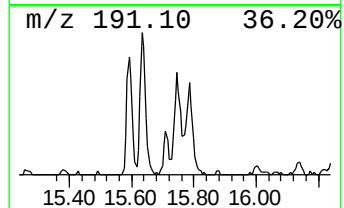
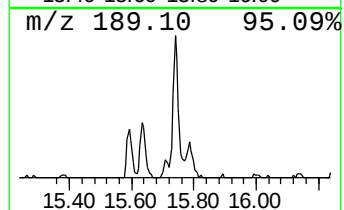
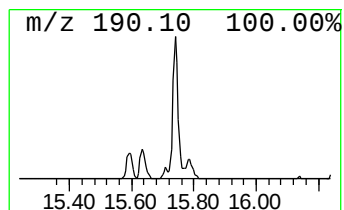
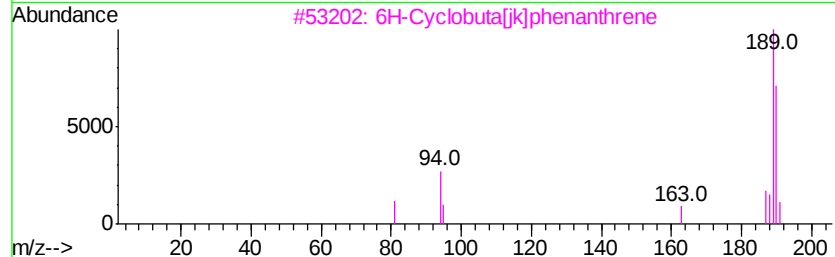
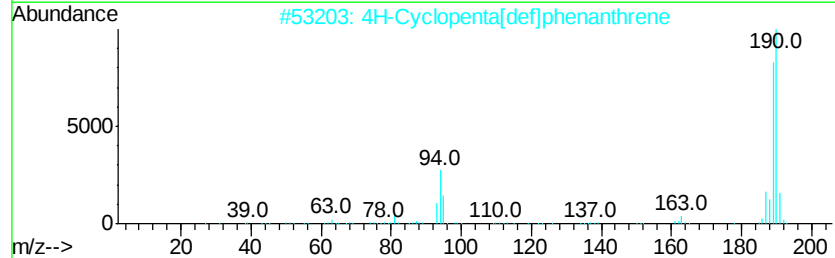
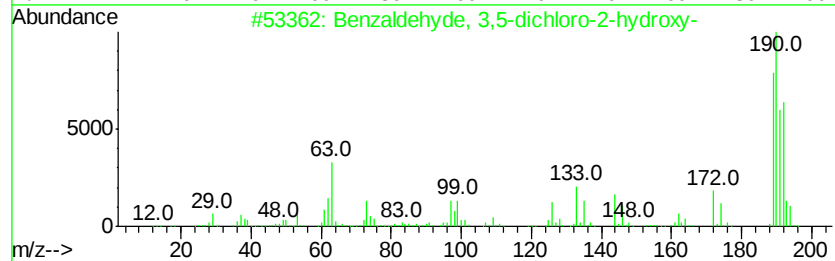
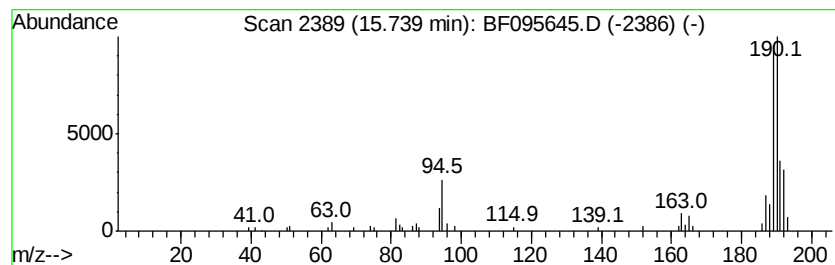
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzaldehyde, 3,5-dichloro-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.74	5.89 ng	129857	Phenanthrene-d10	14.80

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzaldehydve, 3,5-dichloro-2-hvd...	190	C7H4Cl2O2	000090-60-8	72
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
3		6H-Cyclobuta[fik]phenanthrene	190	C15H10	083469-43-6	49
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	46
5		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060217\
Data File : BF095645.D
Acq On : 2 Jun 2017 21:09
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ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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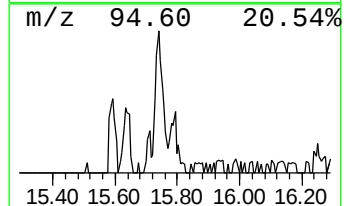
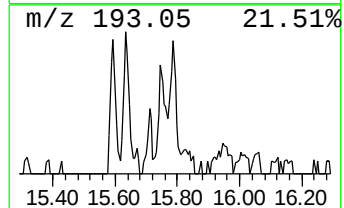
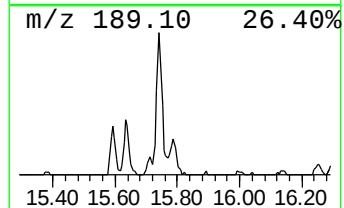
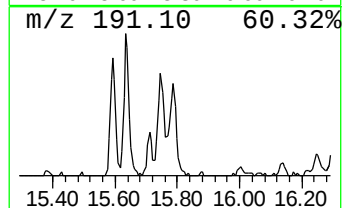
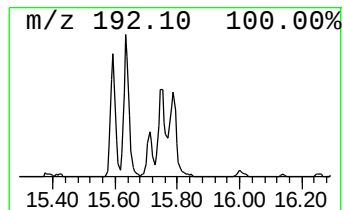
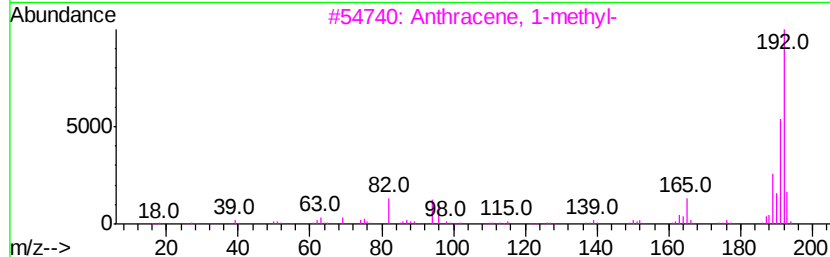
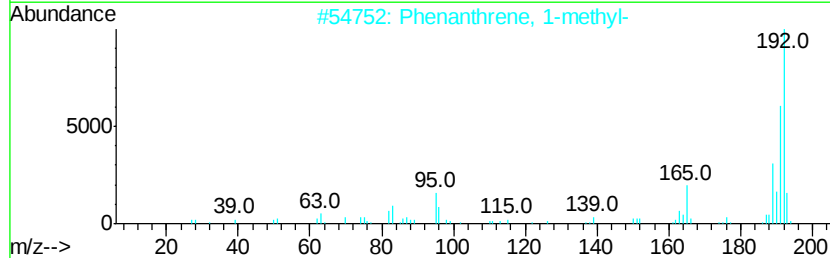
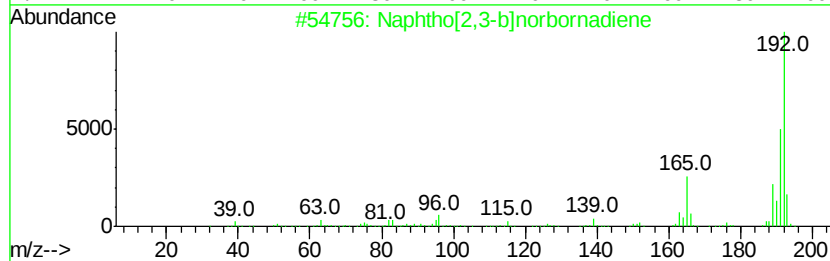
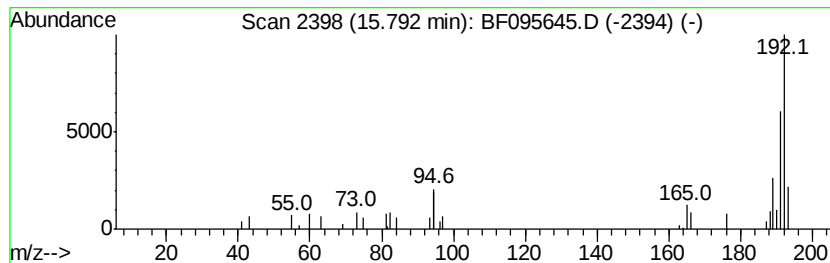
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Naphtho[2,3-b]norbornadiene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.79	3.33 ng	73357	Phenanthrene-d10	14.80

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	70
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	64
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	62
4		5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	62
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	58



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060217\
Data File : BF095645.D
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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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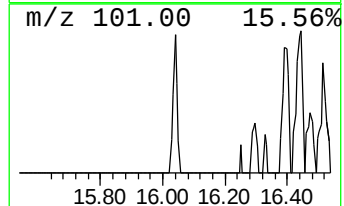
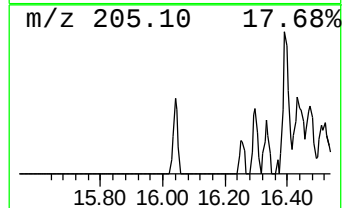
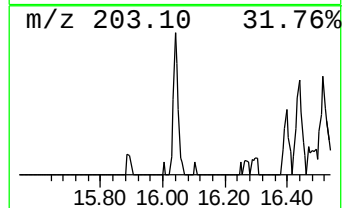
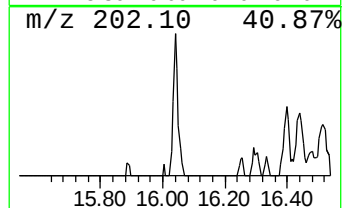
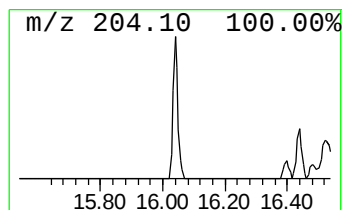
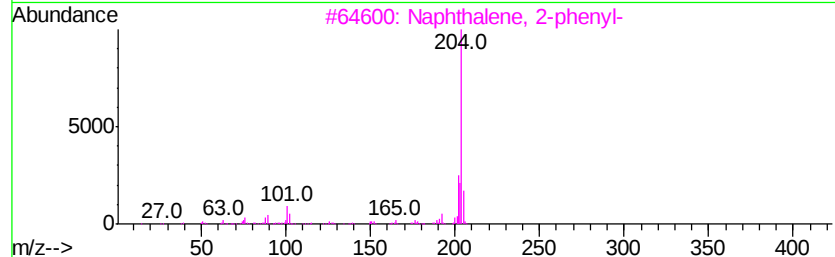
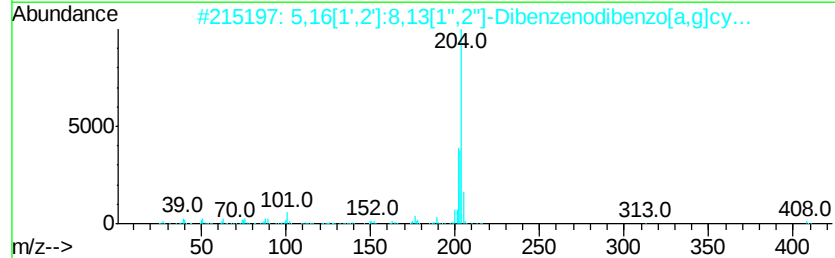
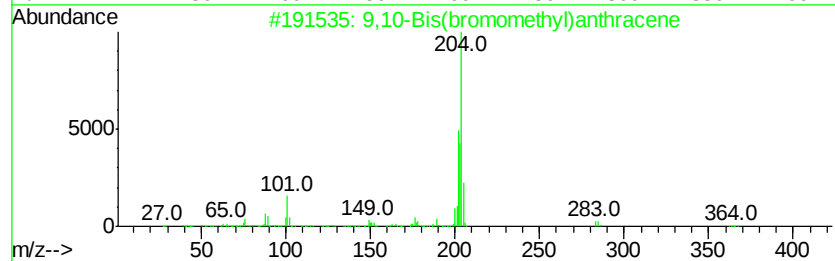
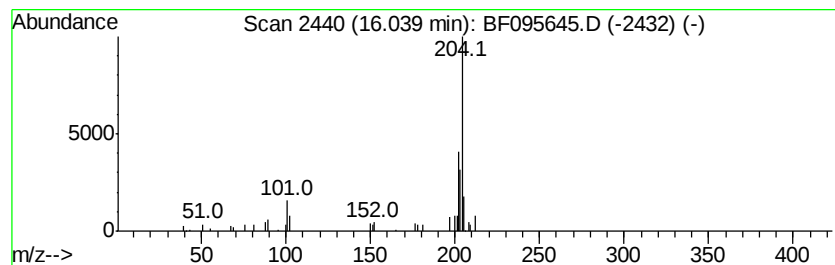
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 9,10-Bis(bromomethyl)anthra... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.04	2.38 ng	52572	Phenanthrene-d10	14.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	87
2		5,16[1',2'] : 8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	78
3		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	70
4		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	62
5		Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	53



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060217\
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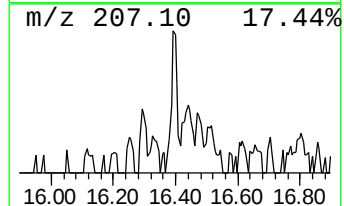
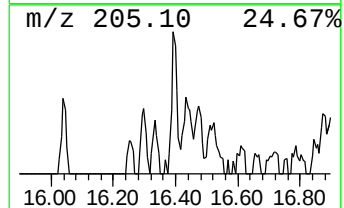
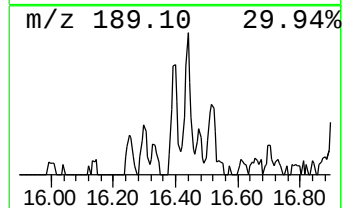
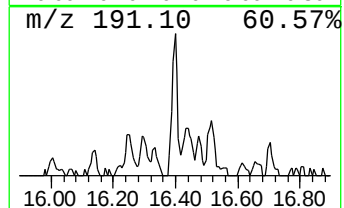
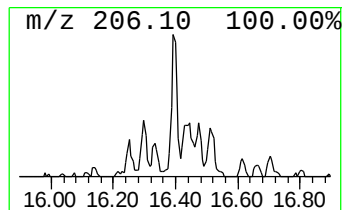
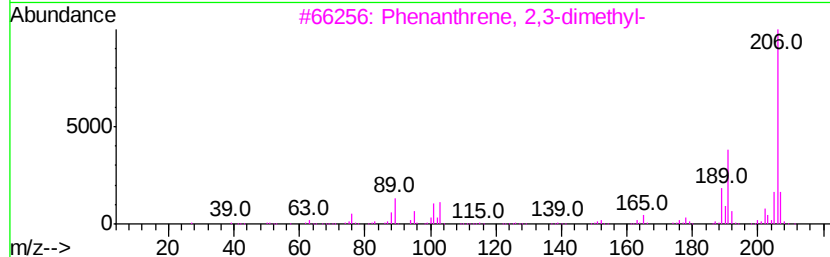
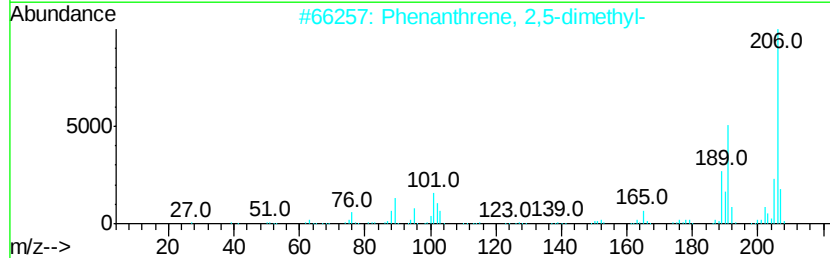
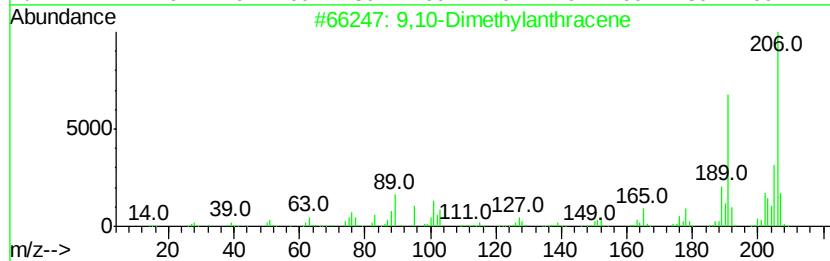
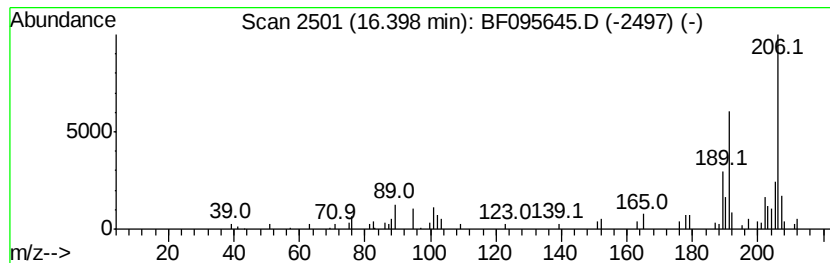
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 9,10-Dimethylantracene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.40	3.43 ng	75554	Phenanthrene-d10	14.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Dimethylantracene	206	C16H14	000781-43-1	98
2		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	95
3		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	90
4		1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	90
5		Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	89



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060217\
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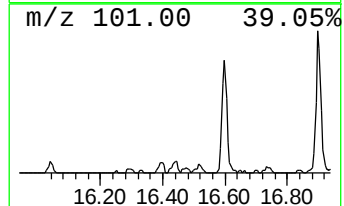
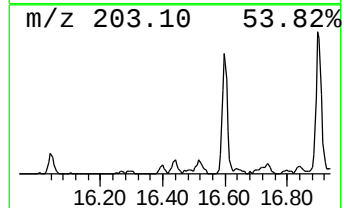
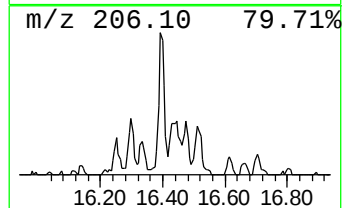
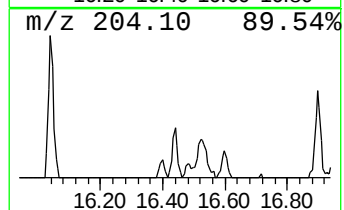
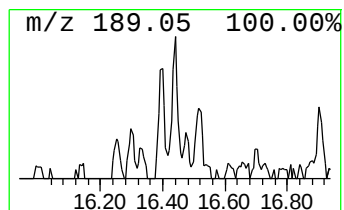
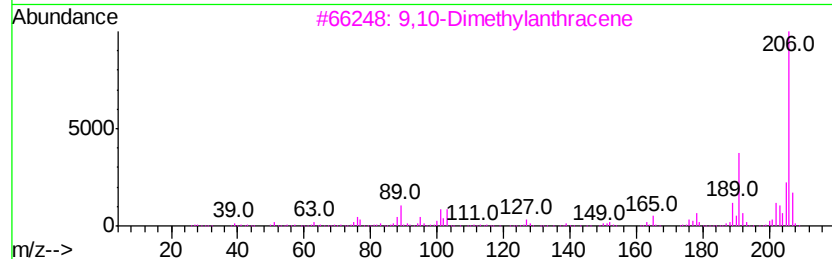
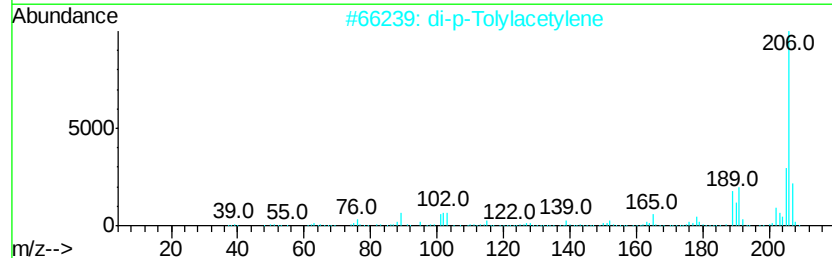
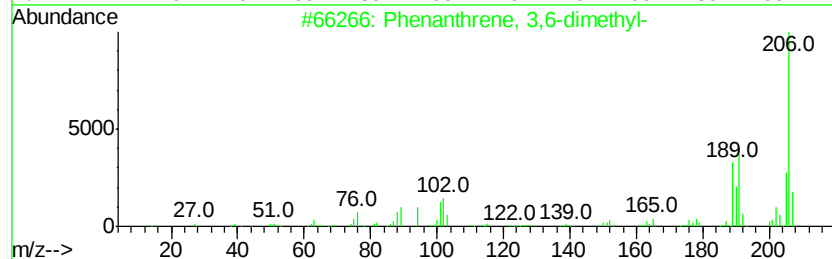
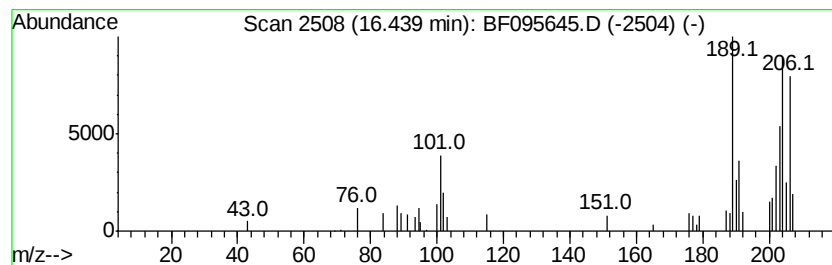
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown16.44 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.44	2.21 ng	48822	Phenanthrene-d10	14.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	46
2			di-p-Tolylacetylene	206	C16H14	002789-88-0	42
3			9,10-Dimethylanthracene	206	C16H14	000781-43-1	38
4			Levamisole	204	C11H12N2S	014769-73-4	30
5			Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	30



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060217\
Data File : BF095645.D
Acq On : 2 Jun 2017 21:09
Operator : SJ/MA
Sample : I3441-11 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OR-RO-17273-060117

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.22	7.22	17.8	ng	395736	1	7.54	444807	20.0
Naphthalene, 2,3-...	11.87	2.6	ng	68320	3	12.39	526416	20.0
Naphthalene, 1,3-...	12.05	2.1	ng	55958	3	12.39	526416	20.0
Anthracene, 2-met...	15.59	3.7	ng	81841	4	14.80	441124	20.0
Anthracene, 1-met...	15.63	4.0	ng	88966	4	14.80	441124	20.0
Benzaldehyde, 3,5...	15.74	5.9	ng	129857	4	14.80	441124	20.0
Naphtho[2,3-b]nor...	15.79	3.3	ng	73357	4	14.80	441124	20.0
9,10-Bis(bromomet...	16.04	2.4	ng	52572	4	14.80	441124	20.0
9,10-Dimethylanth...	16.40	3.4	ng	75554	4	14.80	441124	20.0
unknown16.44	16.44	2.2	ng	48822	4	14.80	441124	20.0