

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060217\
 Data File : BF095651.D
 Acq On : 3 Jun 2017 00:22
 Operator : SJ/MA
 Sample : I3441-03 5X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-RO-5004-060117

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	5.487	643	646	672	rBV	61481	226186	33.74%	2.513%
2	7.128	922	925	936	rBV	65922	184277	27.49%	2.048%
3	7.216	936	940	956	rVB	154389	357371	53.31%	3.971%
4	7.540	990	995	1007	rBV	332093	444997	66.39%	4.944%
5	7.775	1031	1035	1051	rBV	159480	240522	35.88%	2.672%
6	8.469	1150	1153	1168	rBV	47679	131810	19.66%	1.465%
7	9.569	1336	1340	1345	rBV	441992	569257	84.92%	6.325%
8	10.739	1535	1539	1551	rBV	46765	69251	10.33%	0.769%
9	10.886	1558	1564	1575	rBV	42430	65095	9.71%	0.723%
10	11.345	1638	1642	1655	rBV	286624	343327	51.22%	3.815%
11	11.657	1691	1695	1700	rBV	6544	10690	1.59%	0.119%
12	11.757	1709	1712	1713	rBV	15720	15605	2.33%	0.173%
13	11.775	1713	1715	1719	rVB2	14036	15918	2.37%	0.177%
14	11.875	1727	1732	1736	rBV	33881	49027	7.31%	0.545%
15	11.916	1736	1739	1748	rVB	20438	30364	4.53%	0.337%
16	12.051	1758	1762	1766	rBV2	20884	38658	5.77%	0.430%
17	12.180	1779	1784	1791	rBV2	23960	39923	5.96%	0.444%
18	12.398	1816	1821	1827	rBV2	433983	557881	83.23%	6.199%
19	12.451	1827	1830	1839	rVB	42285	61338	9.15%	0.682%
20	12.616	1852	1858	1863	rBV3	5756	11200	1.67%	0.124%
21	12.751	1874	1881	1890	rBV2	21873	41891	6.25%	0.465%
22	12.833	1890	1895	1899	rVB3	11080	14497	2.16%	0.161%
23	12.951	1910	1915	1920	rBV3	12790	18783	2.80%	0.209%
24	13.092	1933	1939	1942	rBV3	6032	10232	1.53%	0.114%
25	13.245	1959	1965	1968	rBV4	13175	21062	3.14%	0.234%
26	13.298	1968	1974	1984	rVV	67068	98084	14.63%	1.090%
27	13.386	1984	1989	1992	rVV4	8999	16798	2.51%	0.187%
28	13.439	1994	1998	2001	rVV3	7829	11320	1.69%	0.126%
29	13.480	2001	2005	2008	rVV	6728	9911	1.48%	0.110%
30	13.574	2016	2021	2022	rVV2	9471	9841	1.47%	0.109%
31	13.598	2022	2025	2030	rVB5	8579	14415	2.15%	0.160%
32	13.704	2040	2043	2052	rBV2	68728	114830	17.13%	1.276%
33	14.016	2090	2096	2097	rBV2	10538	12511	1.87%	0.139%
34	14.033	2097	2099	2103	rVB3	9269	12353	1.84%	0.137%

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

35	14.198	2119	2127	2131	rBV3	15172	28838	4.30%	0.320%
36	14.239	2131	2134	2140	rVV4	11679	18845	2.81%	0.209%
37	14.321	2140	2148	2153	rVV4	9206	15842	2.36%	0.176%
38	14.427	2161	2166	2169	rBV4	8520	10914	1.63%	0.121%
39	14.633	2198	2201	2207	rVB	18809	24877	3.71%	0.276%
40	14.745	2215	2220	2224	rBV4	7681	11965	1.78%	0.133%
41	14.798	2224	2229	2232	rVV	389475	453287	67.62%	5.037%
42	14.839	2232	2236	2245	rVB	258338	356886	53.24%	3.965%
43	14.927	2248	2251	2258	rBV	63115	89055	13.29%	0.990%
44	15.021	2262	2267	2271	rBV3	10575	17755	2.65%	0.197%
45	15.233	2300	2303	2307	rBV	13498	20283	3.03%	0.225%
46	15.333	2317	2320	2327	rBV	18930	23216	3.46%	0.258%
47	15.415	2331	2334	2338	rVB3	8845	10249	1.53%	0.114%
48	15.463	2338	2342	2348	rBV3	11729	22144	3.30%	0.246%
49	15.592	2361	2364	2368	rVB	57704	63281	9.44%	0.703%
50	15.639	2368	2372	2379	rVB	67363	83850	12.51%	0.932%
51	15.710	2380	2384	2386	rBV2	20804	23224	3.46%	0.258%
52	15.745	2386	2390	2394	rVV3	77953	110841	16.54%	1.232%
53	15.786	2394	2397	2403	rVB	46737	61596	9.19%	0.684%
54	15.945	2421	2424	2428	rBV3	7207	11241	1.68%	0.125%
55	16.015	2431	2436	2438	rBV3	12150	17266	2.58%	0.192%
56	16.039	2438	2440	2451	rVB3	24059	44562	6.65%	0.495%
57	16.251	2473	2476	2480	rVB2	12412	14029	2.09%	0.156%
58	16.298	2480	2484	2488	rBV	24335	29757	4.44%	0.331%
59	16.333	2488	2490	2494	rVB	12643	13929	2.08%	0.155%
60	16.398	2494	2501	2504	rBV2	63904	82942	12.37%	0.922%
61	16.445	2504	2509	2512	rVV3	30079	47545	7.09%	0.528%
62	16.474	2512	2514	2517	rVB3	15665	15207	2.27%	0.169%
63	16.533	2517	2524	2527	rBV5	32709	64486	9.62%	0.717%
64	16.568	2527	2530	2532	rBV3	30788	27863	4.16%	0.310%
65	16.598	2532	2535	2541	rVB	270141	317732	47.40%	3.530%
66	16.645	2541	2543	2548	rVB3	8669	9142	1.36%	0.102%
67	16.704	2548	2553	2556	rBV5	18953	25780	3.85%	0.286%
68	16.739	2556	2559	2566	rVB3	13932	19747	2.95%	0.219%
69	16.804	2566	2570	2573	rBV2	13547	16825	2.51%	0.187%
70	16.904	2583	2587	2592	rBV	347288	376689	56.20%	4.185%
71	16.974	2597	2599	2604	rVB3	10665	13466	2.01%	0.150%

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72	17.027	2604	2608	2612	rVB3	22569	27487	4.10%	0.305%
73	17.068	2612	2615	2617	rBV3	15874	17137	2.56%	0.190%
74	17.104	2617	2621	2624	rBV2	12624	12839	1.92%	0.143%
75	17.145	2624	2628	2636	rVB	269139	304352	45.40%	3.382%
76	17.233	2637	2643	2646	rBV3	29606	51305	7.65%	0.570%
77	17.345	2658	2662	2664	rBV4	28091	42478	6.34%	0.472%
78	17.374	2664	2667	2672	rVB	64879	77499	11.56%	0.861%
79	17.457	2678	2681	2686	rVB2	47619	55068	8.22%	0.612%
80	17.504	2686	2689	2693	rBV2	49999	68021	10.15%	0.756%
81	17.598	2699	2705	2706	rBV6	10091	17210	2.57%	0.191%
82	17.621	2706	2709	2712	rVV	43101	44583	6.65%	0.495%
83	17.662	2713	2716	2722	rVV	27844	36282	5.41%	0.403%
84	17.809	2736	2741	2743	rBV6	10451	14518	2.17%	0.161%
85	17.986	2768	2771	2773	rBV2	16050	16360	2.44%	0.182%
86	18.015	2773	2776	2781	rBV6	18008	33950	5.06%	0.377%
87	18.121	2792	2794	2796	rBV3	10239	10314	1.54%	0.115%
88	18.180	2800	2804	2806	rVV2	25851	38150	5.69%	0.424%
89	18.204	2806	2808	2812	rVB2	28726	28448	4.24%	0.316%
90	18.239	2812	2814	2817	rBV	21651	21142	3.15%	0.235%
91	18.351	2830	2833	2836	rVB4	22808	26492	3.95%	0.294%
92	18.404	2838	2842	2847	rBV7	32835	54959	8.20%	0.611%
93	18.498	2852	2858	2861	rBV2	512738	670310	100.00%	7.448%
94	18.533	2861	2864	2870	rVB	114580	179521	26.78%	1.995%
95	18.627	2878	2880	2883	rVB3	34695	33052	4.93%	0.367%
96	19.009	2942	2945	2948	rVB2	29187	29769	4.44%	0.331%
97	19.774	3071	3075	3087	rBV	90933	178515	26.63%	1.984%
98	20.062	3122	3124	3127	rBV	59732	54500	8.13%	0.606%
99	20.121	3132	3134	3139	rVV	82479	83280	12.42%	0.925%
100	20.174	3139	3143	3147	rBV	273416	308043	45.96%	3.423%

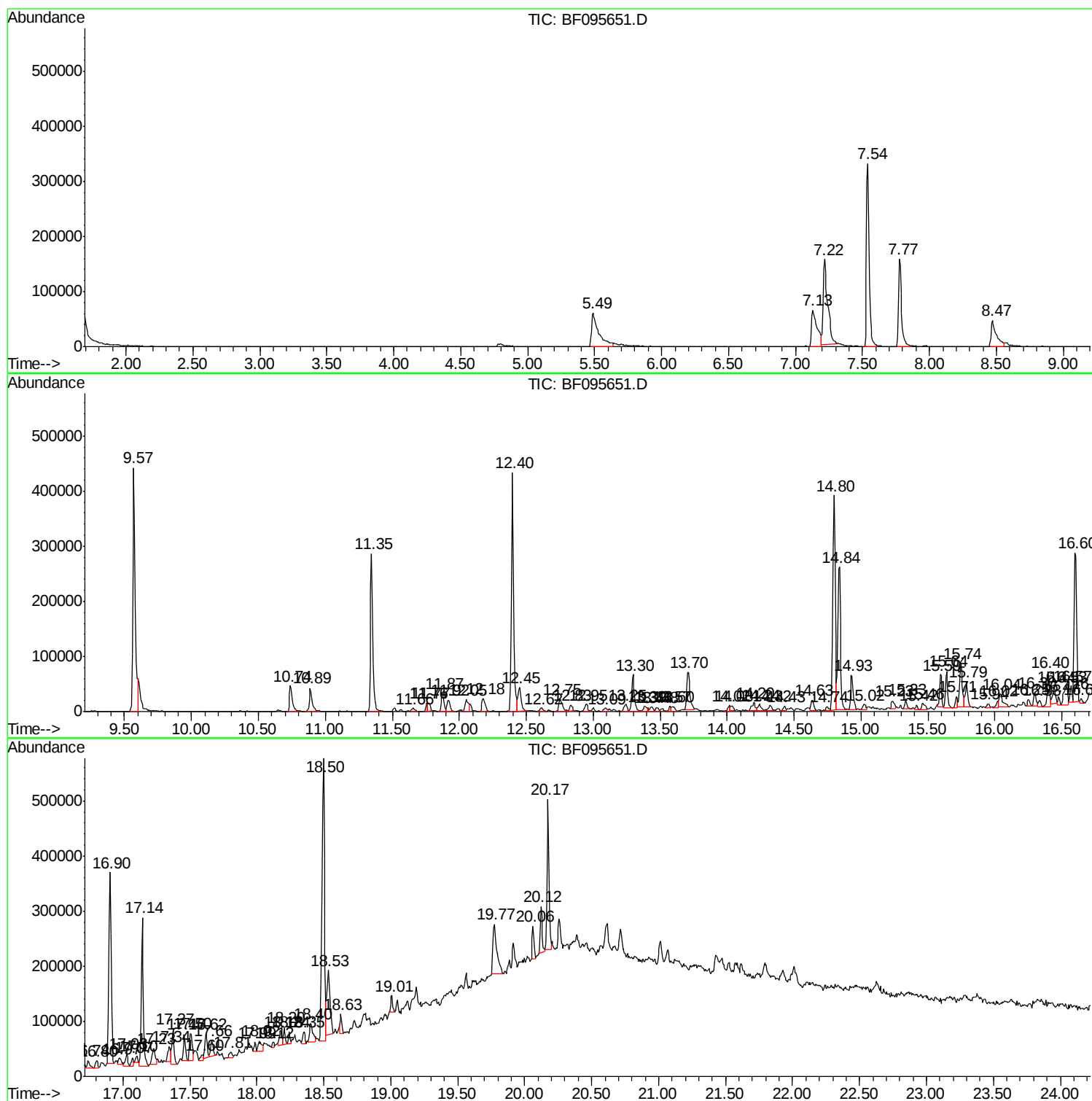
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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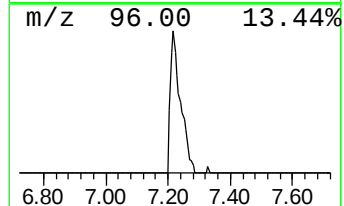
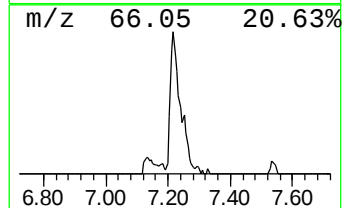
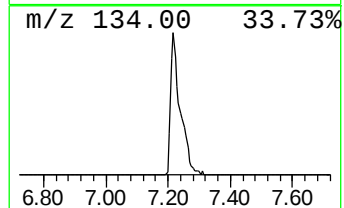
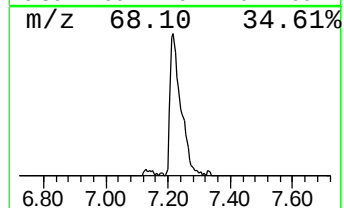
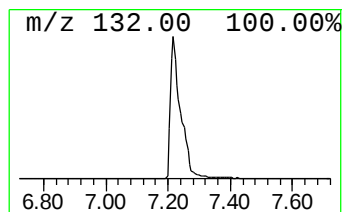
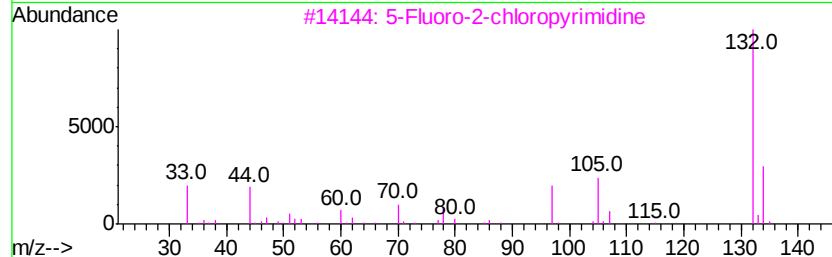
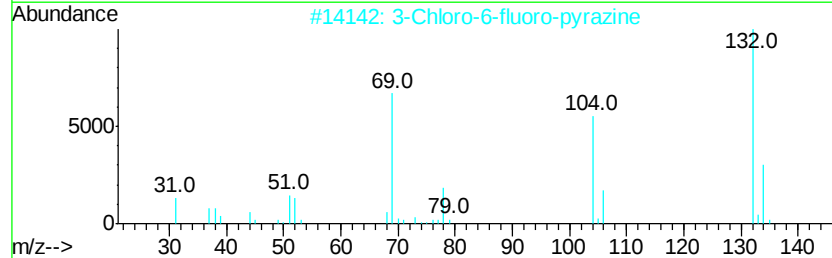
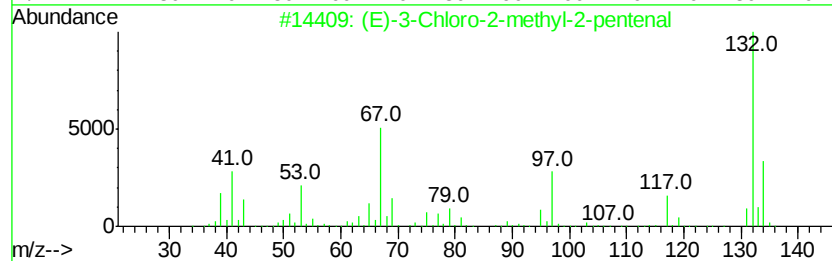
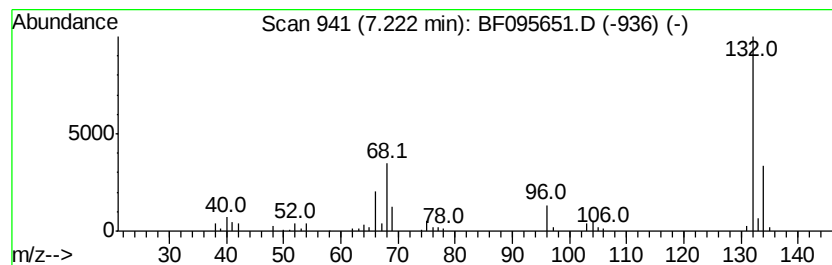
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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	16.06 ng	357371	1,4-Dichlorobenzene-d4	7.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	17
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
5		Xenon	132	Xe	007440-63-3	9



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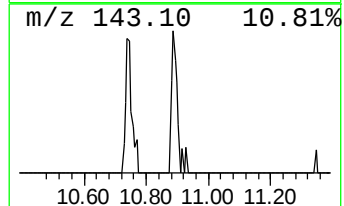
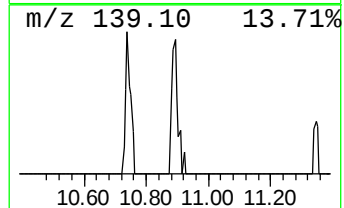
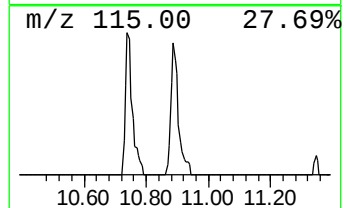
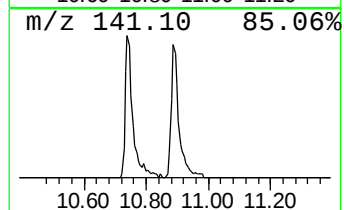
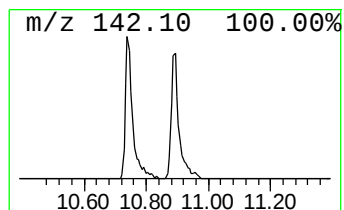
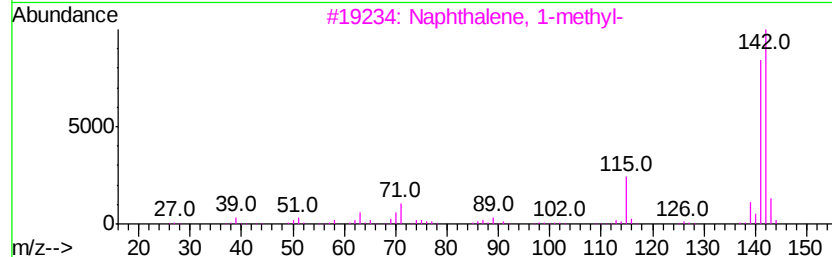
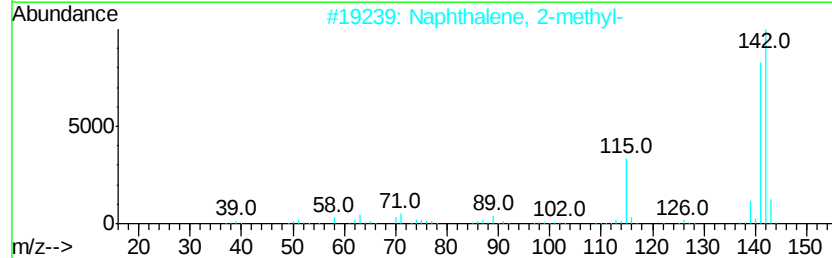
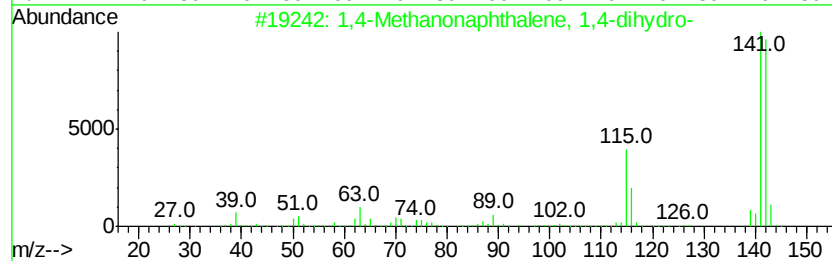
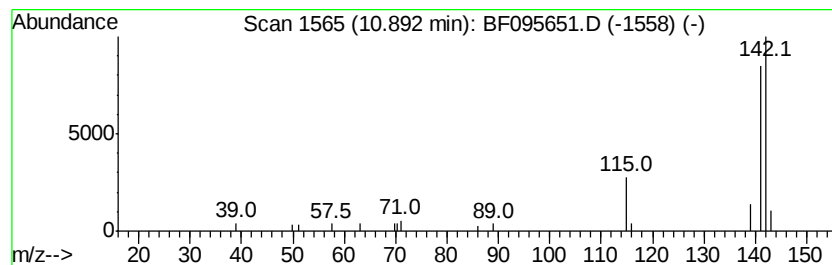
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1,4-Methanonaphthalene, 1,4... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.89	2.29 ng	65095	Naphthalene-d8	9.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	91
3		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	87
4		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	83
5		Benzocycloheptatriene	142	C11H10	000264-09-5	64



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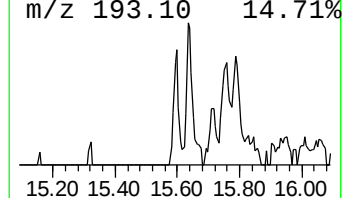
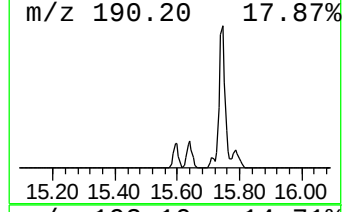
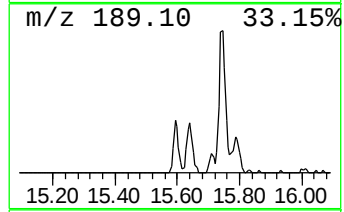
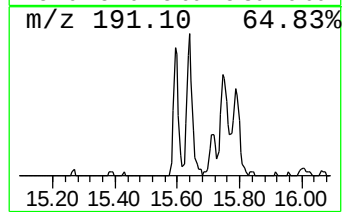
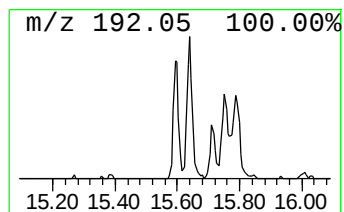
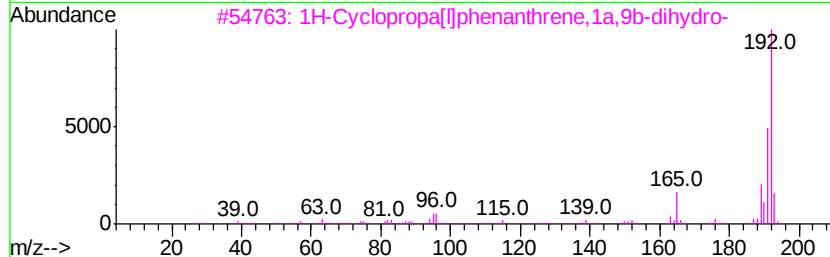
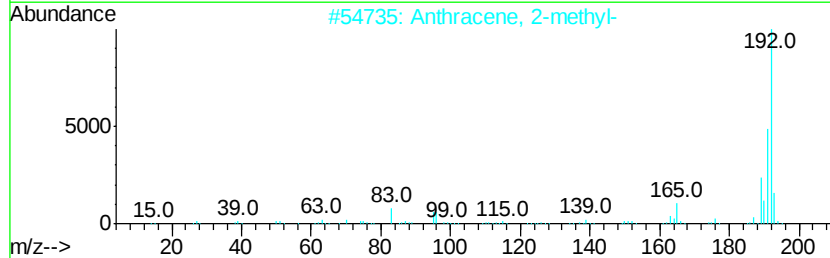
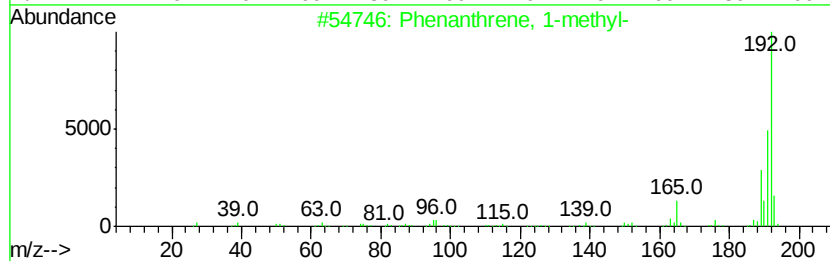
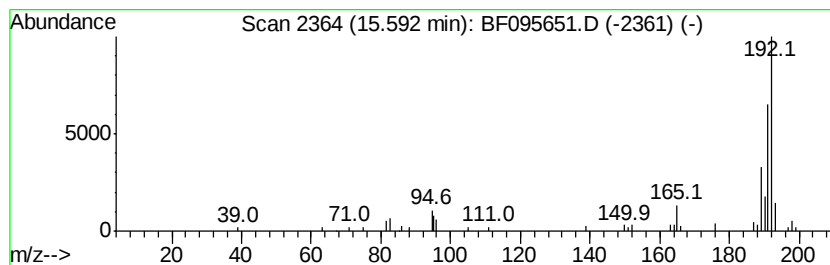
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Peak Number 5 Phenanthrene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.59	2.79 ng	63281	Phenanthrene-d10	14.80

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
3		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	94
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060217\
Data File : BF095651.D
Acq On : 3 Jun 2017 00:22
Operator : SJ/MA
Sample : I3441-03 5X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
OR-RO-5004-060117

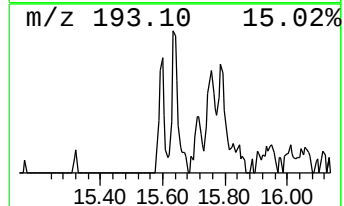
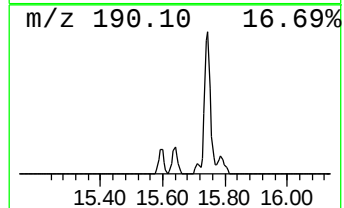
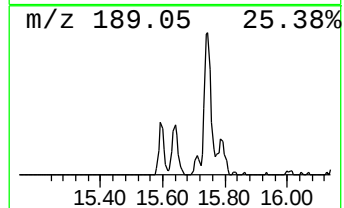
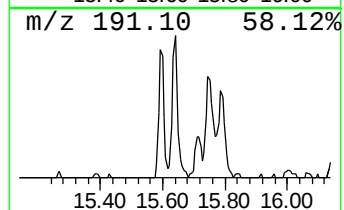
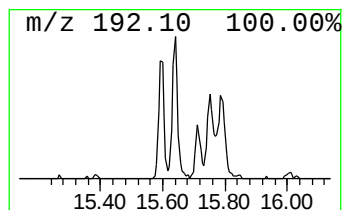
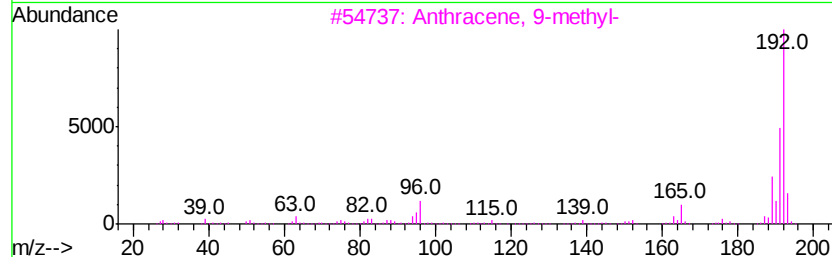
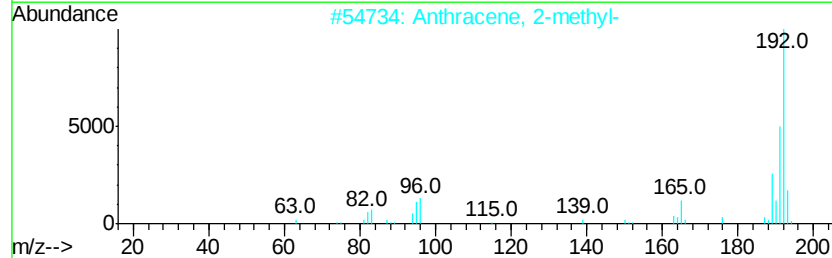
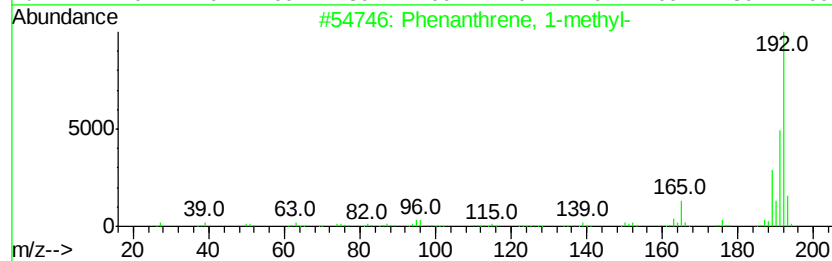
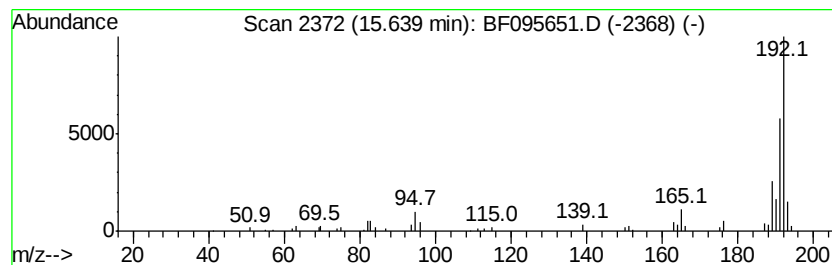
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Anthracene, 9-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.64	3.70 ng	83850	Phenanthrene-d10	14.80

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3		Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	92
5		Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	90



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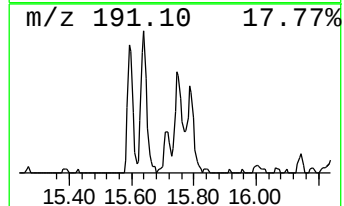
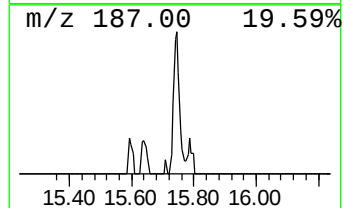
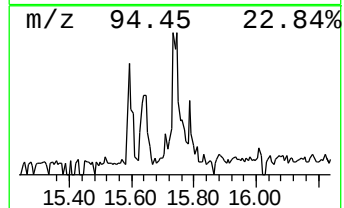
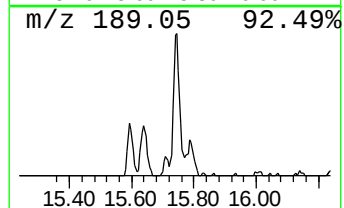
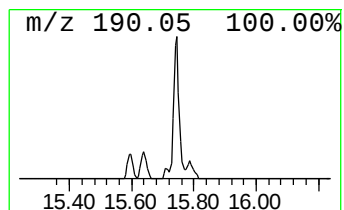
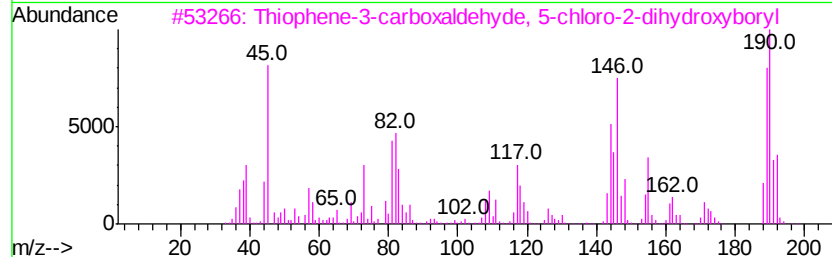
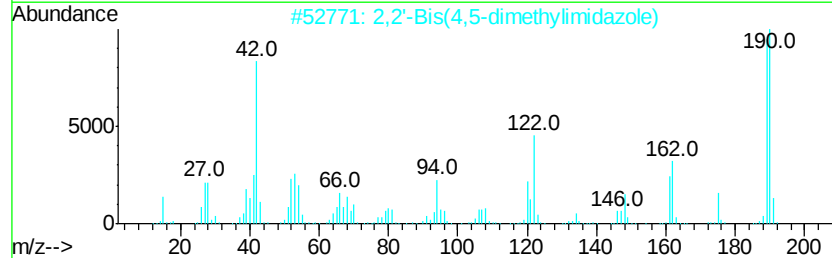
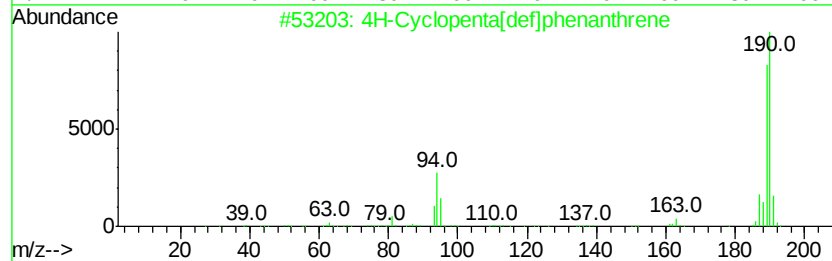
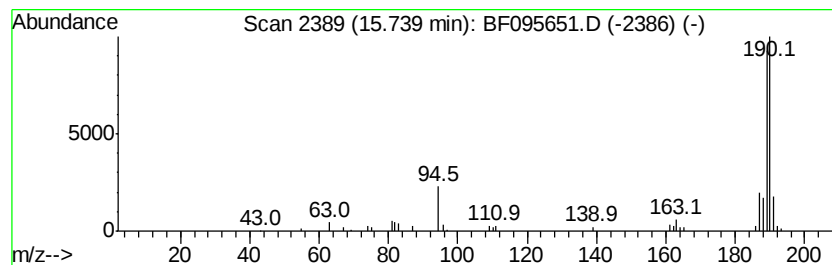
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.74	4.89 ng	110841	Phenanthrene-d10	14.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
2		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45
3		Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	40
4		6H-Cyclobuta[flk]phenanthrene	190	C15H10	083469-43-6	33
5		6,7-Methylenedioxy-4(3H)-quinazo...	190	C9H6N2O3	028310-11-4	33



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060217\
Data File : BF095651.D
Acq On : 3 Jun 2017 00:22
Operator : SJ/MA
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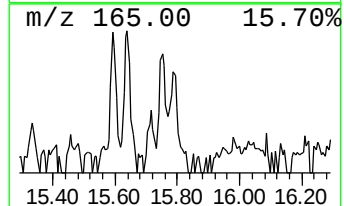
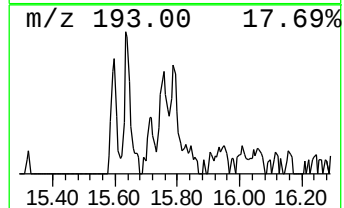
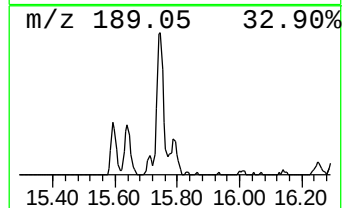
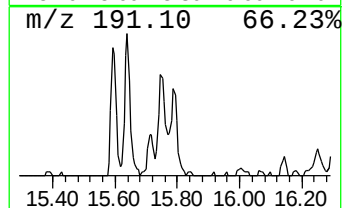
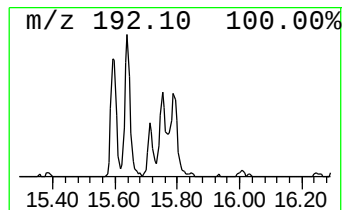
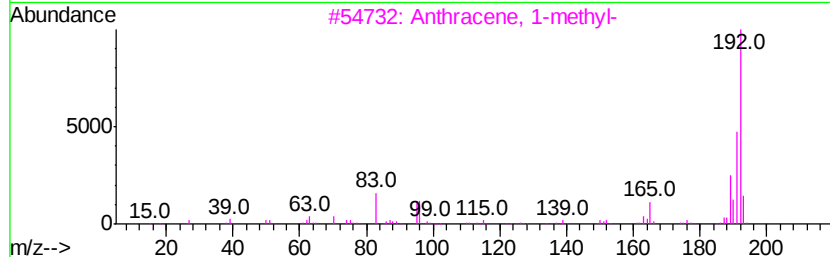
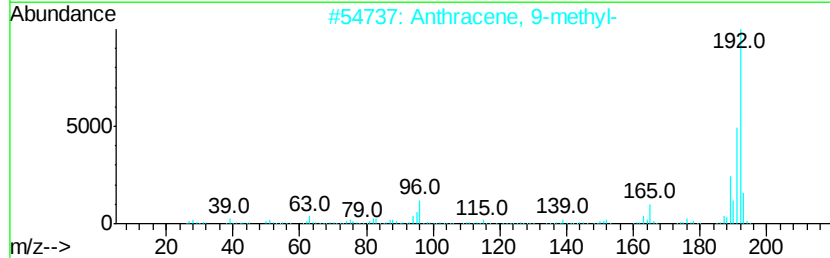
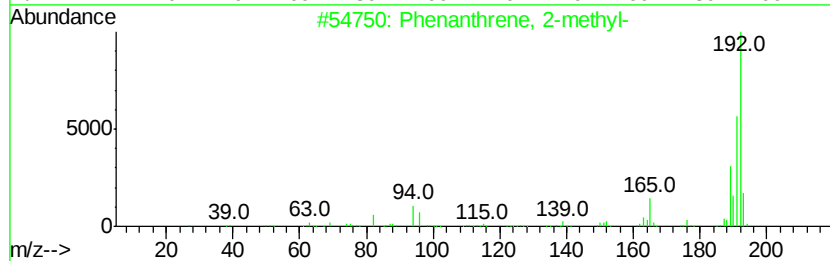
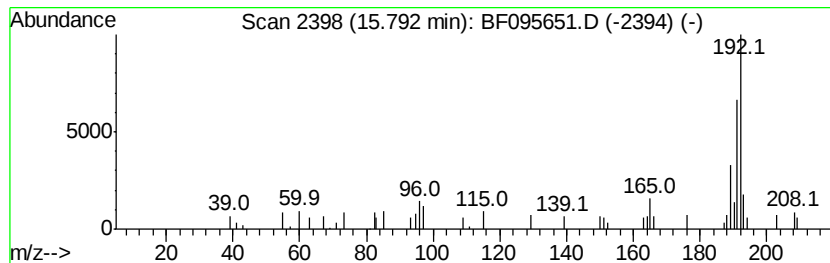
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenanthrene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.79	2.72 ng	61596	Phenanthrene-d10	14.80

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	81
2		Anthracene, 9-methyl-	192	C15H12	000779-02-2	76
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	62
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	62
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	58



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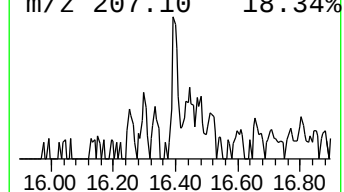
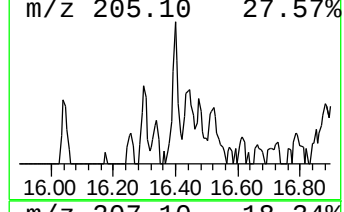
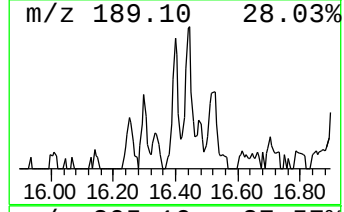
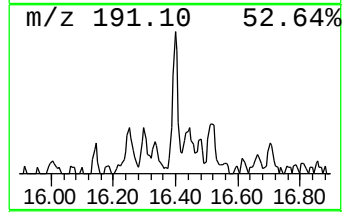
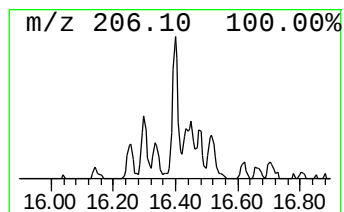
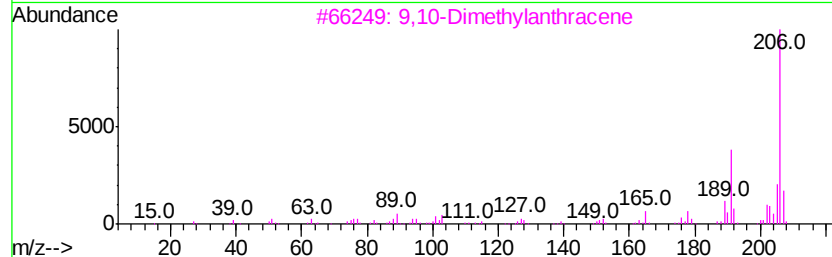
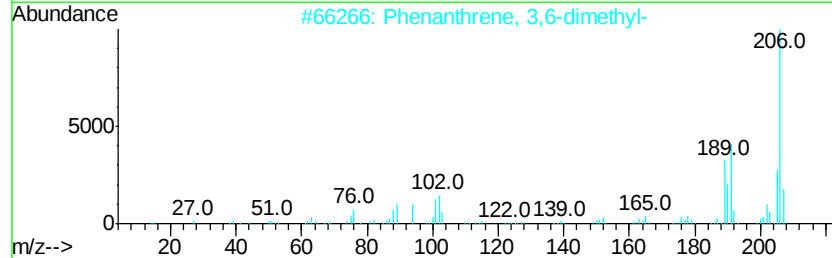
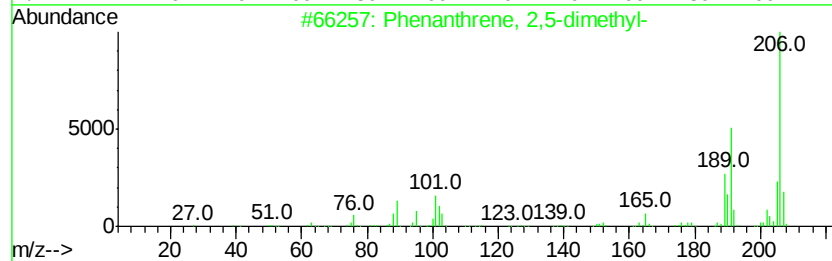
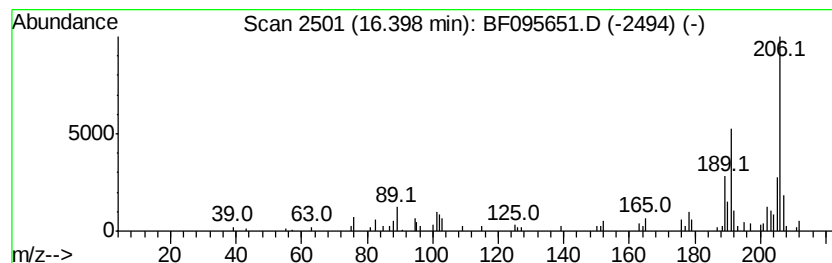
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenanthrene, 2,5-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.40	3.66 ng	82942	Phenanthrene-d10	14.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	95
2		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	94
3		9,10-Dimethylanthracene	206	C16H14	000781-43-1	94
4		Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	92
5		1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	92



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060217\
Data File : BF095651.D
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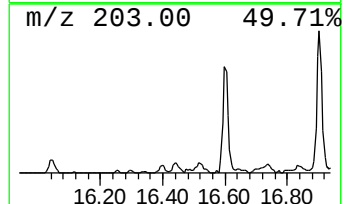
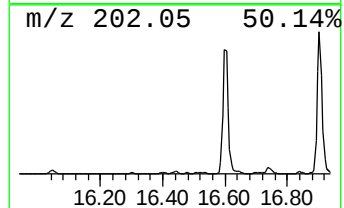
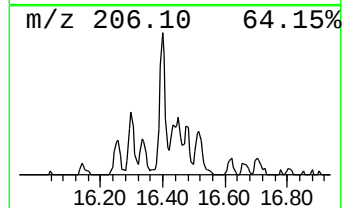
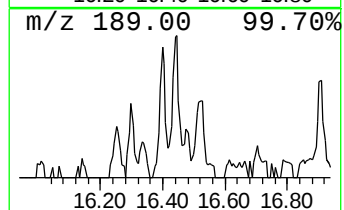
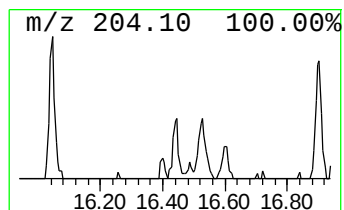
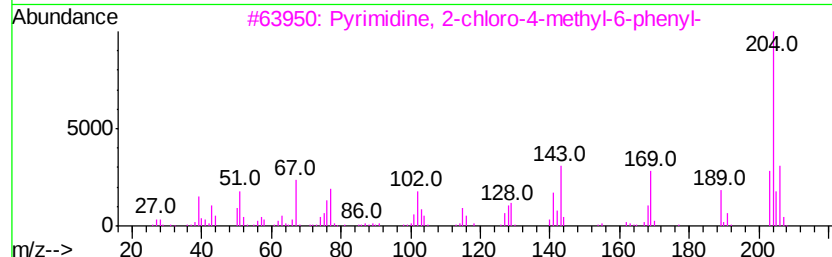
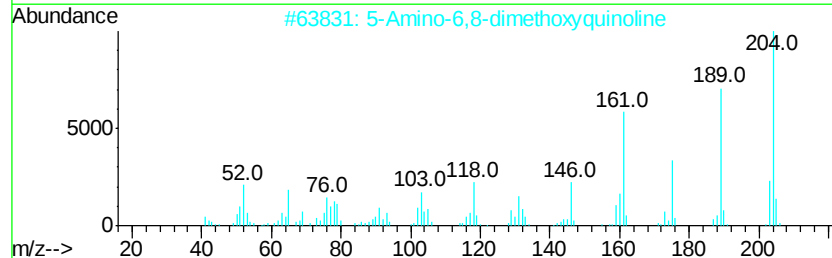
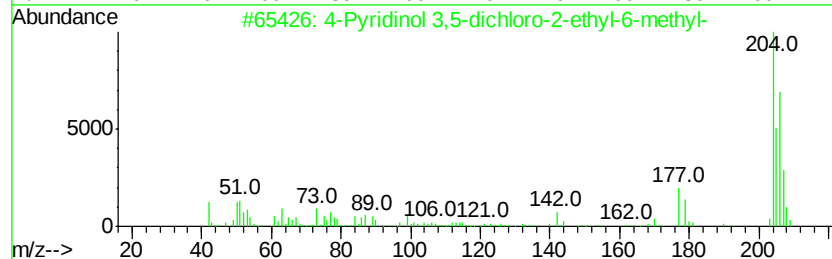
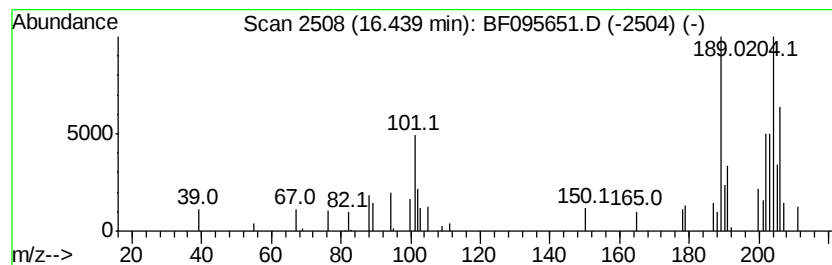
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown16.44 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.44	2.10 ng	47545	Phenanthrene-d10	14.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Pyridinol 3,5-dichloro-2-ethyl...	205	C8H9Cl2NO	1000253-55-0	38
2			5-Amino-6,8-dimethoxyquinoline	204	C11H12N2O2	1000213-95-2	38
3			Pyrimidine, 2-chloro-4-methyl-6-...	204	C11H9ClN2	032785-40-3	35
4			Naphthalene, 1,2,4a,5,8,8a-hexah...	204	C15H24	005951-61-1	35
5			5-Amino-7,8-dimethoxyisoquinoline	204	C11H12N2O2	1000213-88-5	35



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060217\
Data File : BF095651.D
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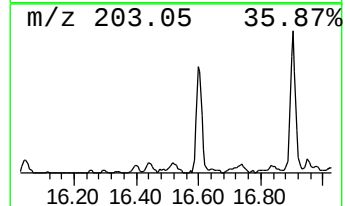
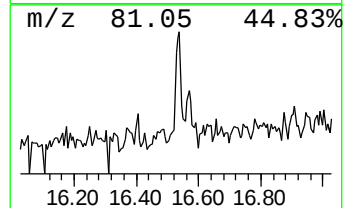
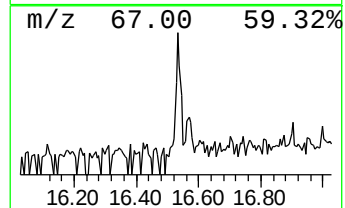
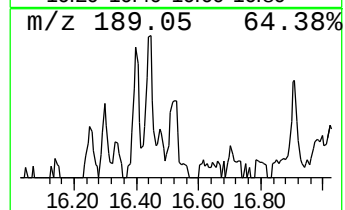
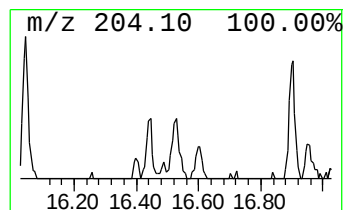
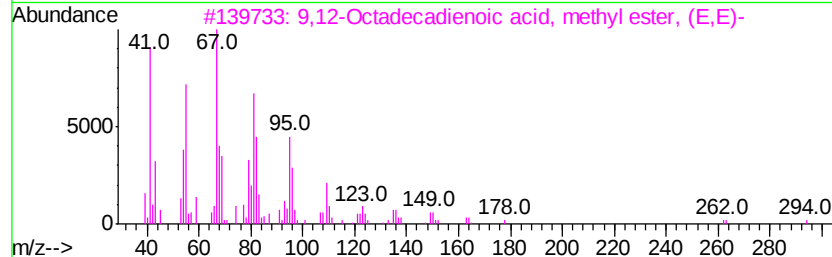
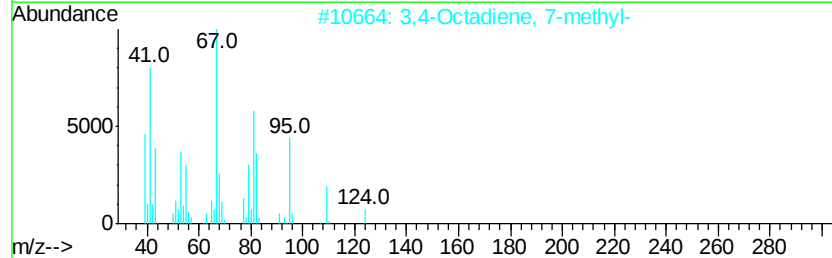
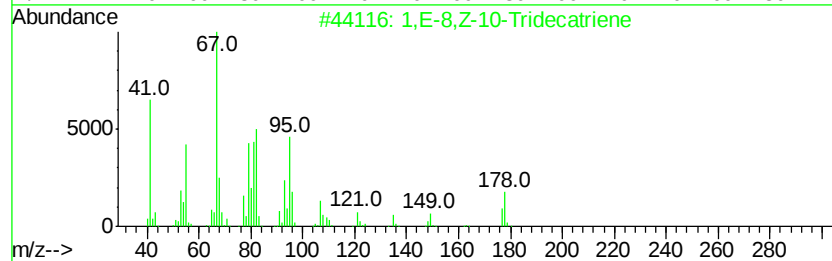
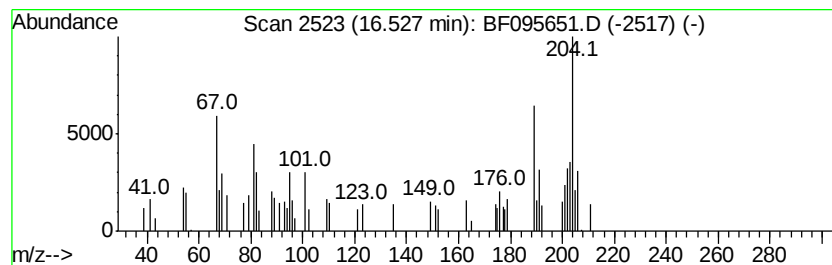
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown16.53 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.53	2.85 ng	64486	Phenanthrene-d10	14.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,E-8,Z-10-Tridecatriene	178	C13H22	080625-30-5	46
2			3,4-Octadiene, 7-methyl-	124	C9H16	037050-05-8	38
3			9,12-Octadecadienoic acid, methv...	294	C19H34O2	002566-97-4	38
4			Bicyclo[5.1.0]octane	110	C8H14	000286-43-1	30
5			Cyclopentane, (1-methylethenyl)-	110	C8H14	055661-02-4	30



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060217\
Data File : BF095651.D
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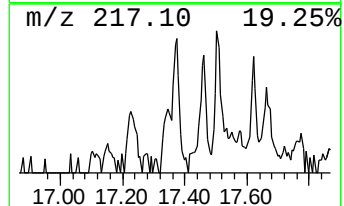
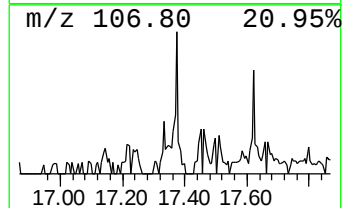
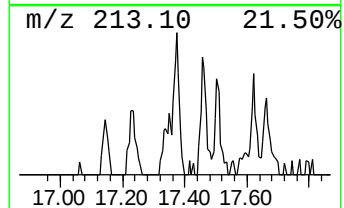
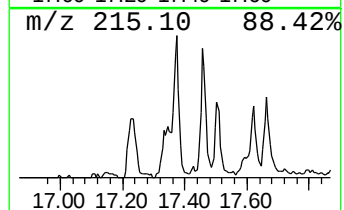
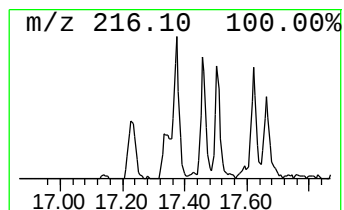
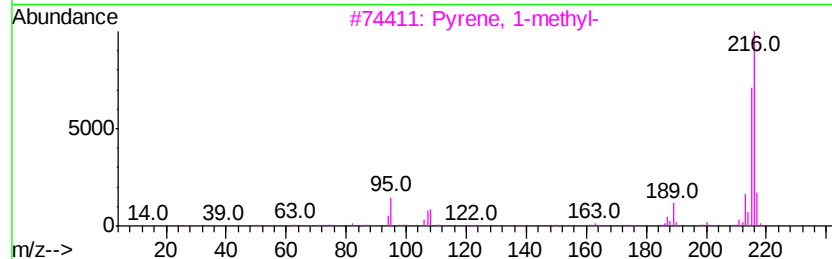
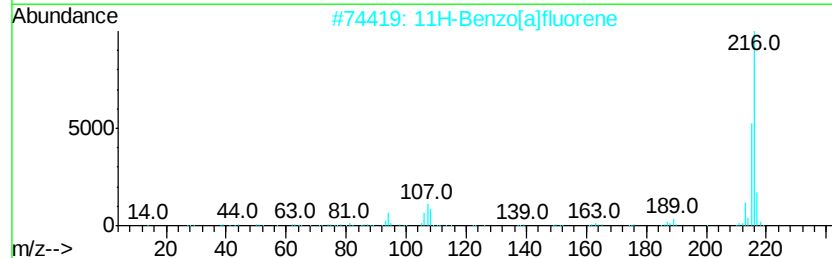
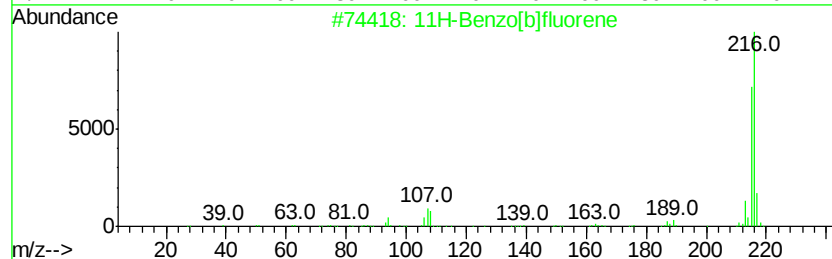
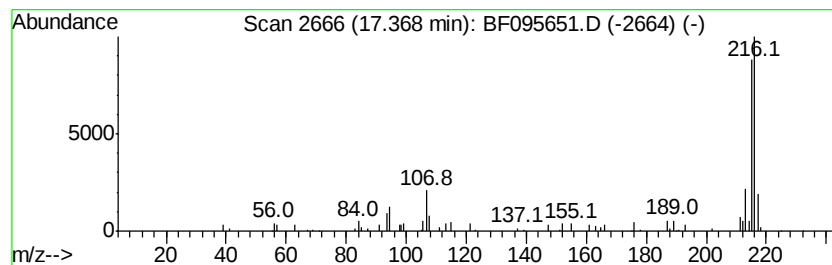
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 11H-Benzo[b]fluorene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.37	2.31 ng	77499	Chrysene-d12	18.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
2		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
3		Pvrene, 1-methyl-	216	C17H12	002381-21-7	87
4		7H-Benzo[c]fluorene	216	C17H12	000205-12-9	81
5		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060217\
Data File : BF095651.D
Acq On : 3 Jun 2017 00:22
Operator : SJ/MA
Sample : I3441-03 5X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OR-RO-5004-060117

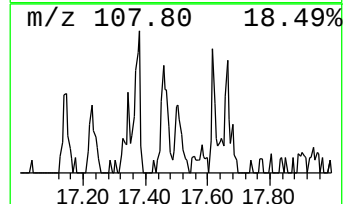
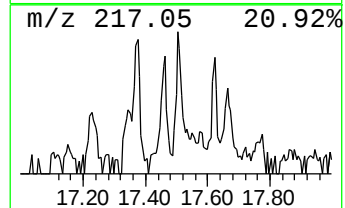
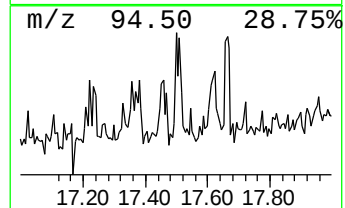
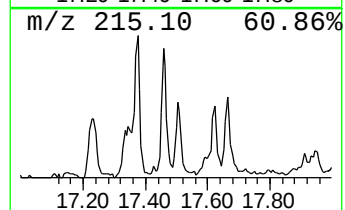
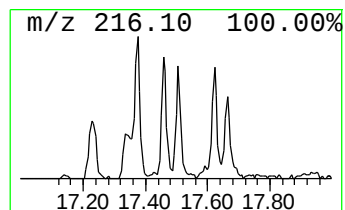
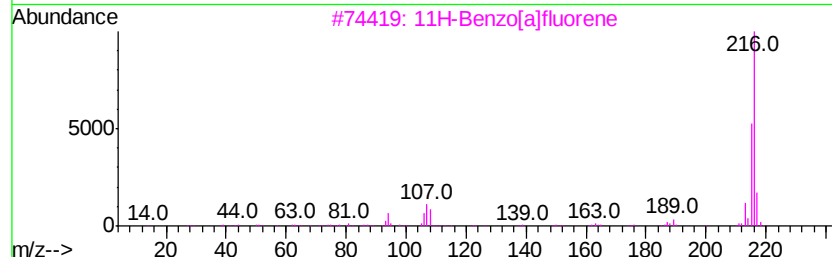
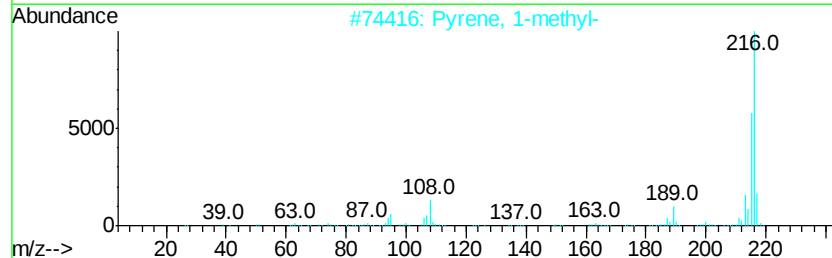
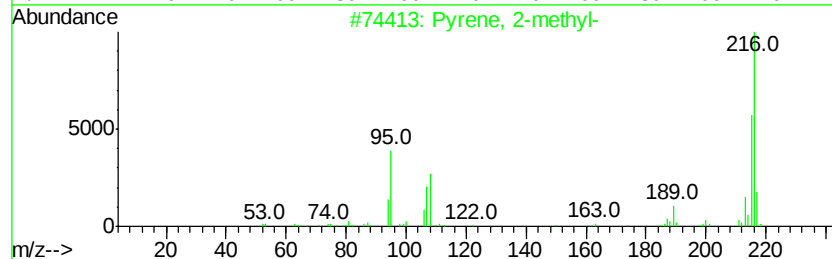
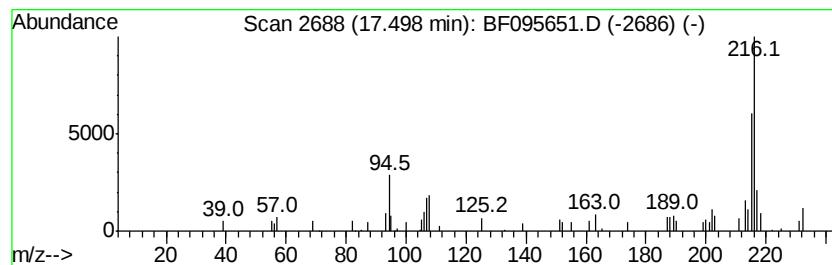
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Pyrene, 2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.50	2.03 ng	68021	Chrysene-d12	18.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	93
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
3		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	80
4		4,6-Dimethoxy-1-naphthaldehyde	216	C13H12O3	065565-33-5	59
5		1,3-Dihydro-5,6-dimethyl-1-(2-pr...	216	C12H12N2S	080276-22-8	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060217\
Data File : BF095651.D
Acq On : 3 Jun 2017 00:22
Operator : SJ/MA
Sample : I3441-03 5X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OR-RO-5004-060117

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	16.1	ng	357371	1	7.54	444997	20.0
1,4-Methanonaphth...	10.89	2.3	ng	65095	2	9.57	569257	20.0
Phenanthrene, 1-m...	15.59	2.8	ng	63281	4	14.80	453287	20.0
Anthracene, 9-met...	15.64	3.7	ng	83850	4	14.80	453287	20.0
4H-Cyclopenta[def...	15.74	4.9	ng	110841	4	14.80	453287	20.0
Phenanthrene, 2-m...	15.79	2.7	ng	61596	4	14.80	453287	20.0
Phenanthrene, 2,5...	16.40	3.7	ng	82942	4	14.80	453287	20.0
unknown16.44	16.44	2.1	ng	47545	4	14.80	453287	20.0
unknown16.53	16.53	2.9	ng	64486	4	14.80	453287	20.0
11H-Benzo[b]fluorene	17.37	2.3	ng	77499	5	18.50	670310	20.0
Pyrene, 2-methyl-	17.50	2.0	ng	68021	5	18.50	670310	20.0