

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060717\
 Data File : BF095771.D
 Acq On : 8 Jun 2017 5:23
 Operator : SJ/MA
 Sample : I3469-12
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-3(0-2)20170603

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-BF060517.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	4.592	507	511	536	rBV2	53974	159012	8.06%	0.669%
2	5.275	623	627	658	rBV	1039276	1782970	90.36%	7.505%
3	6.945	906	911	923	rBV	1112152	1710642	86.70%	7.200%
4	7.039	923	927	943	rVB	1343646	1960562	99.37%	8.252%
5	7.386	981	986	1002	rBV	309635	407172	20.64%	1.714%
6	7.622	1021	1026	1040	rBV	1122094	1436161	72.79%	6.045%
7	8.304	1136	1142	1160	rBV	730961	1046138	53.02%	4.403%
8	9.416	1326	1331	1345	rBV	425269	532240	26.97%	2.240%
9	11.198	1628	1634	1652	rVB	1577534	1973089	100.00%	8.305%
10	11.886	1746	1751	1759	rBV	202411	242640	12.30%	1.021%
11	12.239	1805	1811	1816	rBV2	434414	521461	26.43%	2.195%
12	12.286	1816	1819	1825	rVB2	108976	140961	7.14%	0.593%
13	12.586	1863	1870	1873	rBV2	16338	23336	1.18%	0.098%
14	13.098	1950	1957	1960	rBV	41123	50166	2.54%	0.211%
15	13.133	1960	1963	1974	rVB2	73232	93322	4.73%	0.393%
16	13.268	1983	1986	1991	rVB3	17141	22081	1.12%	0.093%
17	13.457	2012	2018	2022	rVV3	15483	28866	1.46%	0.121%
18	13.533	2026	2031	2039	rVV	661871	870962	44.14%	3.666%
19	13.868	2084	2088	2090	rBV2	40587	49612	2.51%	0.209%
20	14.033	2110	2116	2119	rBV2	14959	22418	1.14%	0.094%
21	14.421	2178	2182	2186	rBV5	11372	23709	1.20%	0.100%
22	14.462	2186	2189	2195	rVB	30079	40720	2.06%	0.171%
23	14.598	2207	2212	2214	rBV	26244	30924	1.57%	0.130%
24	14.633	2214	2218	2221	rVV	356706	451982	22.91%	1.902%
25	14.668	2221	2224	2232	rVB	594726	717249	36.35%	3.019%
26	14.756	2232	2239	2246	rVB	198745	255157	12.93%	1.074%
27	14.856	2250	2256	2260	rBV4	9953	20107	1.02%	0.085%
28	15.280	2323	2328	2331	rBV2	13698	23267	1.18%	0.098%
29	15.433	2349	2354	2358	rBV	73695	87019	4.41%	0.366%
30	15.474	2358	2361	2366	rVV	86650	105612	5.35%	0.445%
31	15.551	2370	2374	2376	rVV2	51394	56384	2.86%	0.237%
32	15.586	2376	2380	2385	rVV3	125858	186768	9.47%	0.786%
33	15.627	2385	2387	2389	rVV2	51602	65902	3.34%	0.277%
34	15.651	2389	2391	2400	rVB	95175	105102	5.33%	0.442%

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35	15.886	2427	2431	2437	rVB	65360	71239	3.61%	0.300%
36	16.098	2463	2467	2471	rBV4	13944	20846	1.06%	0.088%
37	16.151	2471	2476	2479	rBV3	14092	21953	1.11%	0.092%
38	16.245	2488	2492	2496	rBV	47858	59507	3.02%	0.250%
39	16.286	2496	2499	2503	rVV5	27108	36496	1.85%	0.154%
40	16.368	2509	2513	2520	rVB4	31202	46551	2.36%	0.196%
41	16.450	2520	2527	2532	rBV	823033	984532	49.90%	4.144%
42	16.498	2532	2535	2541	rVB4	19584	24733	1.25%	0.104%
43	16.550	2541	2544	2548	rBV4	27550	38362	1.94%	0.161%
44	16.645	2556	2560	2569	rBV2	28786	56475	2.86%	0.238%
45	16.750	2572	2578	2583	rVV	921181	1041409	52.78%	4.383%
46	16.798	2583	2586	2589	rVV	47146	53354	2.70%	0.225%
47	16.839	2589	2593	2596	rVV3	27163	47609	2.41%	0.200%
48	16.868	2596	2598	2606	rVB5	59731	72502	3.67%	0.305%
49	16.945	2607	2611	2614	rBV	59729	64835	3.29%	0.273%
50	17.003	2614	2621	2625	rVV	1106992	1363508	69.11%	5.739%
51	17.080	2631	2634	2638	rVV2	40427	56414	2.86%	0.237%
52	17.127	2639	2642	2646	rVB3	19343	22509	1.14%	0.095%
53	17.192	2646	2653	2655	rBV7	46178	88064	4.46%	0.371%
54	17.221	2655	2658	2662	rVB	147239	169133	8.57%	0.712%
55	17.274	2663	2667	2669	rBV4	19711	25142	1.27%	0.106%
56	17.309	2669	2673	2677	rVV	149996	175889	8.91%	0.740%
57	17.350	2677	2680	2686	rVV5	77677	157226	7.97%	0.662%
58	17.415	2688	2691	2695	rVV3	39470	59801	3.03%	0.252%
59	17.468	2697	2700	2703	rVV	42546	49747	2.52%	0.209%
60	17.509	2703	2707	2713	rVV	36706	59603	3.02%	0.251%
61	17.568	2713	2717	2720	rVV5	18267	28036	1.42%	0.118%
62	17.797	2753	2756	2760	rVB3	54425	62493	3.17%	0.263%
63	17.862	2763	2767	2772	rBV4	51724	71143	3.61%	0.299%
64	18.021	2791	2794	2797	rBV	56320	66131	3.35%	0.278%
65	18.056	2797	2800	2802	rVV	60524	60395	3.06%	0.254%
66	18.086	2802	2805	2809	rVV	43981	54671	2.77%	0.230%
67	18.121	2809	2811	2815	rVB2	50799	58546	2.97%	0.246%
68	18.262	2829	2835	2843	rBV5	96172	237058	12.01%	0.998%
69	18.344	2843	2849	2852	rVV2	519180	834990	42.32%	3.515%
70	18.380	2852	2855	2861	rVB	370654	414800	21.02%	1.746%
71	18.474	2867	2871	2875	rVB	98611	106281	5.39%	0.447%

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72	18.627	2894	2897	2902	rBV2	46519	49385	2.50%	0.208%
73	18.680	2902	2906	2911	rVB3	65340	94420	4.79%	0.397%
74	18.856	2933	2936	2939	rVV	55607	58501	2.96%	0.246%
75	18.892	2940	2942	2947	rVB2	41756	48913	2.48%	0.206%
76	18.968	2952	2955	2958	rVB2	41931	46617	2.36%	0.196%
77	19.003	2958	2961	2963	rBV3	31442	42317	2.14%	0.178%
78	19.062	2969	2971	2975	rBV2	45643	49909	2.53%	0.210%
79	19.615	3061	3065	3068	rBV	254266	361775	18.34%	1.523%
80	19.727	3081	3084	3087	rVB	51743	46314	2.35%	0.195%
81	19.903	3111	3114	3118	rVB	139591	145644	7.38%	0.613%
82	19.962	3121	3124	3128	rVB	219031	225987	11.45%	0.951%
83	20.015	3130	3133	3136	rBV	226538	247658	12.55%	1.042%
84	21.197	3331	3334	3341	rVB	85980	136948	6.94%	0.576%
85	21.538	3389	3392	3398	rVB2	64497	98289	4.98%	0.414%

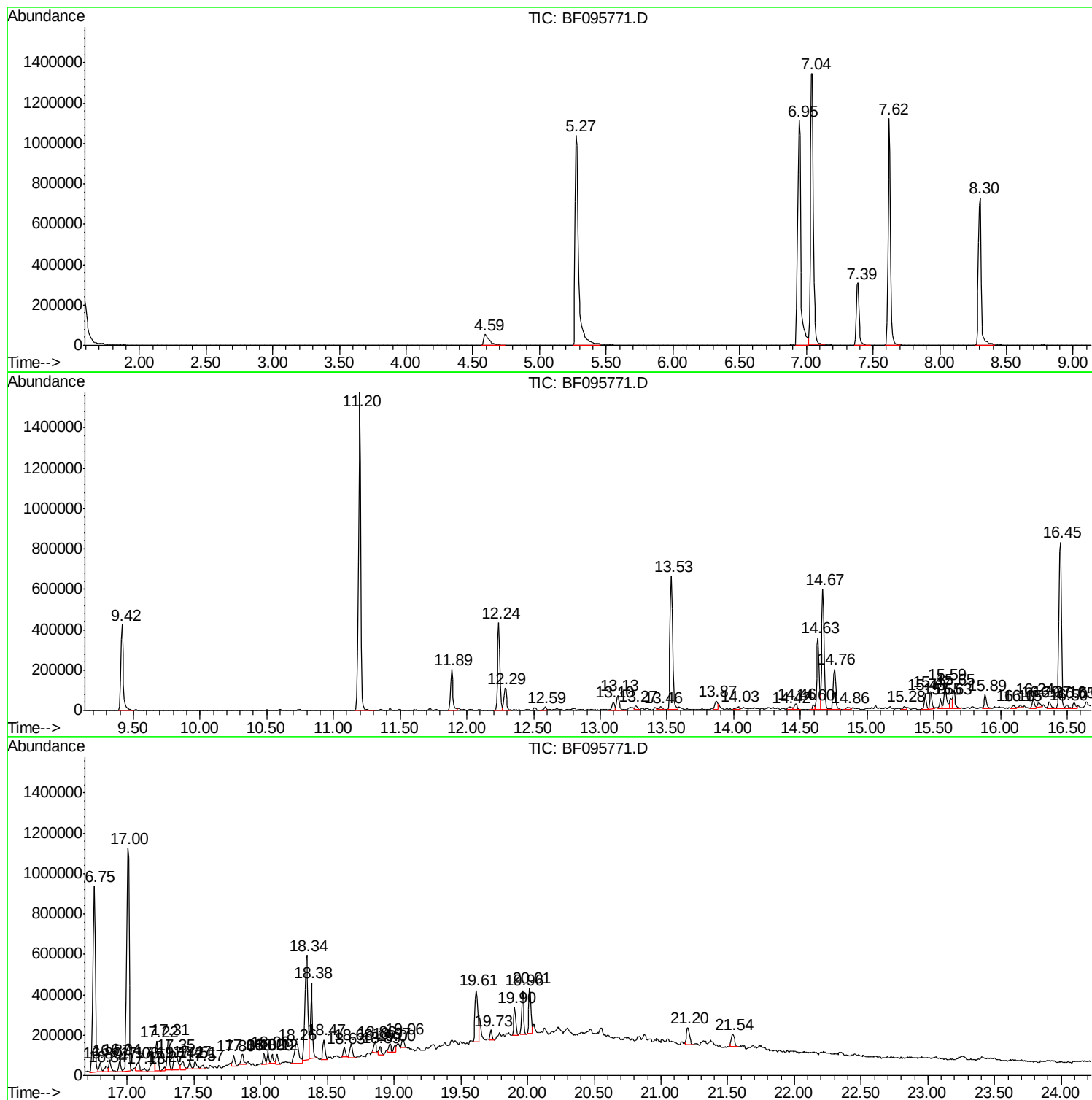
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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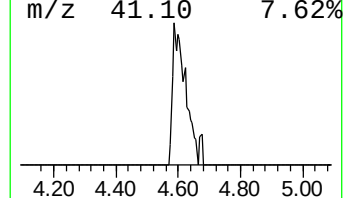
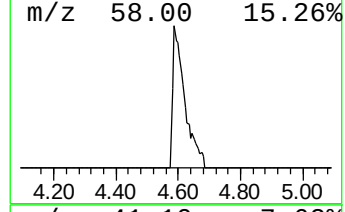
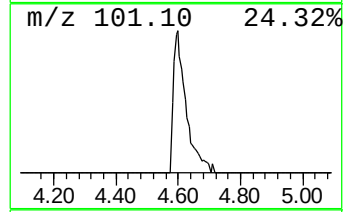
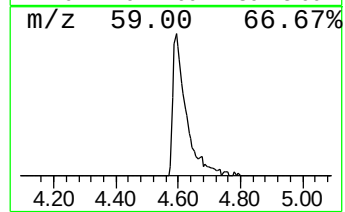
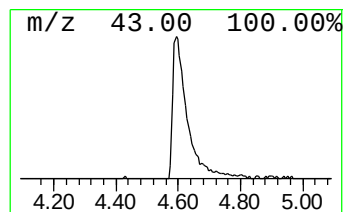
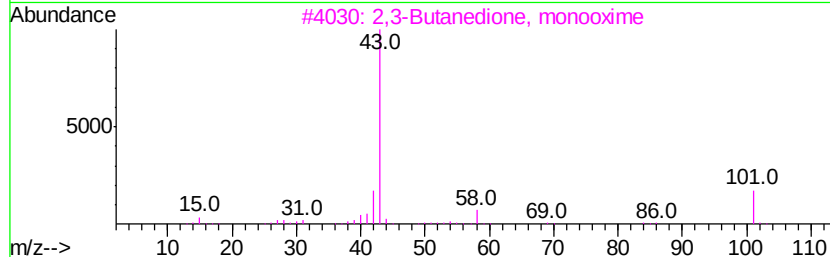
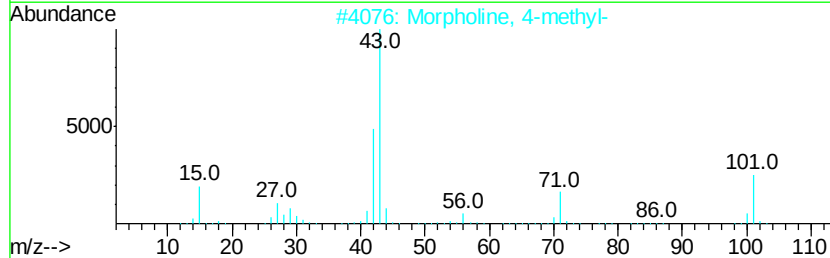
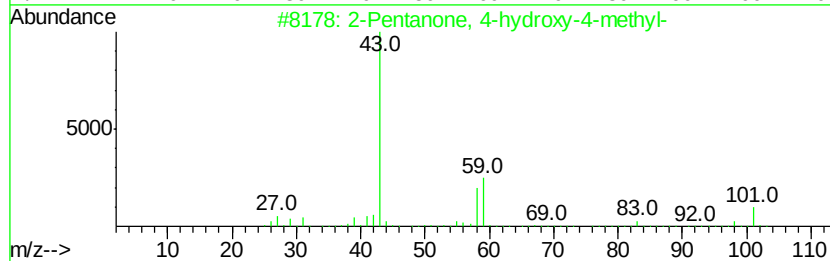
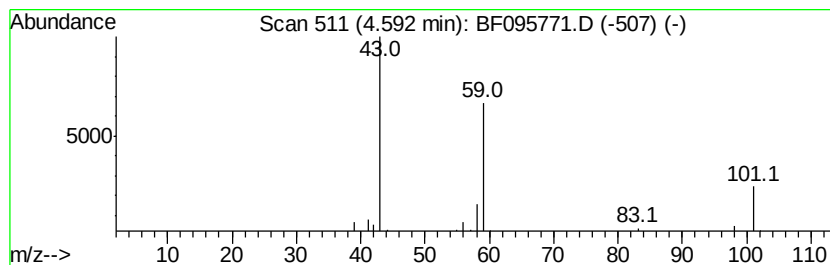
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.59	7.81 ng	159012	1,4-Dichlorobenzene-d4	7.39

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			5-Hexen-2-one	98	C6H10O	000109-49-9	9
5			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9



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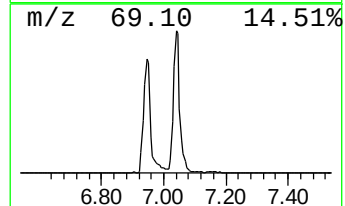
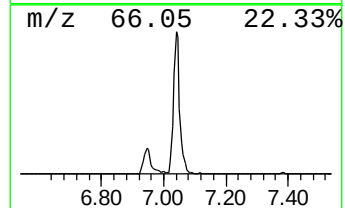
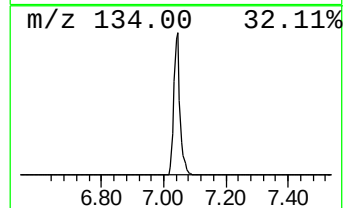
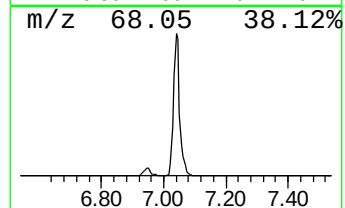
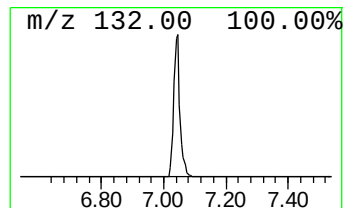
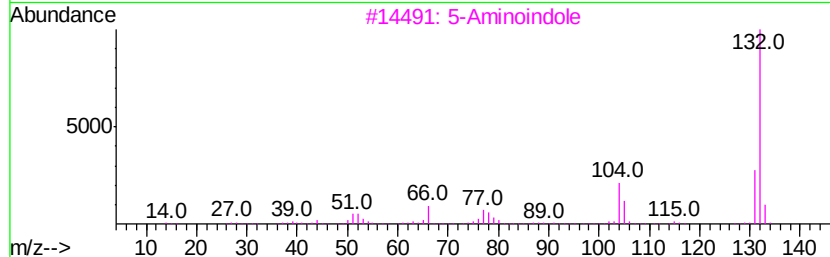
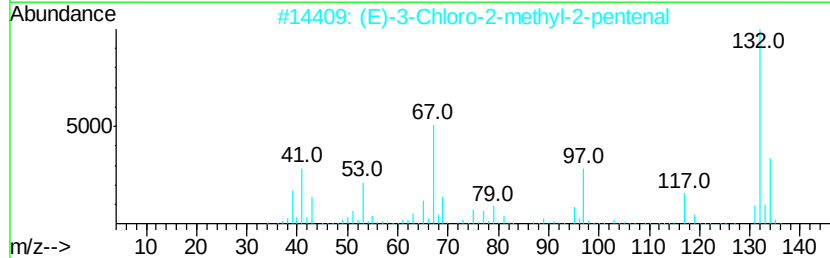
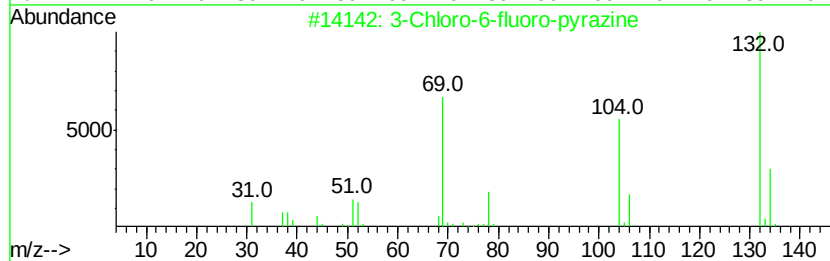
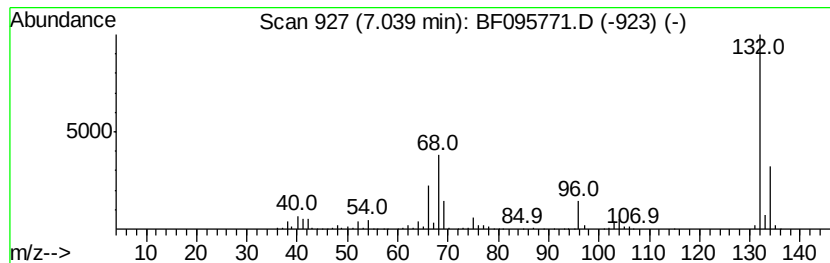
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown7.04 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.04	96.30 ng	1960560	1,4-Dichlorobenzene-d4	7.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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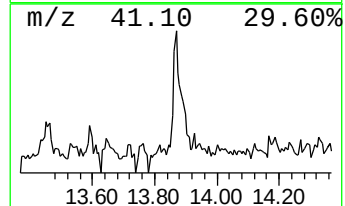
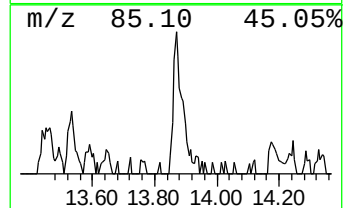
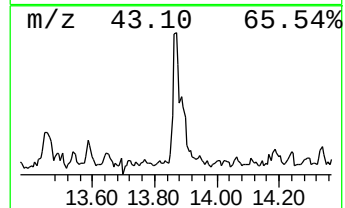
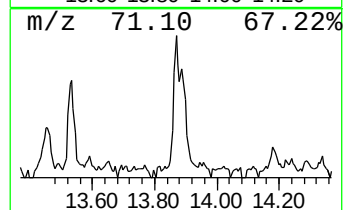
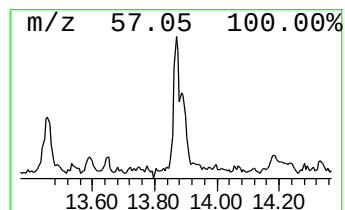
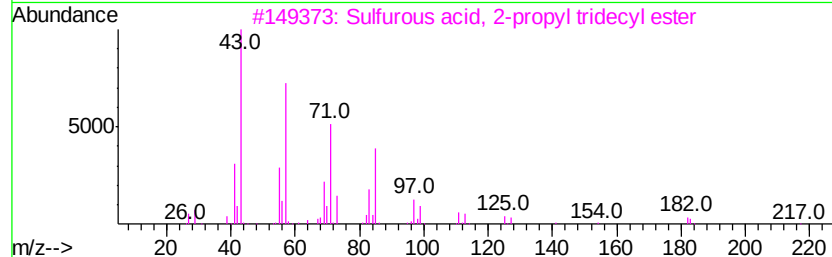
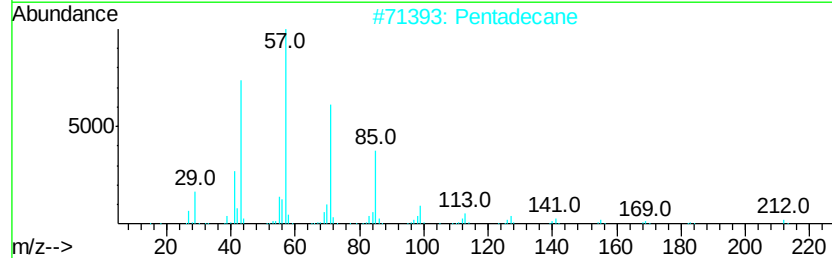
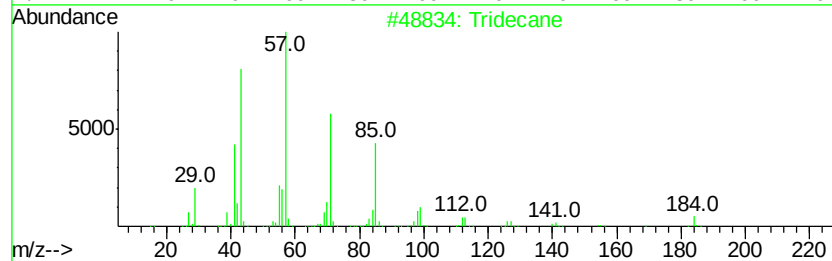
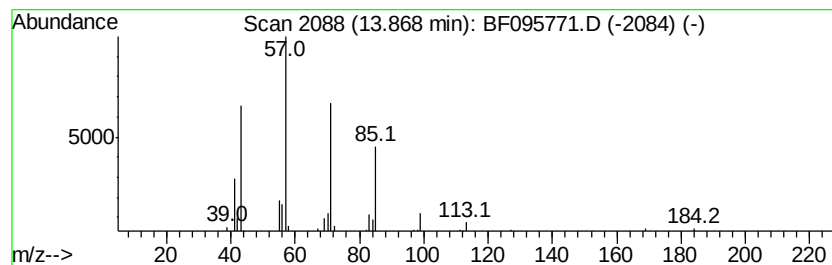
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Tridecane Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.87	2.20 ng	49612	Phenanthrene-d10		14.63
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	93
2	Pentadecane	212	C15H32	000629-62-9	86
3	Sulfurous acid, 2-propyl tridecyl...	306	C16H34O3S	1000309-12-4	86
4	Octadecane	254	C18H38	000593-45-3	86
5	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	86



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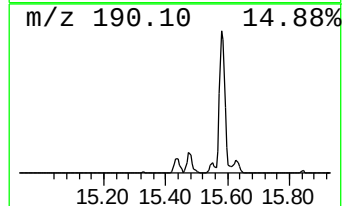
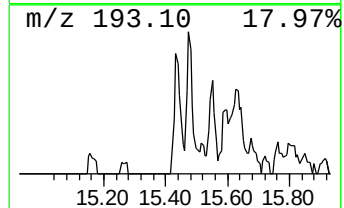
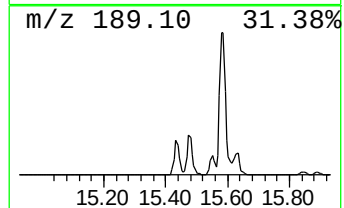
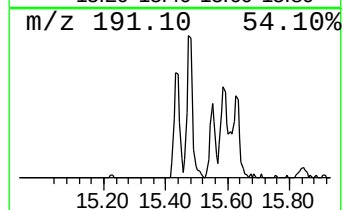
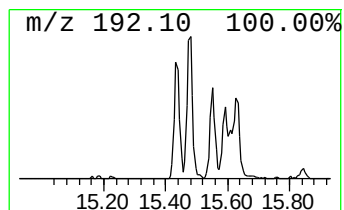
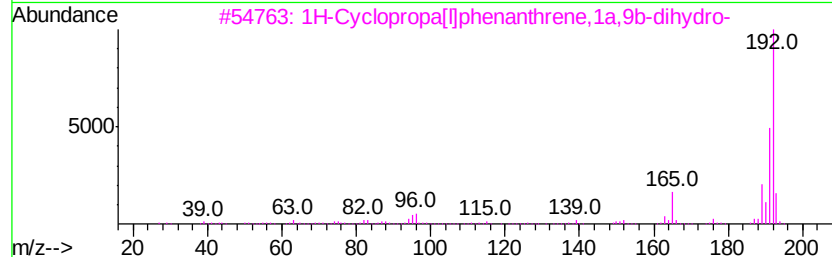
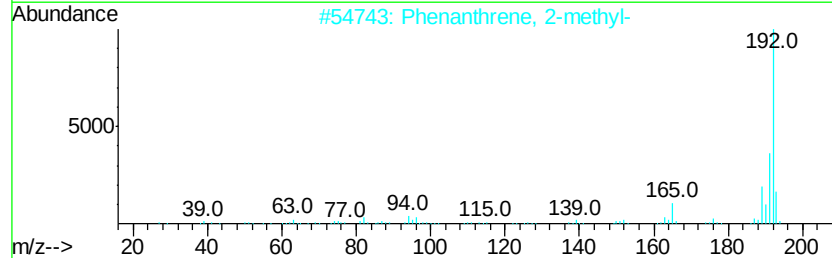
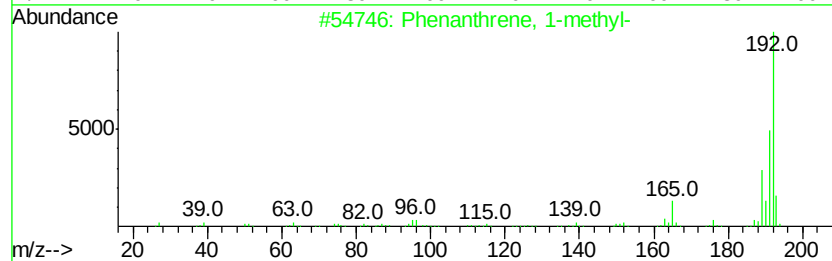
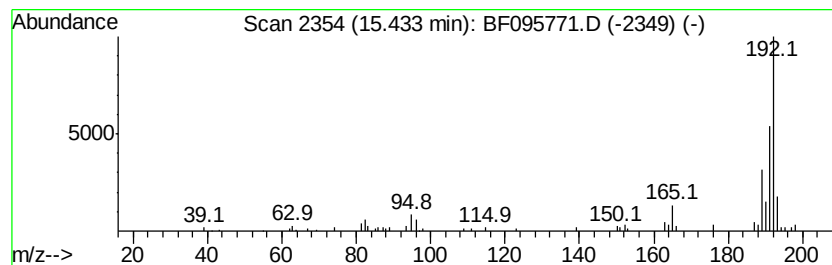
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Phenanthrene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.43	3.85 ng	87019	Phenanthrene-d10	14.63	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
5	1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	95



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060717\
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Acq On : 8 Jun 2017 5:23
Operator : SJ/MA
Sample : I3469-12
Misc :
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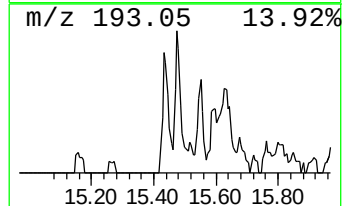
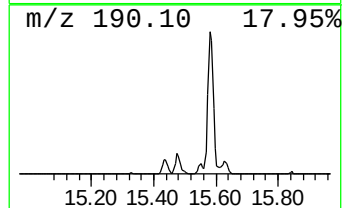
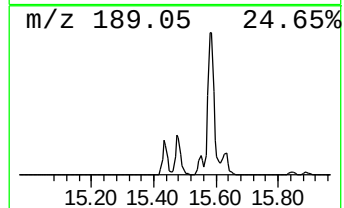
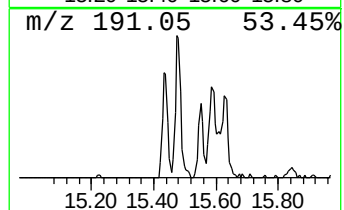
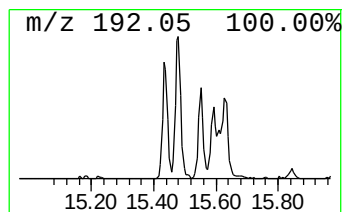
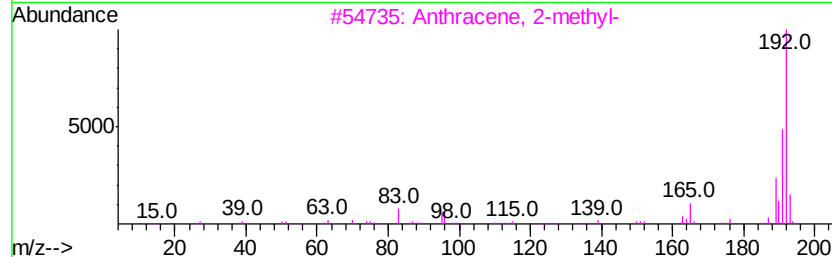
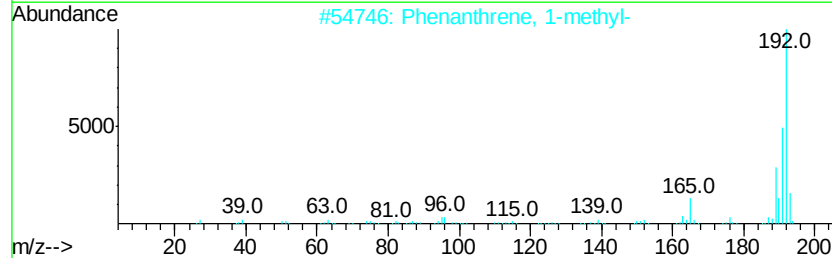
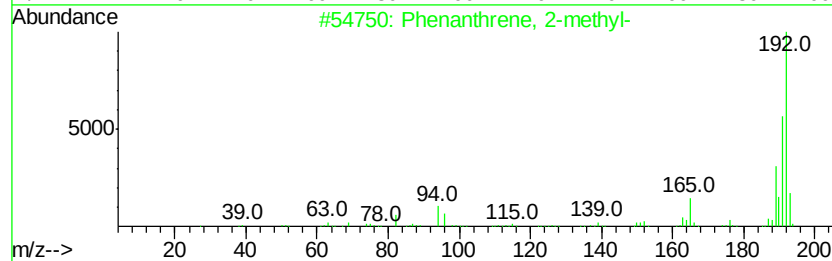
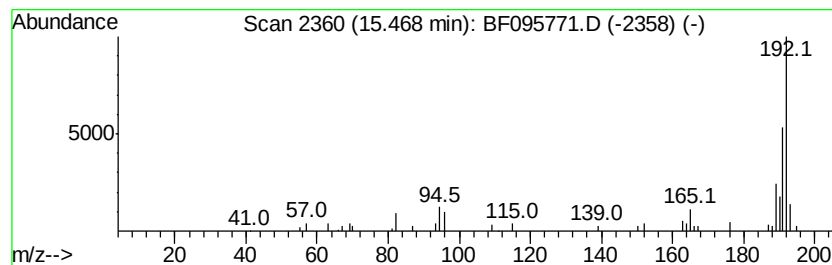
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Phenanthrene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.47	4.67 ng	105612	Phenanthrene-d10	14.63	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	95
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	95



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060717\
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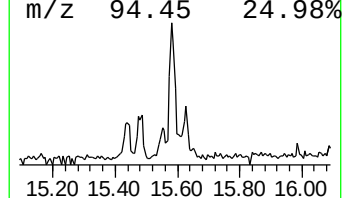
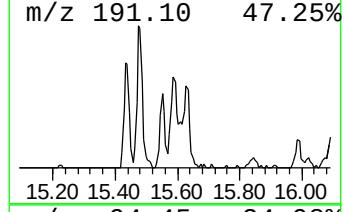
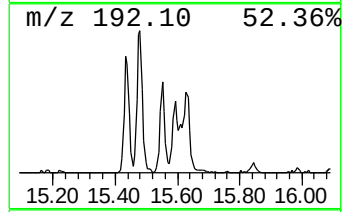
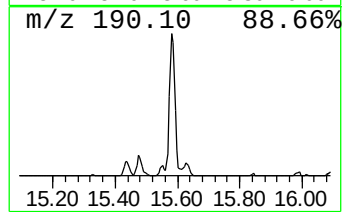
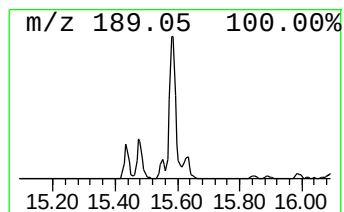
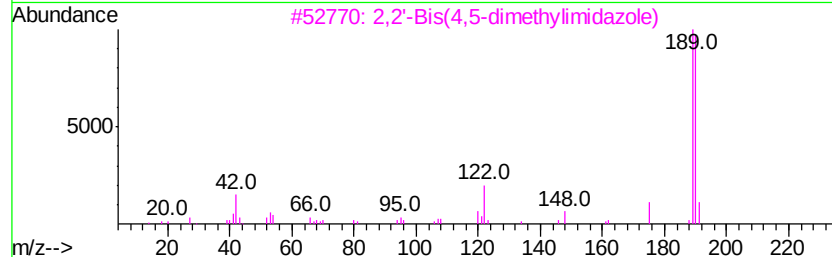
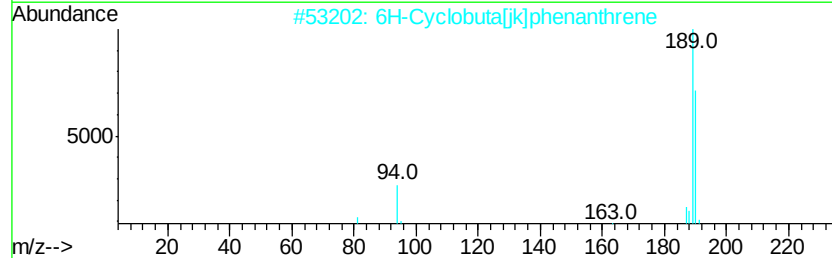
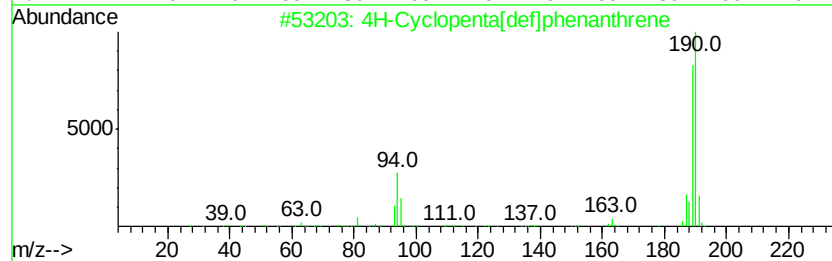
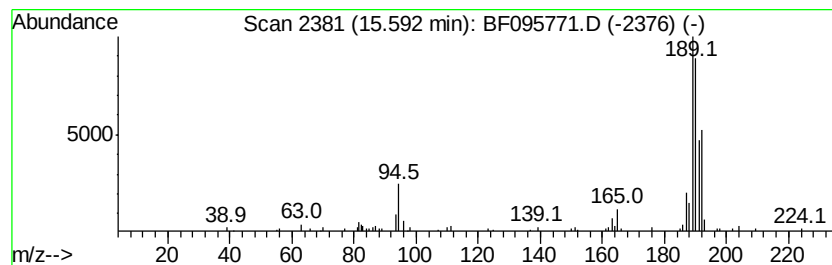
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD		R.T.
15.59	8.26 ng	186768	Phenanthrene-d10		14.63
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	91
2	6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	50
3	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40
4	Methyl diselenide	190	C2H6Se2	007101-31-7	38
5	2,6-Dichlorobenzaldoxime	189	C7H5Cl2NO	025185-95-9	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060717\
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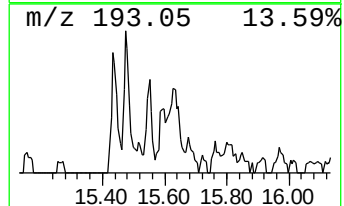
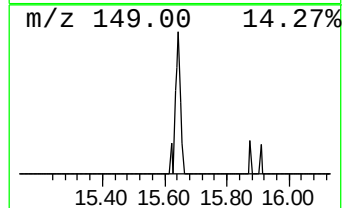
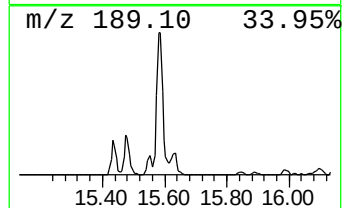
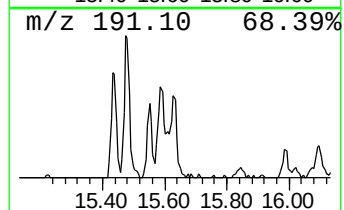
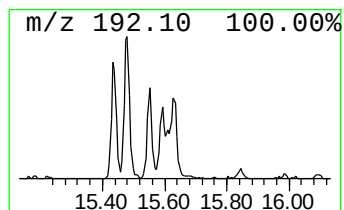
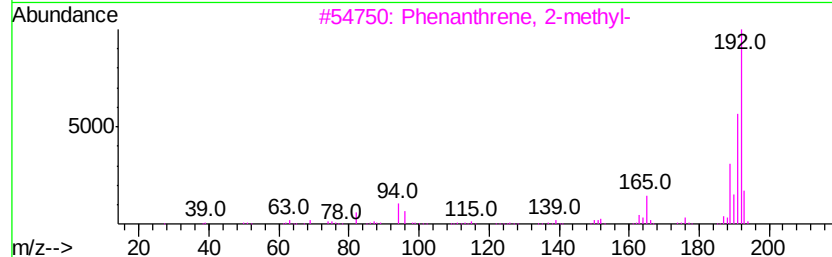
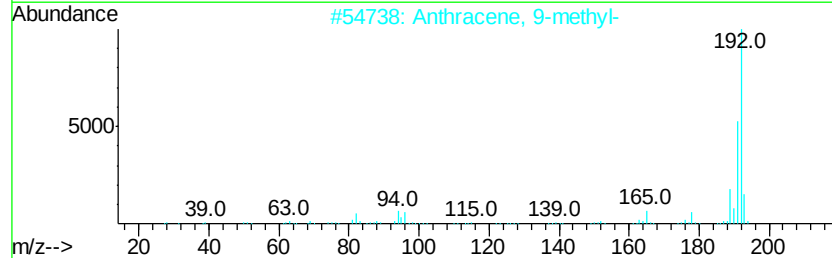
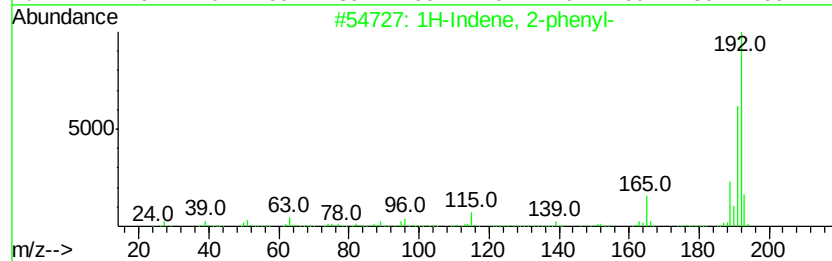
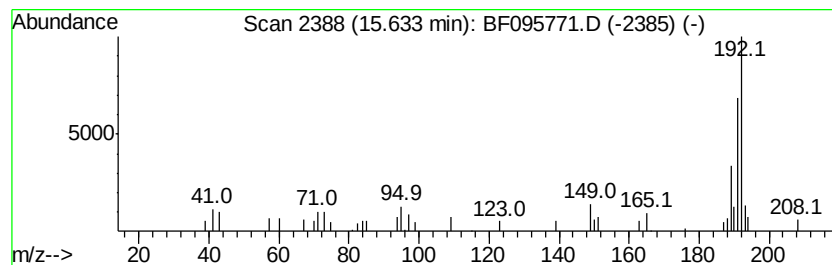
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1H-Indene, 2-phenyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.63	2.92 ng	65902	Phenanthrene-d10	14.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	70
2		Anthracene, 9-methyl-	192	C15H12	000779-02-2	68
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	64
4		Propyne, 1,3-diphenyl-	192	C15H12	004980-70-5	62
5		2-Hydroxy-4-methyl-6-(2-thienyl)...	192	C9H8N2OS	026963-45-1	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060717\
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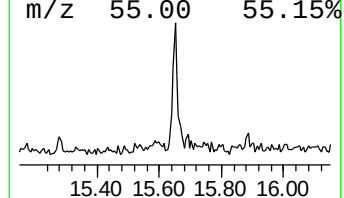
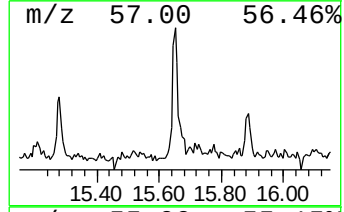
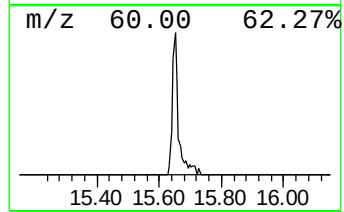
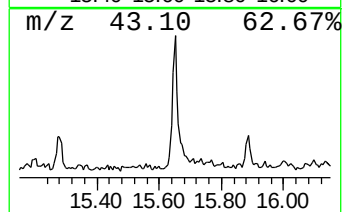
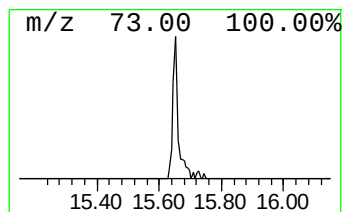
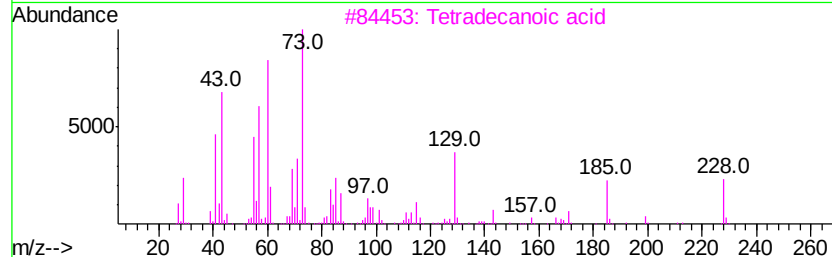
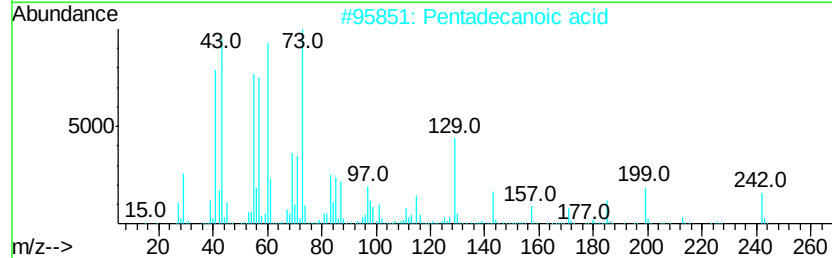
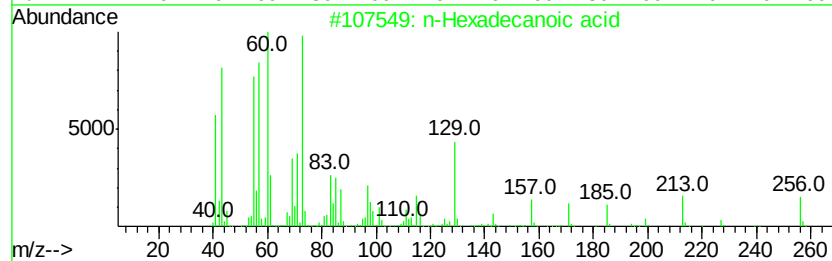
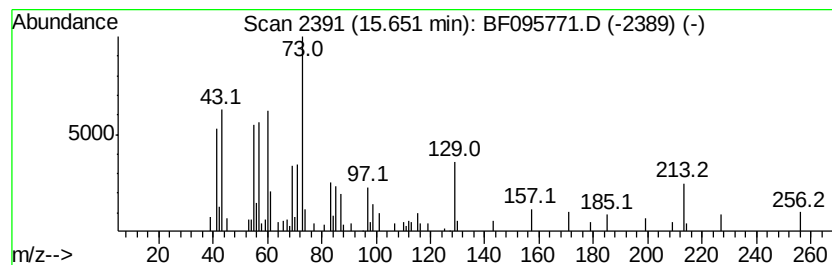
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.65	4.65 ng	105102	Phenanthrene-d10	14.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	64
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	53
4	n-Decanoic acid	172	C10H20O2	000334-48-5	43
5	Dodecanoic acid	200	C12H24O2	000143-07-7	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060717\
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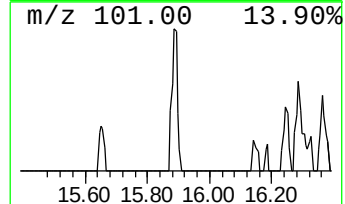
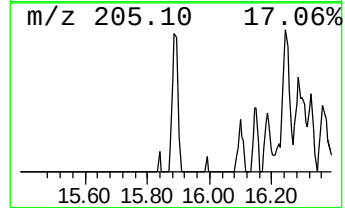
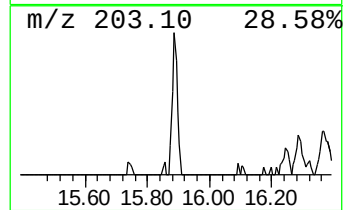
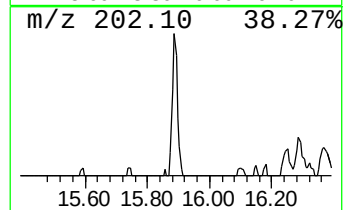
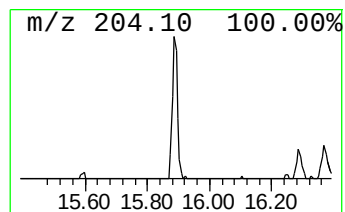
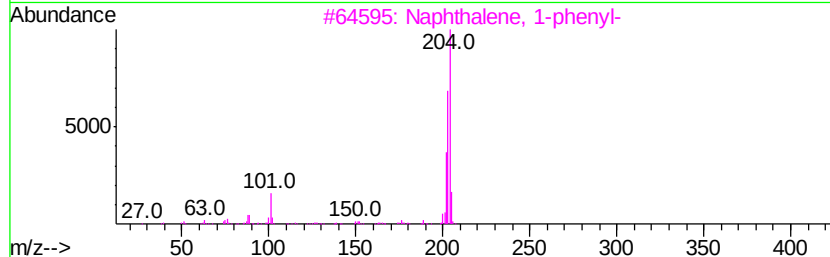
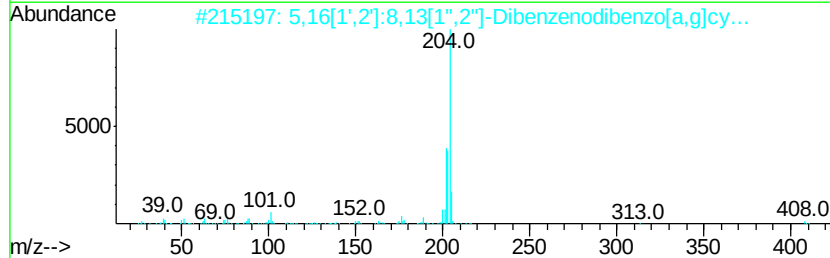
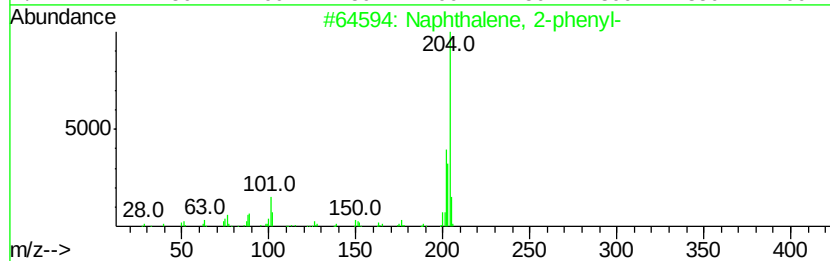
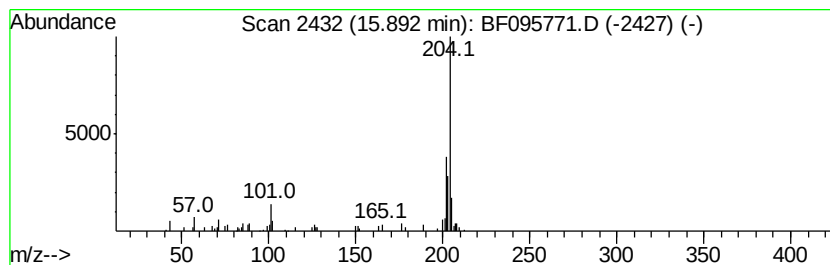
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Naphthalene, 2-phenyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.89	3.15 ng	71239	Phenanthrene-d10	14.63	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	93
2	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	87
3	Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	78
4	9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	62
5	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	62



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060717\
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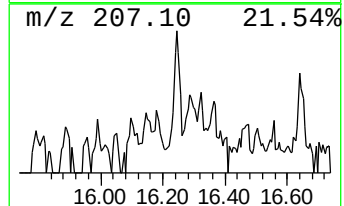
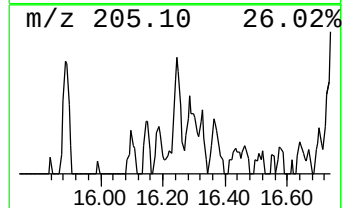
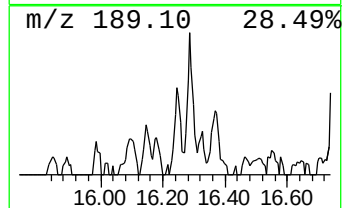
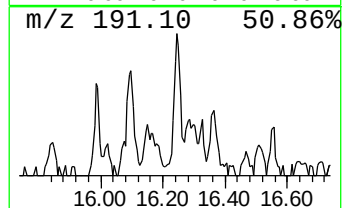
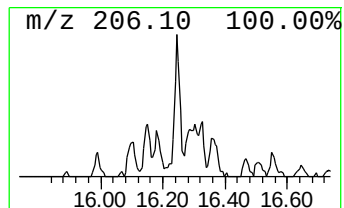
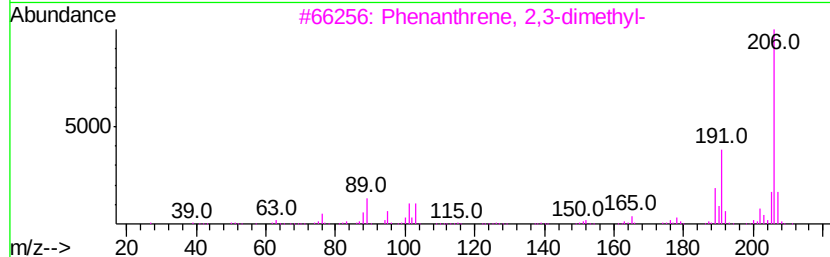
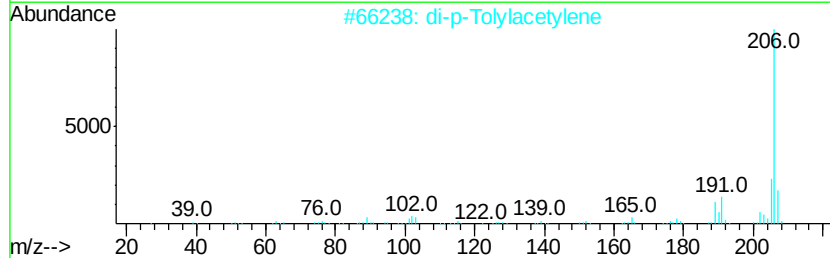
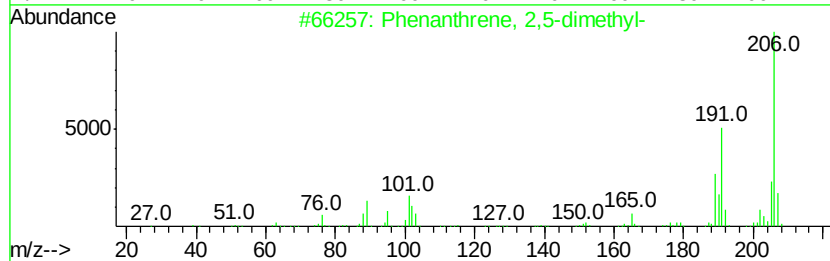
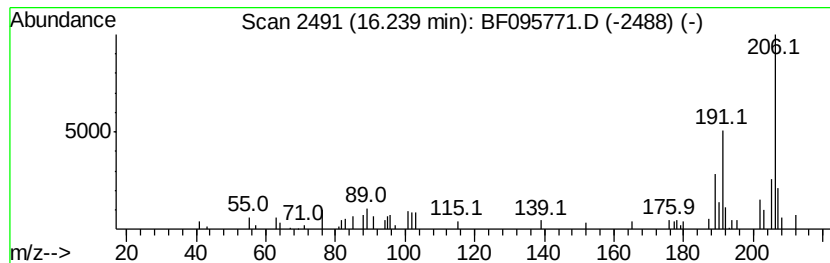
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Phenanthrene, 2,5-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.24	2.63 ng	59507	Phenanthrene-d10		14.63	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2		di-p-Tolylacetylene	206	C16H14	002789-88-0	93
3		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	93
4		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	91
5		9,10-Dimethylantracene	206	C16H14	000781-43-1	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060717\
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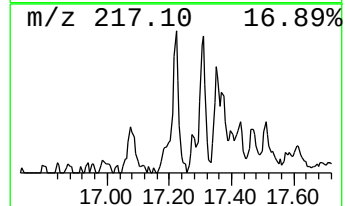
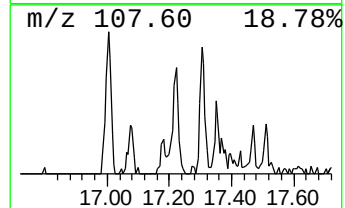
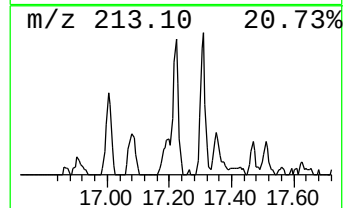
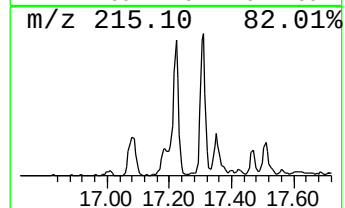
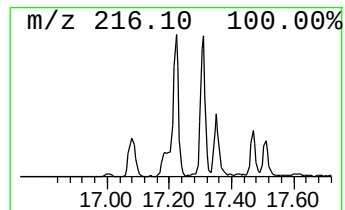
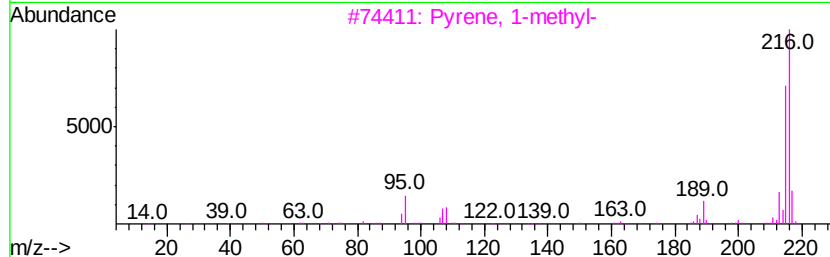
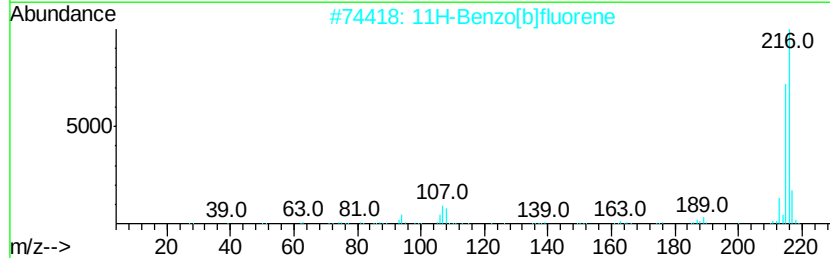
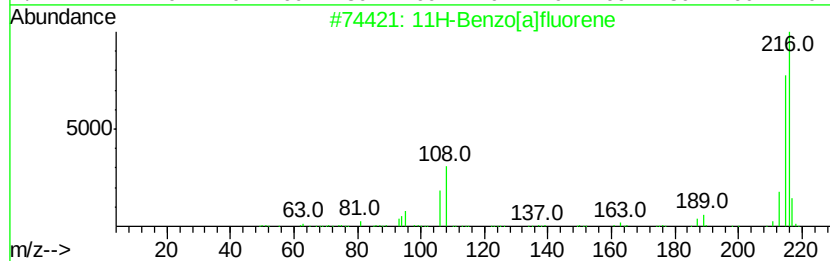
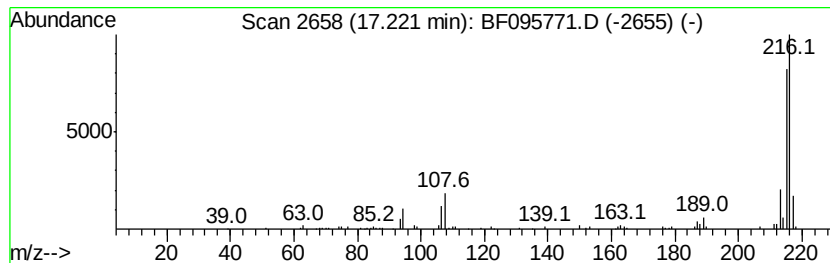
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 11H-Benzo[a]fluorene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.22	4.05 ng	169133	Chrysene-d12	18.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
3			Pvrene, 1-methyl-	216	C17H12	002381-21-7	87
4			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	74
5			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	70



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060717\
Data File : BF095771.D
Acq On : 8 Jun 2017 5:23
Operator : SJ/MA
Sample : I3469-12
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-3(0-2)20170603

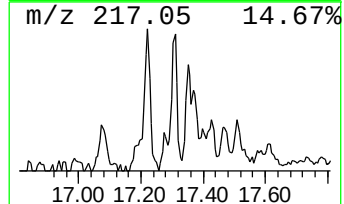
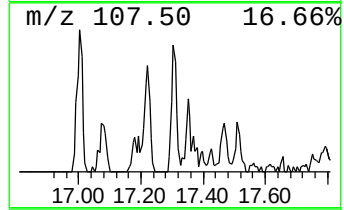
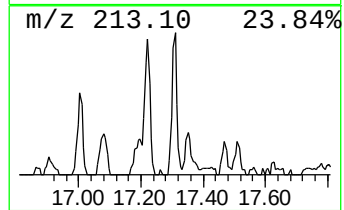
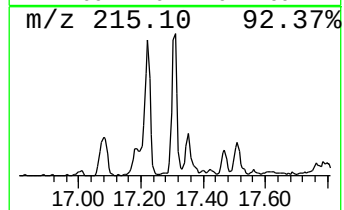
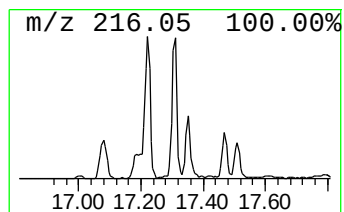
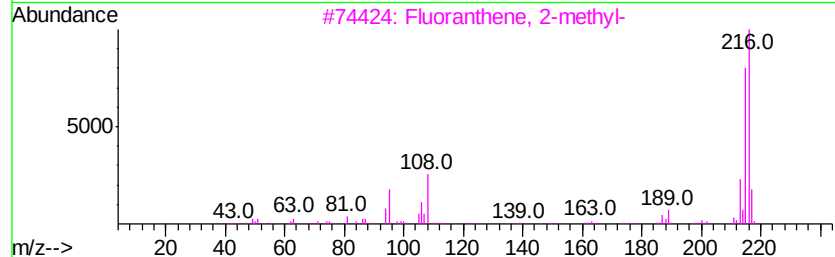
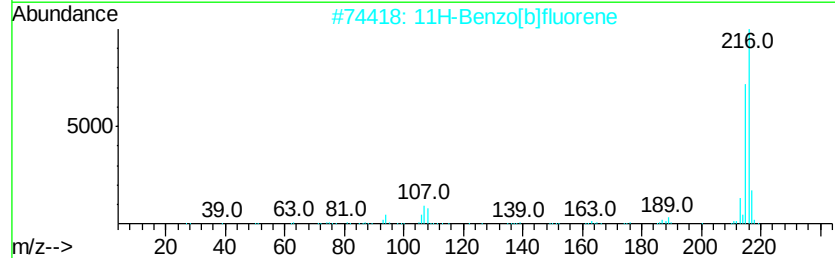
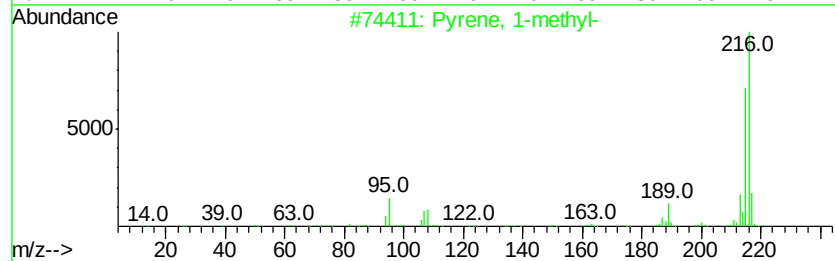
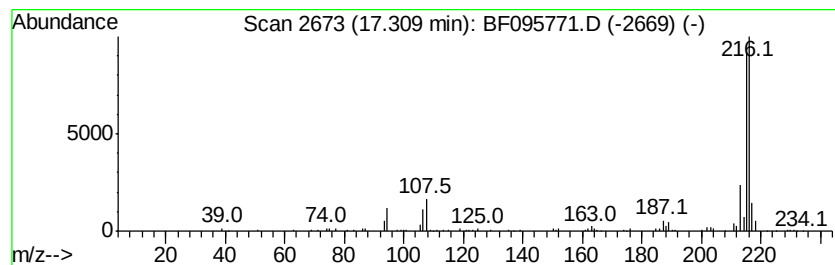
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Pyrene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.31	4.21 ng	175889	Chrysene-d12	18.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
3			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87
4			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	80
5			Pyrene, 2-methyl-	216	C17H12	003442-78-2	76



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060717\
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Sample : I3469-12
Misc :
ALS Vial : 27 Sample Multiplier: 1

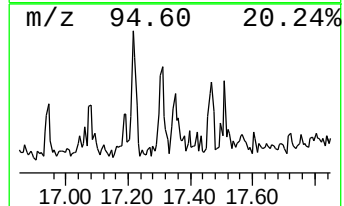
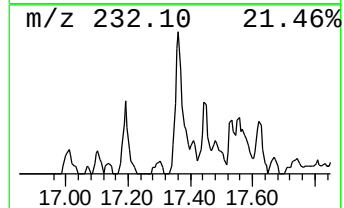
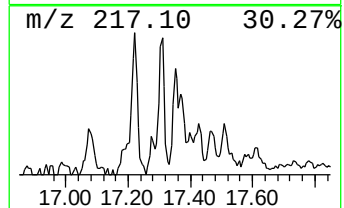
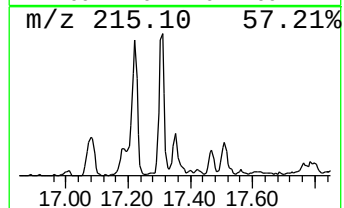
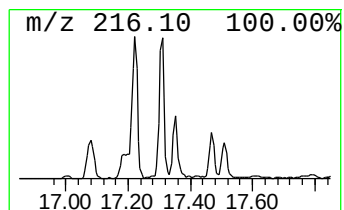
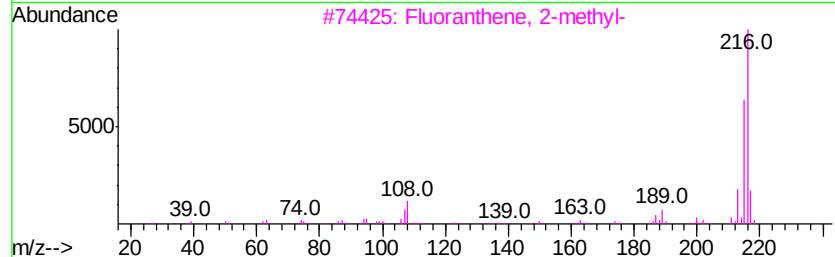
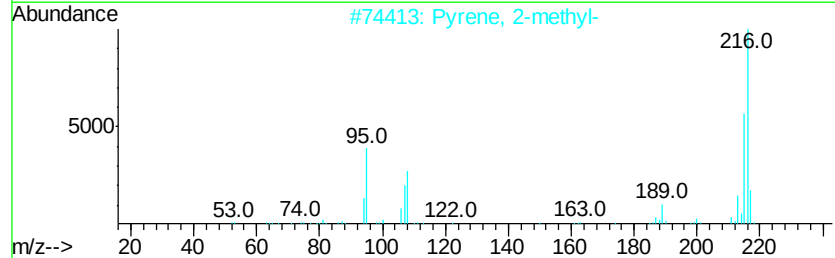
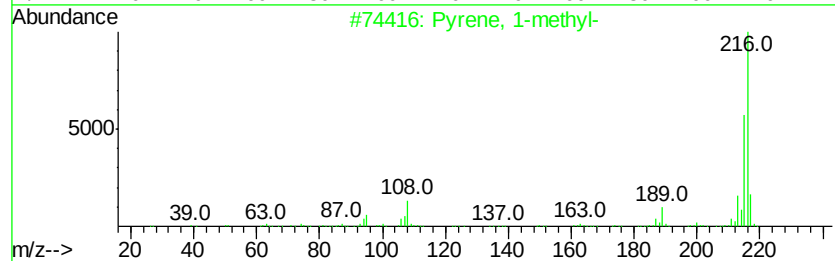
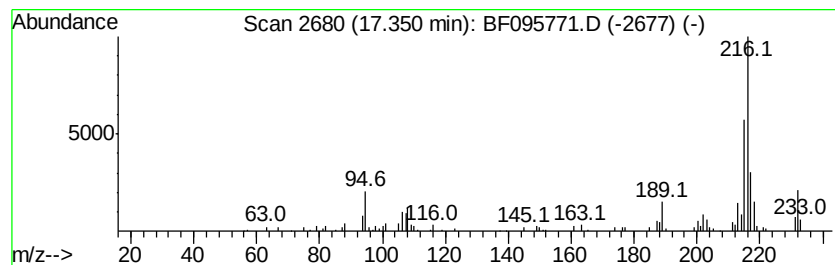
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 11H-Benzo[b]fluorene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.35	3.77 ng	157226	Chrysene-d12	18.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Pyrene, 1-methyl-	216 C17H12	002381-21-7 92
2		Pyrene, 2-methyl-	216 C17H12	003442-78-2 90
3		Fluoranthene, 2-methyl-	216 C17H12	033543-31-6 60
4		11H-Benzo[a]fluorene	216 C17H12	000238-84-6 58
5		11H-Benzo[b]fluorene	216 C17H12	000243-17-4 53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060717\
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Operator : SJ/MA
Sample : I3469-12
Misc :
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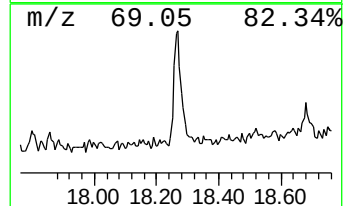
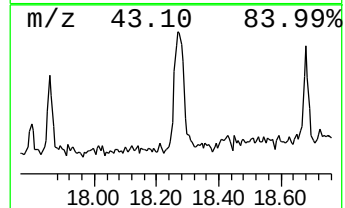
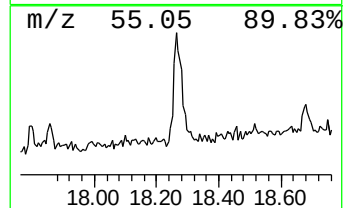
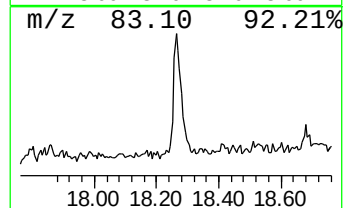
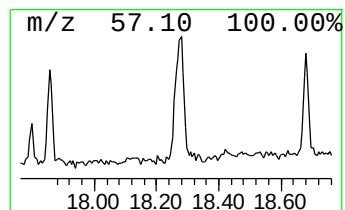
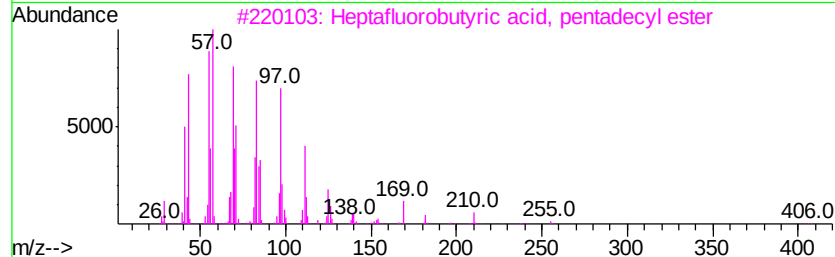
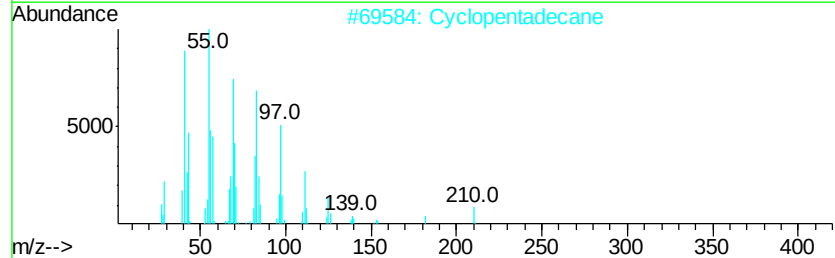
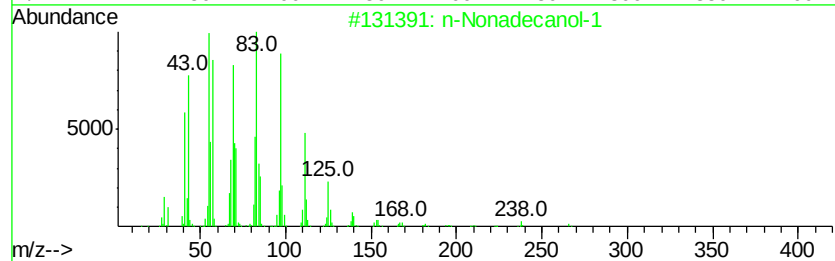
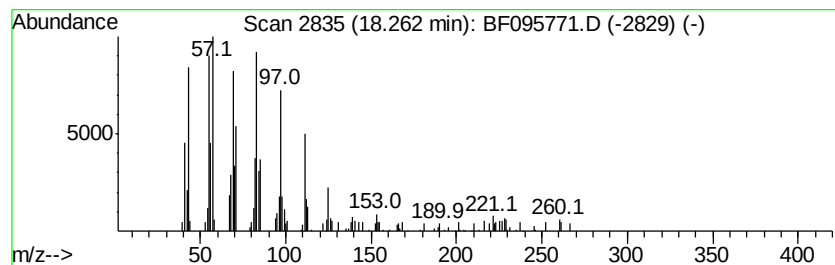
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 n-Nonadecanol-1 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.26	5.68 ng	237058	Chrysene-d12	18.35	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Nonadecanol-1	284	C19H40O	001454-84-8	95
2	Cyclopentadecane	210	C15H30	000295-48-7	94
3	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
4	Behenic alcohol	326	C22H46O	000661-19-8	93
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060717\
Data File : BF095771.D
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Operator : SJ/MA
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ALS Vial : 27 Sample Multiplier: 1

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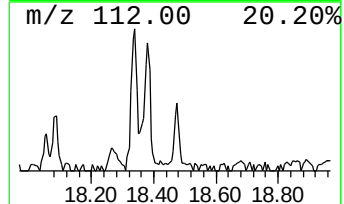
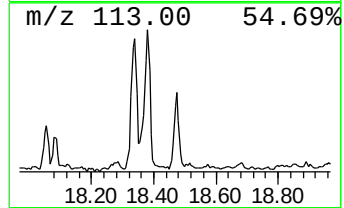
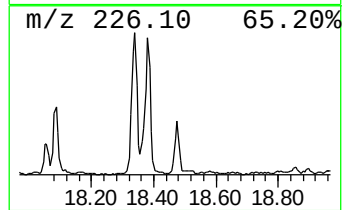
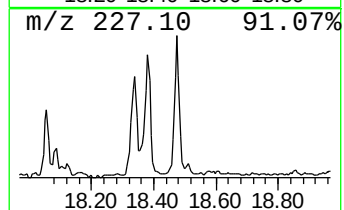
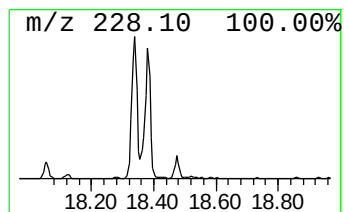
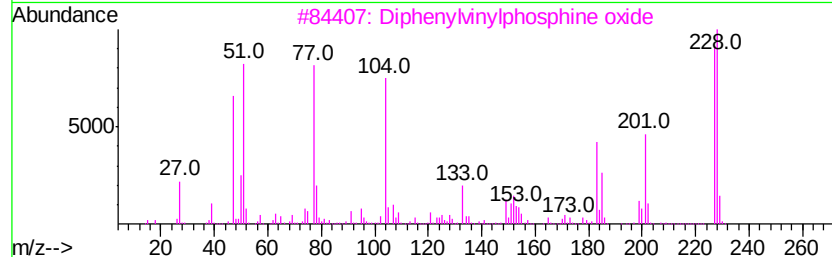
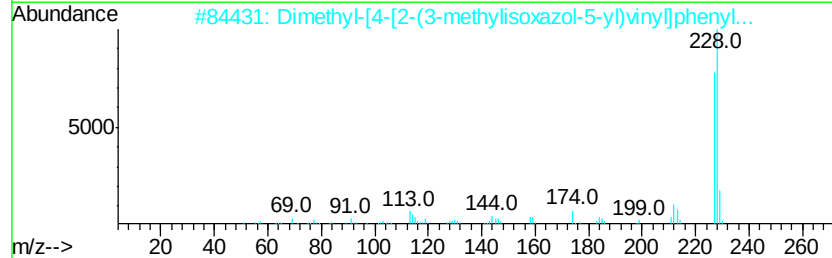
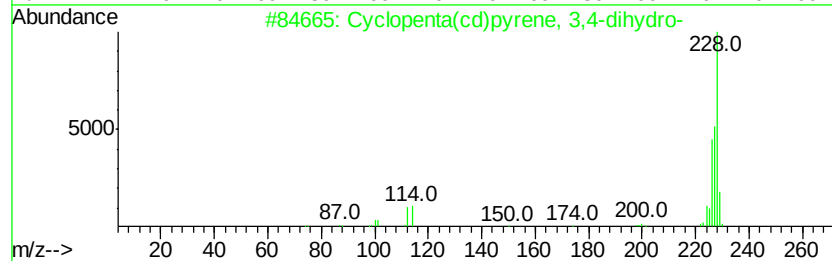
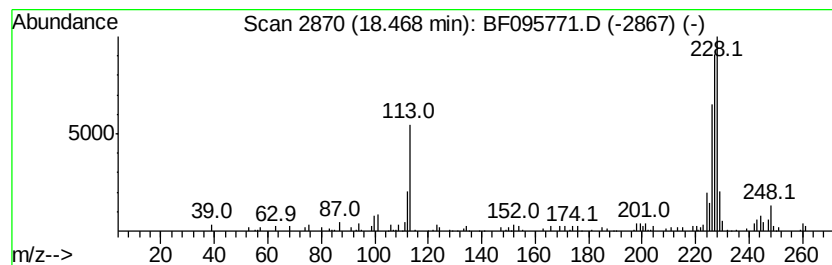
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Cyclopenta(cd)pyrene, 3,4-d... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.47	2.55 ng	106281	Chrysene-d12	18.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(cd)pyrene, 3,4-dihydro-	228	C18H12	025732-74-5	50
2		Dimethyl-[4-[2-(3-methylisoxazol...	228	C14H16N2O	1000306-39-6	49
3		Diphenylvinylphosphine oxide	228	C14H13OP	002096-78-8	46
4		1,3,5-Triazine-2(1H)-thione, 4-(...	227	C9H17N5S	023613-02-7	46
5		Isothiazol-5-amine, 4-bromo-3-ch...	226	C4H4BrClN2S	1000272-59-9	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060717\
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Acq On : 8 Jun 2017 5:23
Operator : SJ/MA
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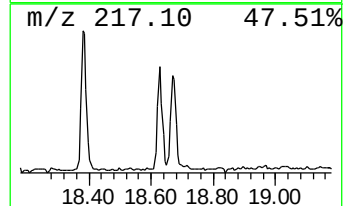
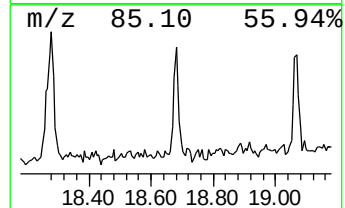
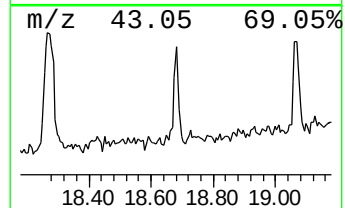
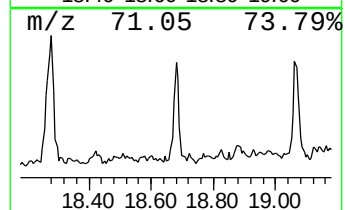
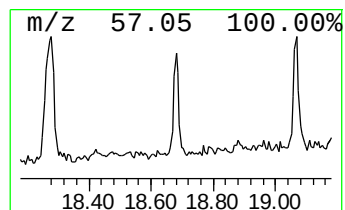
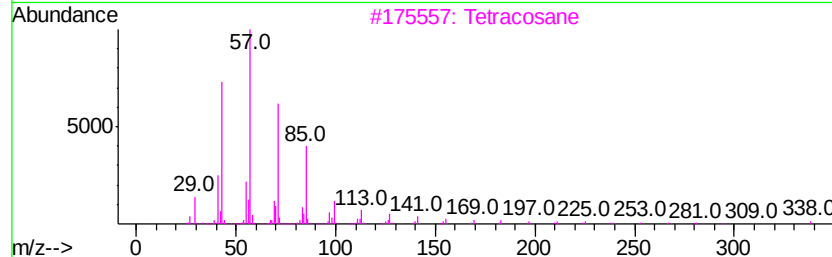
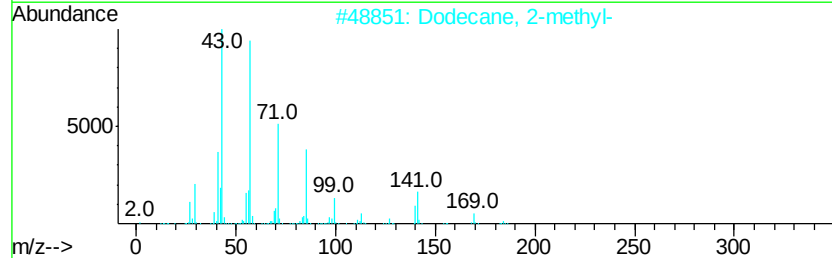
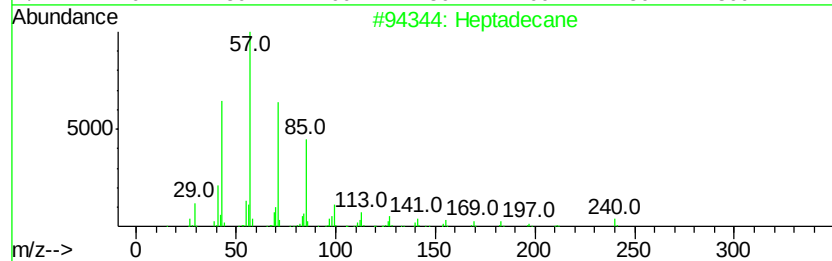
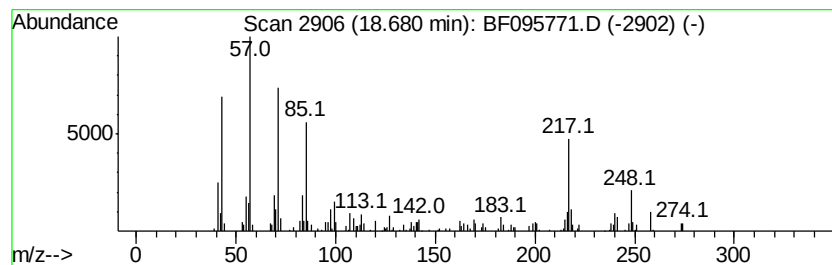
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Heptadecane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.68	2.26 ng	94420	Chrysene-d12	18.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	49
2		Dodecane, 2-methyl-	184	C13H28	001560-97-0	45
3		Tetracosane	338	C24H50	000646-31-1	45
4		Tridecane	184	C13H28	000629-50-5	45
5		Nonadecane	268	C19H40	000629-92-5	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060717\
 Data File : BF095771.D
 Acq On : 8 Jun 2017 5:23
 Operator : SJ/MA
 Sample : I3469-12
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 ALS Vial : 27 Sample Multiplier: 1

Instrument :
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 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
2-Pentanone, 4-hy...	4.59	7.8	ng	159012	1	7.39	407172	20.0
unknown7.04	7.04	96.3	ng	1960560	1	7.39	407172	20.0
Tridecane	13.87	2.2	ng	49612	4	14.63	451982	20.0
Phenanthrene, 1-m...	15.43	3.9	ng	87019	4	14.63	451982	20.0
Phenanthrene, 2-m...	15.47	4.7	ng	105612	4	14.63	451982	20.0
4H-Cyclopenta[def...	15.59	8.3	ng	186768	4	14.63	451982	20.0
1H-Indene, 2-phenyl-	15.63	2.9	ng	65902	4	14.63	451982	20.0
n-Hexadecanoic acid	15.65	4.7	ng	105102	4	14.63	451982	20.0
Naphthalene, 2-ph...	15.89	3.1	ng	71239	4	14.63	451982	20.0
Phenanthrene, 2,5...	16.24	2.6	ng	59507	4	14.63	451982	20.0
11H-Benzo[a]fluorene	17.22	4.0	ng	169133	5	18.35	834990	20.0
Pyrene, 1-methyl-	17.31	4.2	ng	175889	5	18.35	834990	20.0
11H-Benzo[b]fluorene	17.35	3.8	ng	157226	5	18.35	834990	20.0
n-Nonadecanol-1	18.26	5.7	ng	237058	5	18.35	834990	20.0
Cyclopenta(cd)pyr...	18.47	2.5	ng	106281	5	18.35	834990	20.0
Heptadecane	18.68	2.3	ng	94420	5	18.35	834990	20.0