

Data Path : Z:\HPCHEM1\BNA F\DATA\BF053117\
 Data File : BF095573.D
 Acq On : 31 May 2017 20:08
 Operator : SJ/MA
 Sample : I3329-07
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-8(0-3)

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	4.807	501	505	525	rBV2	76206	215554	7.90%	0.695%
2	5.466	614	617	648	rBV	1042486	2182633	80.03%	7.033%
3	7.101	891	895	909	rBV	1344453	2203623	80.80%	7.101%
4	7.207	909	913	936	rVB	1638507	2613852	95.84%	8.423%
5	7.554	968	972	983	rBV	450579	633498	23.23%	2.041%
6	7.789	1007	1012	1037	rVB	1356352	1884353	69.09%	6.072%
7	8.466	1122	1127	1145	rBV	954290	1461531	53.59%	4.710%
8	9.583	1312	1317	1329	rBV	622588	902818	33.10%	2.909%
9	11.360	1613	1619	1630	rBV	2135607	2727376	100.00%	8.789%
10	11.895	1705	1710	1714	rBV	20096	30220	1.11%	0.097%
11	12.048	1732	1736	1742	rBV	280959	367050	13.46%	1.183%
12	12.407	1792	1797	1804	rBV2	706086	922229	33.81%	2.972%
13	12.460	1804	1806	1816	rVB	50514	72447	2.66%	0.233%
14	13.254	1936	1941	1945	rBV	44367	53091	1.95%	0.171%
15	13.313	1945	1951	1961	rVB2	51466	75939	2.78%	0.245%
16	13.583	1992	1997	2000	rBV	24743	35612	1.31%	0.115%
17	13.707	2013	2018	2029	rBV	1034545	1536201	56.33%	4.950%
18	14.024	2068	2072	2075	rBV	54289	72971	2.68%	0.235%
19	14.212	2098	2104	2108	rBV3	17264	30699	1.13%	0.099%
20	14.442	2138	2143	2146	rBV4	22511	34982	1.28%	0.113%
21	14.548	2156	2161	2166	rBV	32634	48234	1.77%	0.155%
22	14.618	2166	2173	2175	rVV5	20271	39987	1.47%	0.129%
23	14.648	2175	2178	2182	rVB	30909	37369	1.37%	0.120%
24	14.759	2192	2197	2201	rBV	42547	53030	1.94%	0.171%
25	14.812	2201	2206	2209	rVV	786255	991354	36.35%	3.195%
26	14.848	2209	2212	2220	rVV	612441	784765	28.77%	2.529%
27	14.942	2224	2228	2234	rBV	129342	173599	6.37%	0.559%
28	15.036	2240	2244	2248	rBV4	20166	32537	1.19%	0.105%
29	15.242	2276	2279	2287	rBV2	29107	53183	1.95%	0.171%
30	15.430	2307	2311	2315	rVB	28143	34984	1.28%	0.113%
31	15.607	2337	2341	2344	rVB	113681	124610	4.57%	0.402%
32	15.648	2344	2348	2355	rBV	136672	164658	6.04%	0.531%
33	15.718	2357	2360	2363	rBV	42268	48158	1.77%	0.155%
34	15.754	2363	2366	2370	rVV2	122147	174044	6.38%	0.561%

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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

35	15.795	2370	2373	2382	rVB	178029	231465	8.49%	0.746%
36	16.024	2408	2412	2414	rBV4	23920	28564	1.05%	0.092%
37	16.048	2414	2416	2425	rVB	70656	92493	3.39%	0.298%
38	16.118	2425	2428	2431	rBV	34299	41447	1.52%	0.134%
39	16.259	2448	2452	2456	rBV2	24417	32113	1.18%	0.103%
40	16.306	2456	2460	2464	rBV2	33877	42767	1.57%	0.138%
41	16.406	2470	2477	2481	rBV	98420	121850	4.47%	0.393%
42	16.453	2481	2485	2488	rVV3	41019	60775	2.23%	0.196%
43	16.530	2493	2498	2503	rBV4	49991	80965	2.97%	0.261%
44	16.606	2507	2511	2517	rVV	781785	911599	33.42%	2.938%
45	16.648	2517	2518	2523	rVB2	32133	36255	1.33%	0.117%
46	16.712	2525	2529	2534	rBV3	33026	60665	2.22%	0.195%
47	16.806	2542	2545	2549	rVB2	25230	31787	1.17%	0.102%
48	16.912	2556	2563	2569	rVV	873318	1008837	36.99%	3.251%
49	17.036	2581	2584	2588	rVB	25440	27561	1.01%	0.089%
50	17.106	2593	2596	2599	rBV	40275	38575	1.41%	0.124%
51	17.159	2599	2605	2612	rVB	1868818	2188853	80.25%	7.053%
52	17.236	2612	2618	2627	rBV4	48120	109475	4.01%	0.353%
53	17.353	2631	2638	2640	rVV4	43605	81962	3.01%	0.264%
54	17.383	2640	2643	2649	rVB	88245	105442	3.87%	0.340%
55	17.471	2654	2658	2661	rVB	50477	60664	2.22%	0.195%
56	17.512	2661	2665	2670	rBV3	105850	158689	5.82%	0.511%
57	17.630	2679	2685	2689	rBV2	38178	47201	1.73%	0.152%
58	17.671	2689	2692	2698	rVB2	28656	33710	1.24%	0.109%
59	18.071	2755	2760	2764	rBV2	52558	58602	2.15%	0.189%
60	18.183	2772	2779	2781	rBV4	36997	52444	1.92%	0.169%
61	18.212	2781	2784	2788	rVV	75618	83707	3.07%	0.270%
62	18.247	2788	2790	2795	rVV	38011	52495	1.92%	0.169%
63	18.289	2795	2797	2801	rVB	29375	28816	1.06%	0.093%
64	18.342	2804	2806	2812	rVB	44826	49801	1.83%	0.160%
65	18.400	2812	2816	2827	rBV2	143288	248348	9.11%	0.800%
66	18.500	2827	2833	2837	rVV2	739422	1081882	39.67%	3.486%
67	18.536	2837	2839	2846	rVB2	226848	336444	12.34%	1.084%
68	18.636	2851	2856	2859	rVV2	46086	62360	2.29%	0.201%
69	18.736	2867	2873	2876	rBV6	20746	38010	1.39%	0.122%
70	18.795	2880	2883	2888	rVV5	31581	60193	2.21%	0.194%
71	18.836	2888	2890	2897	rVB5	27189	33790	1.24%	0.109%

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72	19.012	2914	2920	2924	rBV2	66617	83780	3.07%	0.270%
73	19.053	2924	2927	2931	rVB3	39522	45797	1.68%	0.148%
74	19.200	2949	2952	2958	rVB6	28514	44588	1.63%	0.144%
75	19.530	3000	3008	3013	rBV2	118200	158486	5.81%	0.511%
76	19.630	3020	3025	3029	rVB4	27311	42481	1.56%	0.137%
77	19.777	3046	3050	3053	rBV	230638	312129	11.44%	1.006%
78	19.806	3053	3055	3060	rVB2	81607	105827	3.88%	0.341%
79	19.889	3066	3069	3072	rBV	40484	43643	1.60%	0.141%
80	20.000	3082	3088	3089	rBV5	21678	34997	1.28%	0.113%
81	20.065	3095	3099	3103	rBV	148742	151562	5.56%	0.488%
82	20.124	3106	3109	3114	rVB	195815	198042	7.26%	0.638%
83	20.177	3114	3118	3123	rBV	451354	566731	20.78%	1.826%
84	20.618	3189	3193	3196	rBV5	19897	29640	1.09%	0.096%
85	20.671	3198	3202	3207	rVB6	22848	40110	1.47%	0.129%
86	21.124	3275	3279	3286	rBV9	15374	29762	1.09%	0.096%
87	21.430	3326	3331	3338	rBV	81422	142594	5.23%	0.459%
88	21.571	3351	3355	3359	rBV2	17926	29291	1.07%	0.094%
89	21.624	3361	3364	3372	rVB4	23546	41356	1.52%	0.133%
90	21.794	3387	3393	3400	rBV	69039	135882	4.98%	0.438%
91	22.018	3425	3431	3442	rVB4	30048	75266	2.76%	0.243%
92	23.641	3702	3707	3717	rVB7	19032	43369	1.59%	0.140%
93	23.812	3729	3736	3740	rBV7	17487	39704	1.46%	0.128%

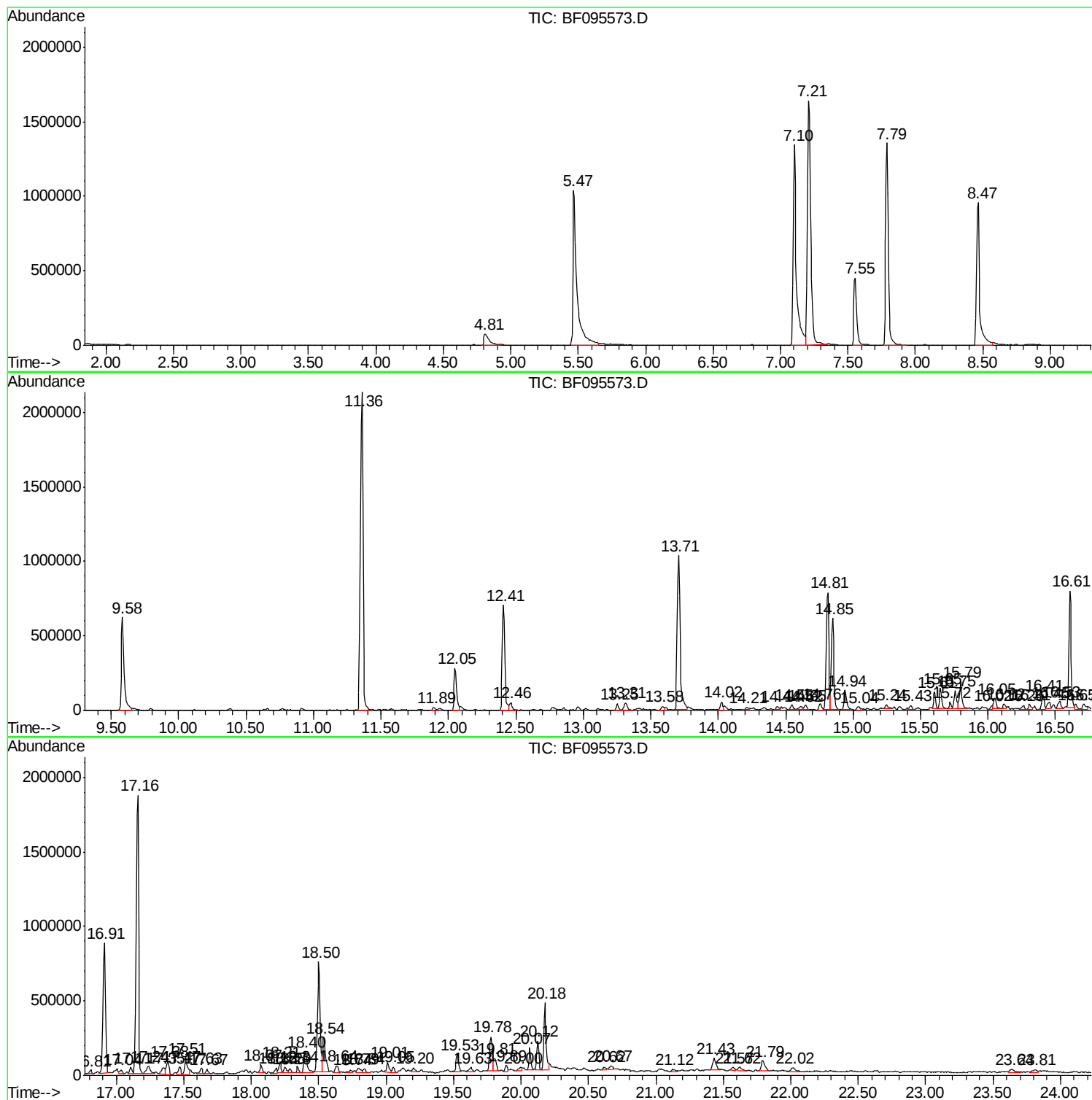
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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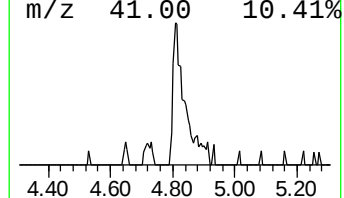
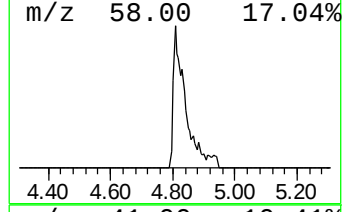
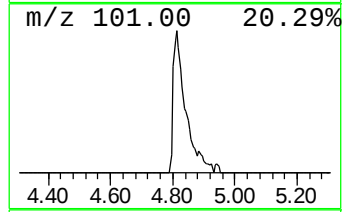
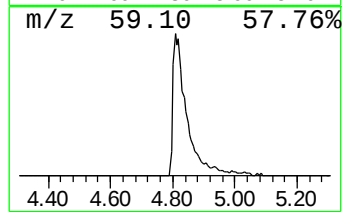
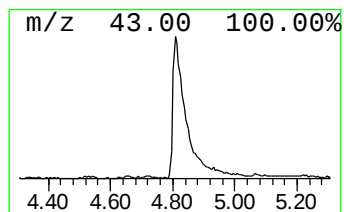
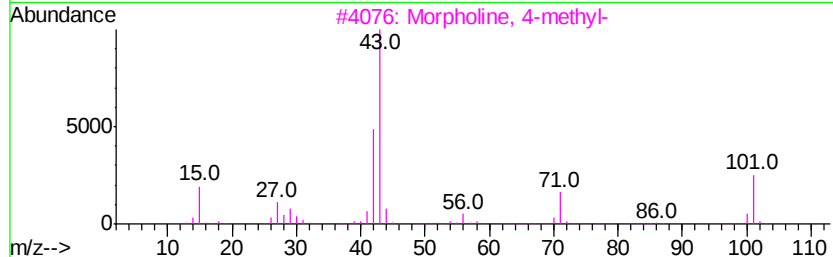
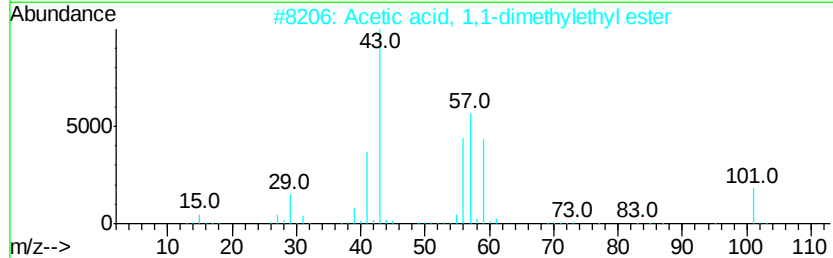
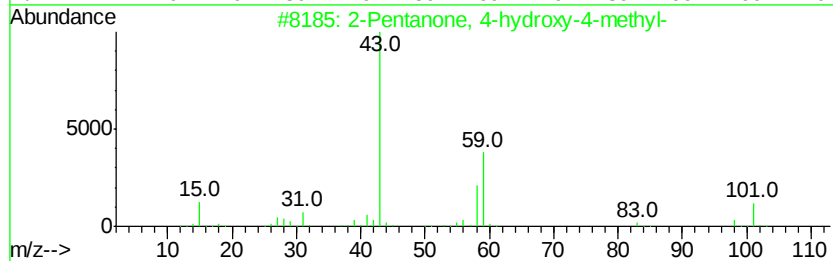
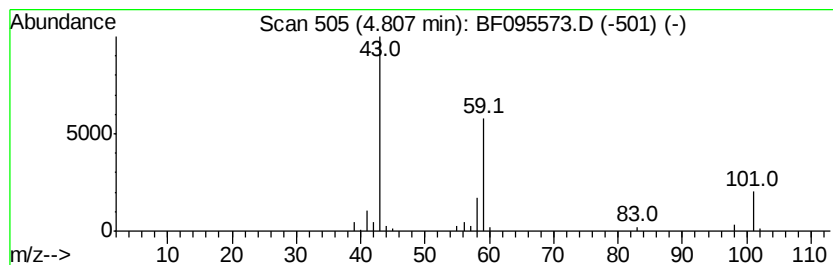
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.81	6.81 ng	215554	1,4-Dichlorobenzene-d4	7.55

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4			Ethanone, 1-(2-methylcyclopropyl)-	98	C6H10O	000930-56-3	9
5			Guanidine	59	CH5N3	000113-00-8	9



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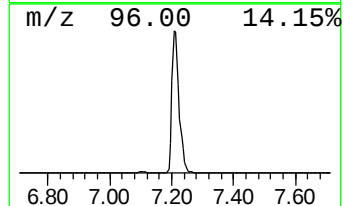
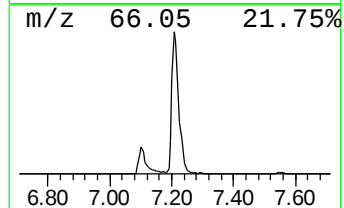
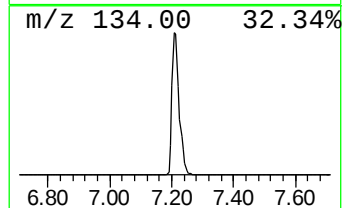
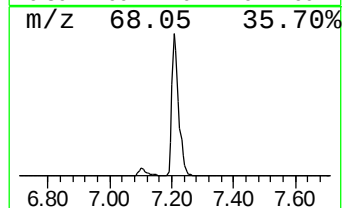
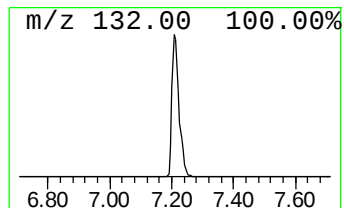
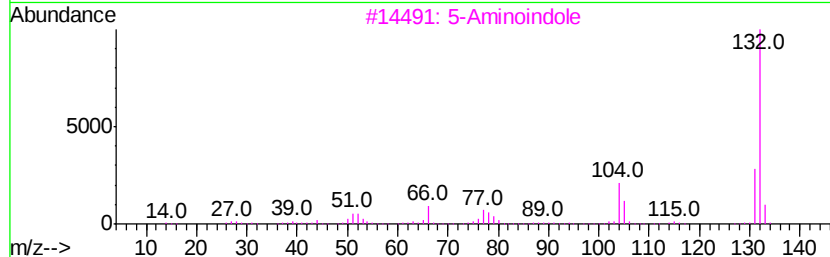
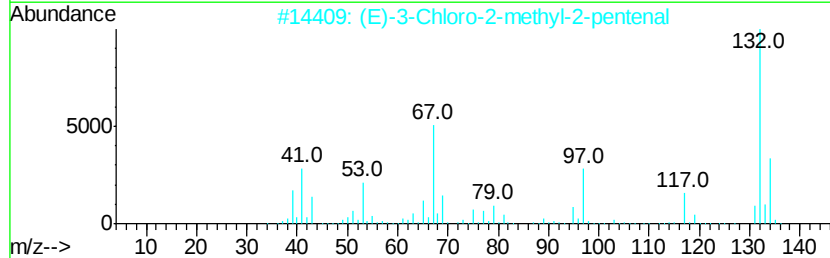
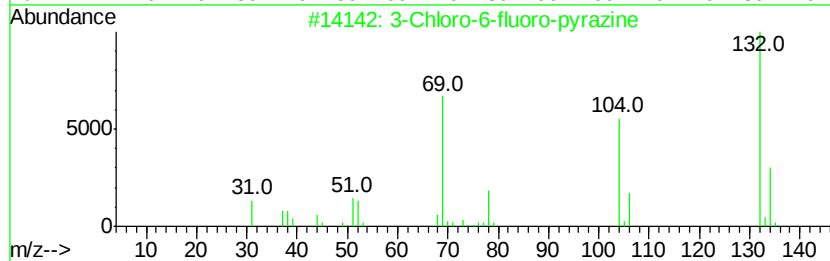
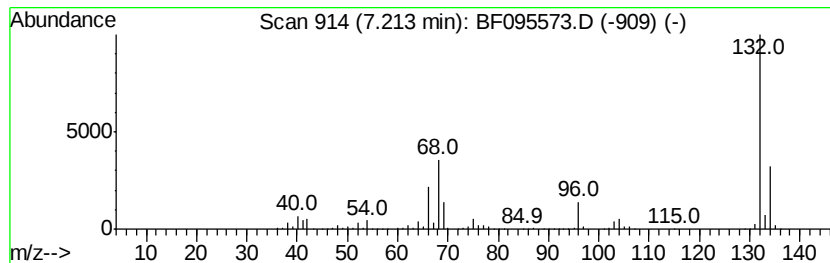
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown7.21 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.21	82.52 ng	2613850	1,4-Dichlorobenzene-d4	7.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		1-Aminoindan	133	C9H11N	034698-41-4	9



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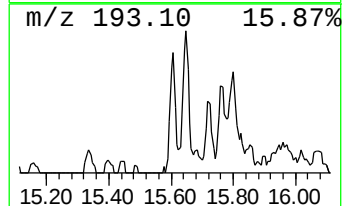
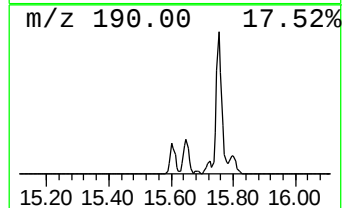
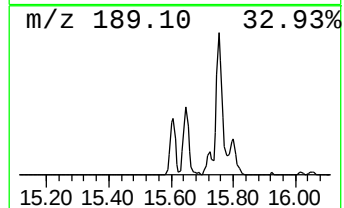
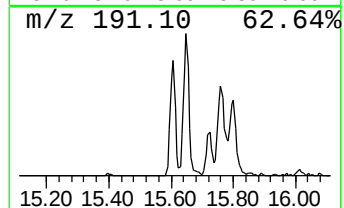
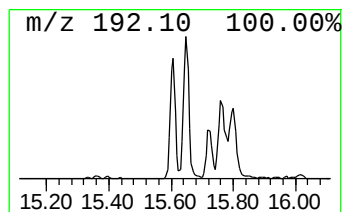
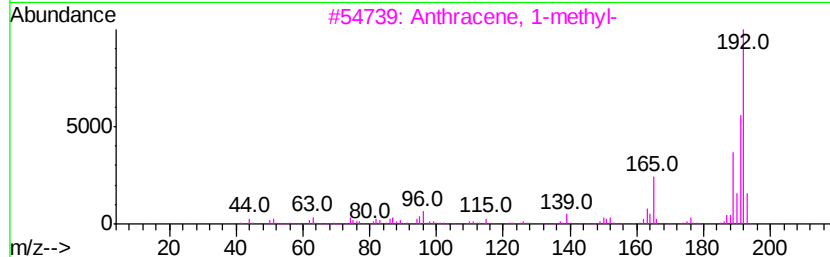
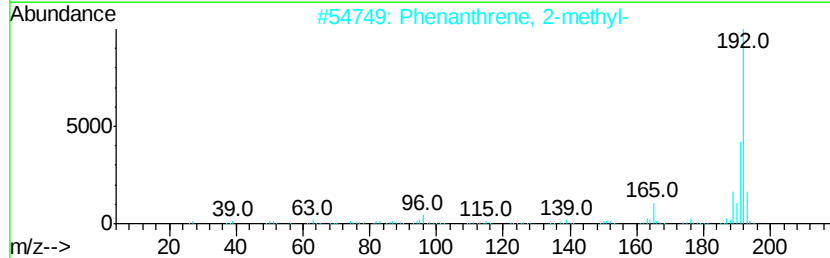
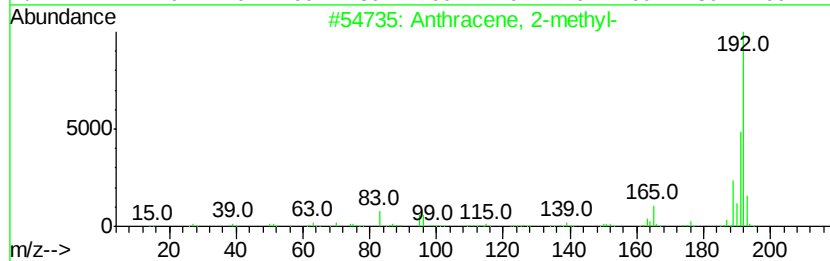
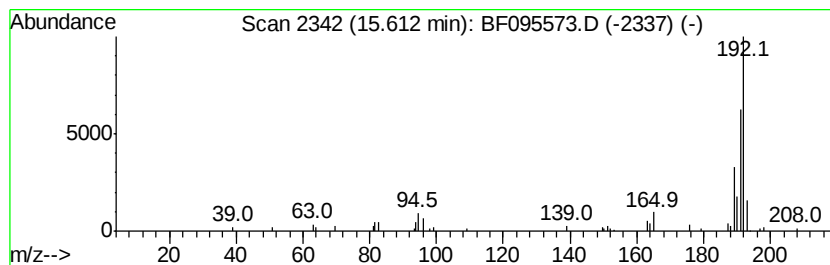
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Anthracene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.61	2.51 ng	124610	Phenanthrene-d10	14.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	95
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
5	Phenanthrene, 9-methyl-	192	C15H12	000883-20-5	91



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ClientSampleId :
B-8(0-3)

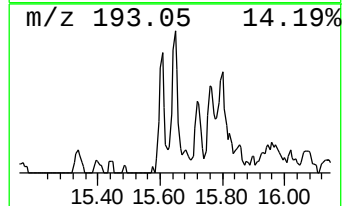
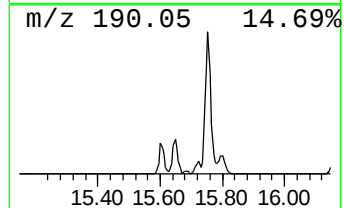
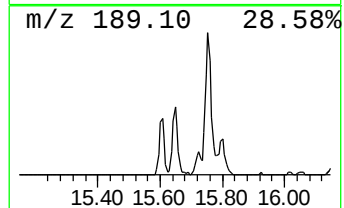
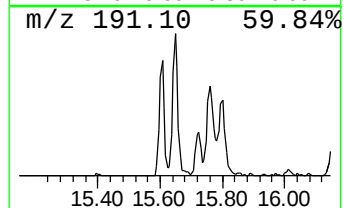
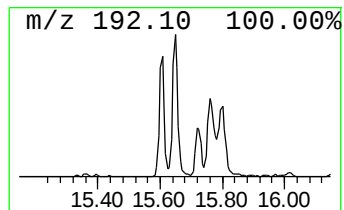
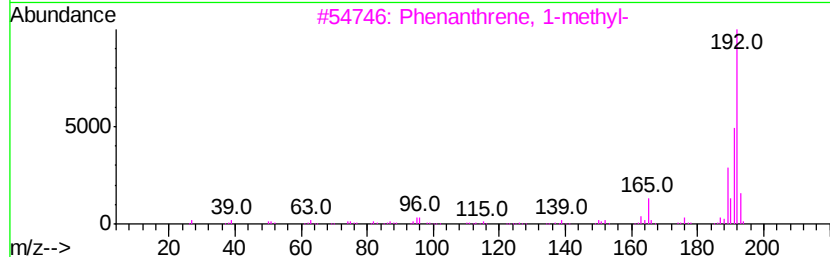
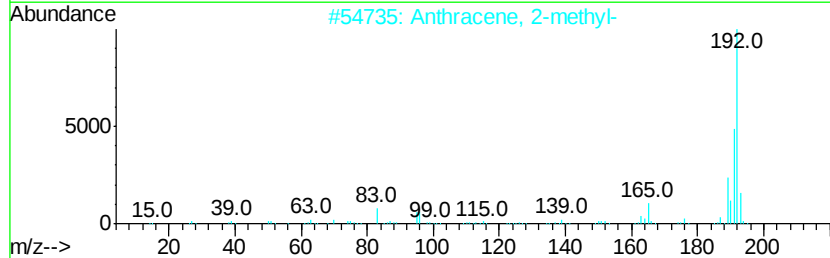
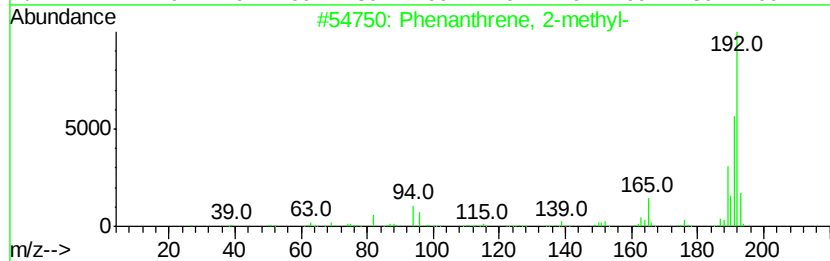
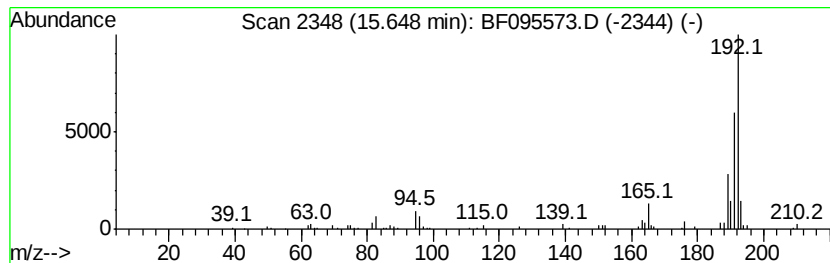
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.65	3.32 ng	164658	Phenanthrene-d10	14.81

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
4		1H-Cyclopropa[1,1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	95
5		5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	95



Data Path : Z:\HPCHEM1\BNA F\DATA\BF053117\
Data File : BF095573.D
Acq On : 31 May 2017 20:08
Operator : SJ/MA
Sample : I3329-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
B-8(0-3)

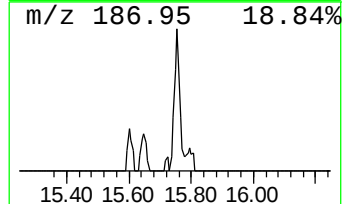
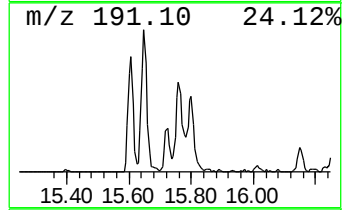
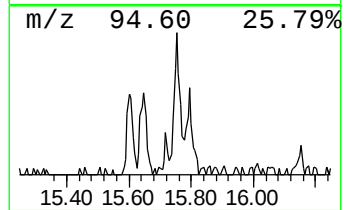
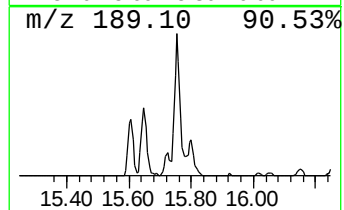
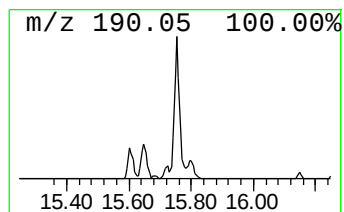
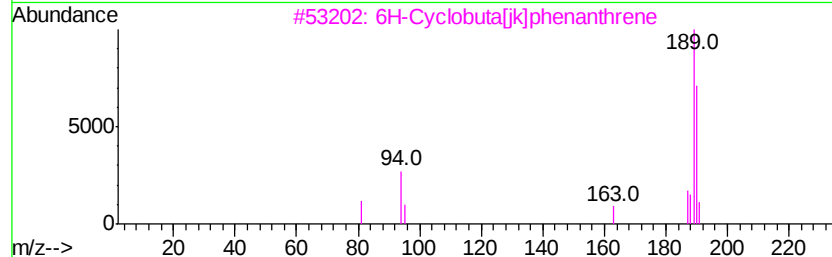
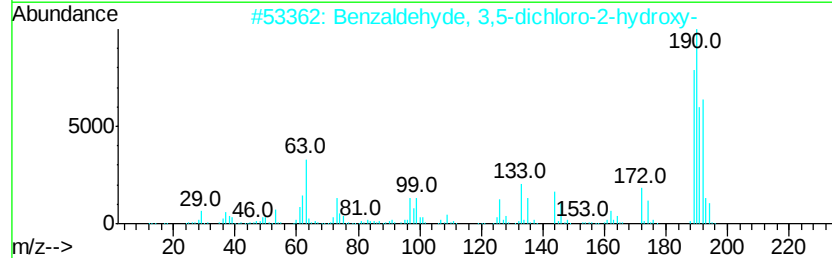
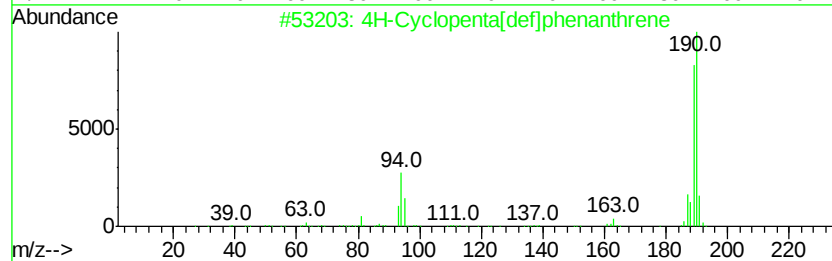
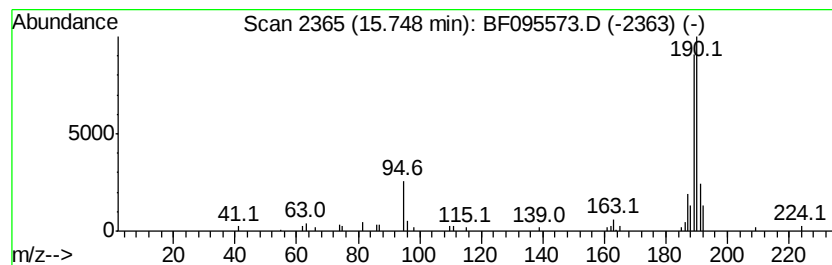
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.75	3.51 ng	174044	Phenanthrene-d10	14.81

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2		Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	50
3		6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	42
4		Methyl diselenide	190	C2H6Se2	007101-31-7	27
5		Benzene, 1,1'-(1,2-propadienyld...	192	C15H12	014251-57-1	22



Data Path : Z:\HPCHEM1\BNA F\DATA\BF053117\
Data File : BF095573.D
Acq On : 31 May 2017 20:08
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Misc :
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ClientSampleId :
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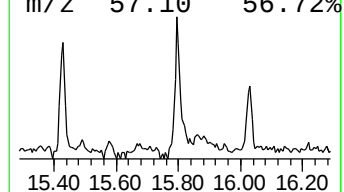
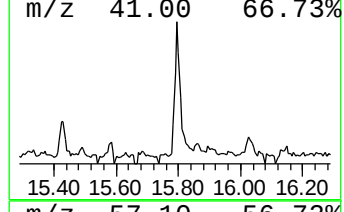
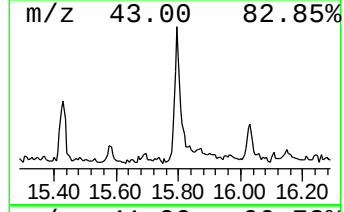
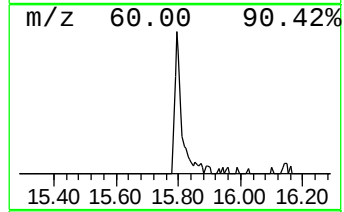
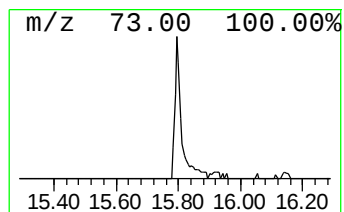
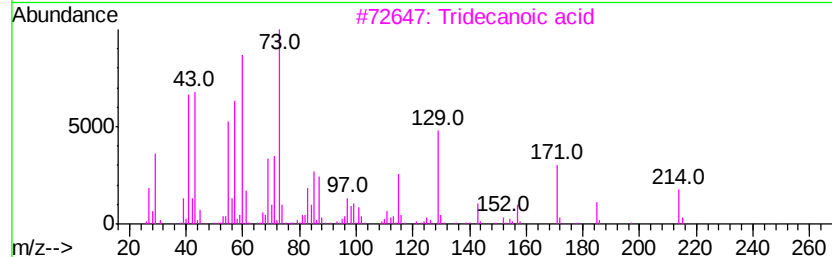
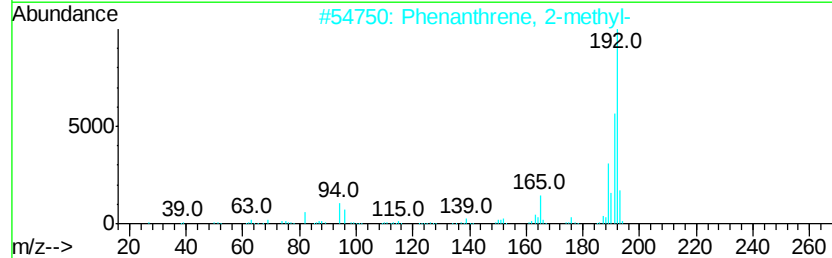
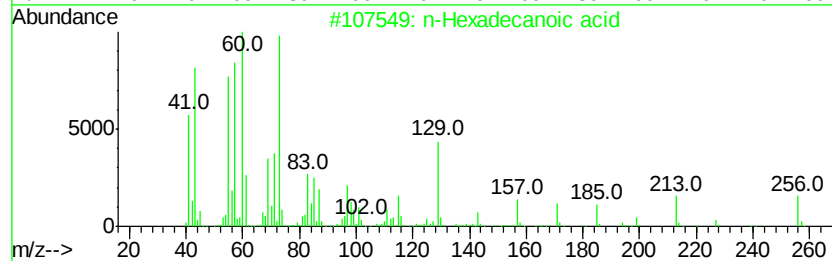
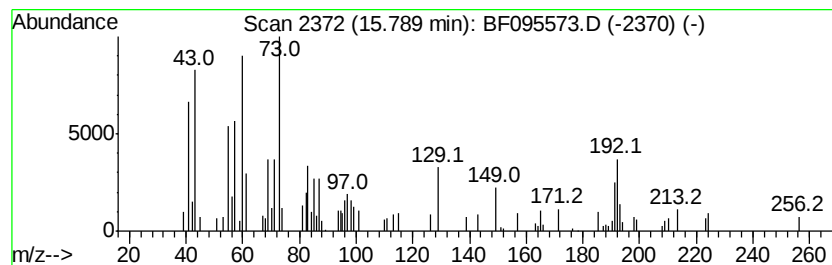
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.79	4.67 ng	231465	Phenanthrene-d10	14.81

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	56
3		Tridecanoic acid	214	C13H26O2	000638-53-9	52
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	42
5		Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	42



Data Path : Z:\HPCHEM1\BNA F\DATA\BF053117\
Data File : BF095573.D
Acq On : 31 May 2017 20:08
Operator : SJ/MA
Sample : I3329-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

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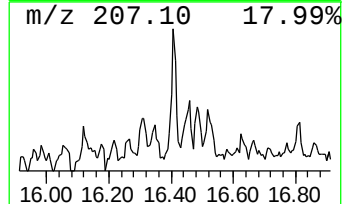
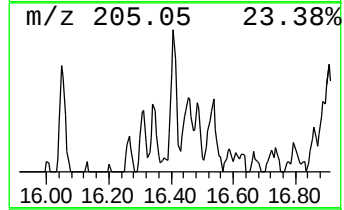
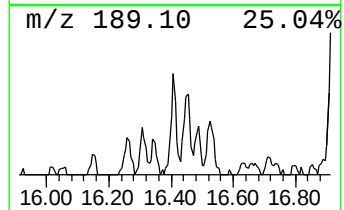
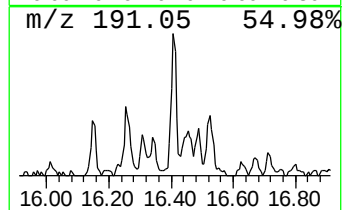
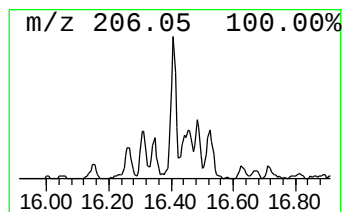
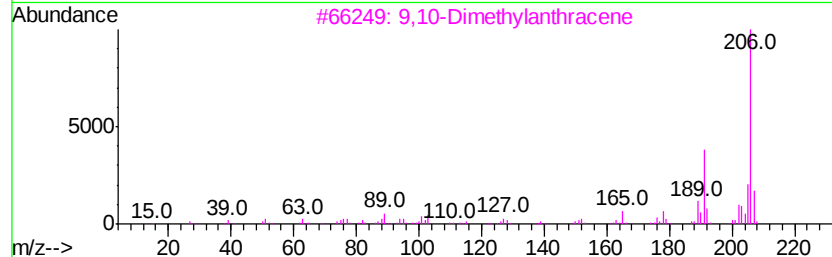
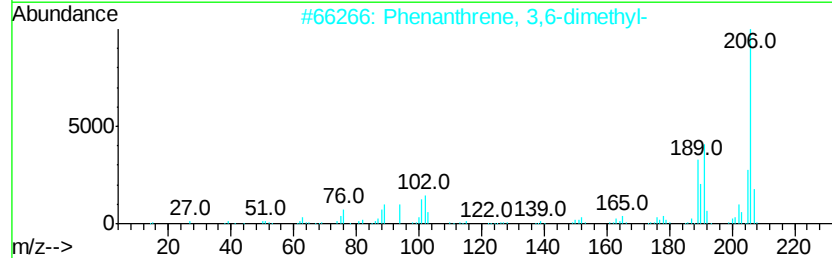
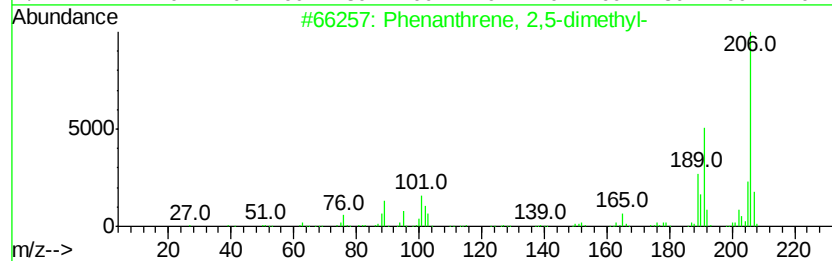
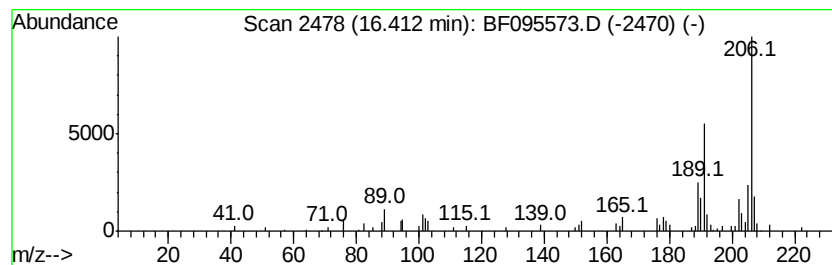
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenanthrene, 2,5-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.41	2.46 ng	121850	Phenanthrene-d10	14.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	97
2		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	93
3		9,10-Dimethylanthracene	206	C16H14	000781-43-1	91
4		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	91
5		Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF053117\
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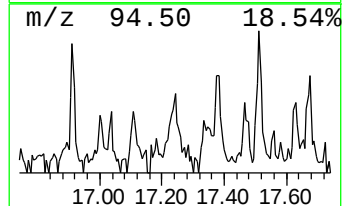
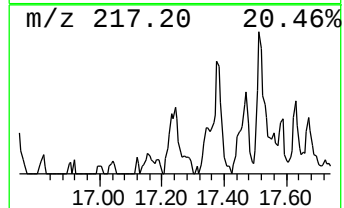
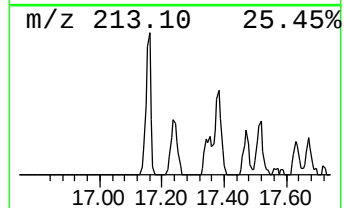
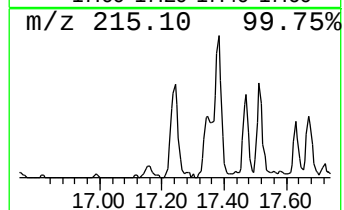
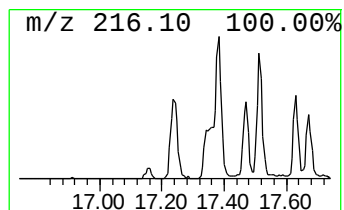
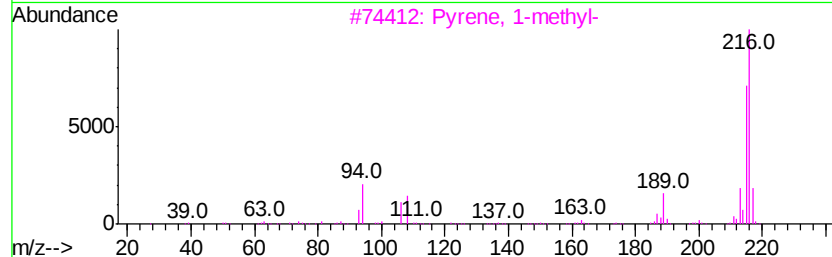
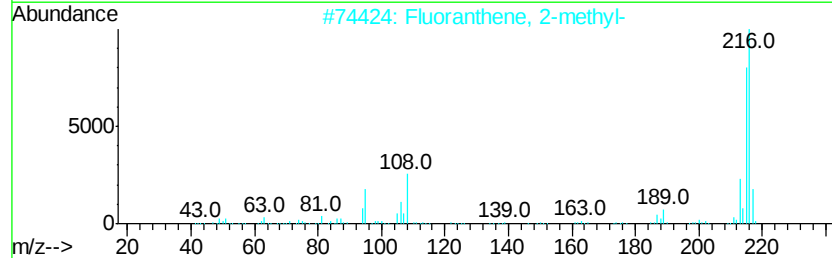
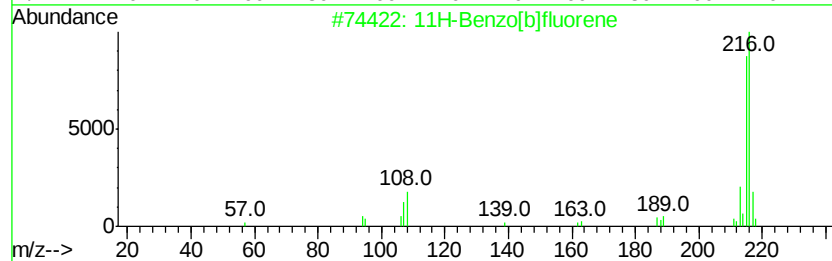
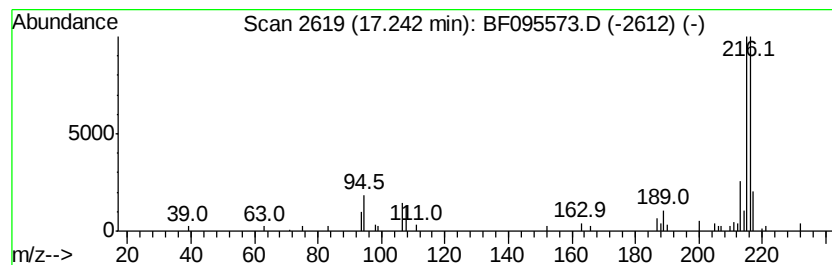
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 11H-Benzo[b]fluorene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.24	2.02 ng	109475	Chrysene-d12	18.51
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4 93
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6 90
3	Pvrene, 1-methyl-	216	C17H12	002381-21-7 89
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6 81
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9 74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF053117\
Data File : BF095573.D
Acq On : 31 May 2017 20:08
Operator : SJ/MA
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Misc :
ALS Vial : 13 Sample Multiplier: 1

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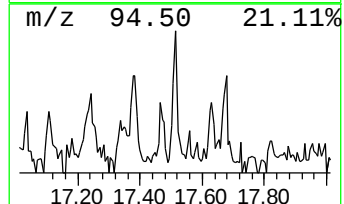
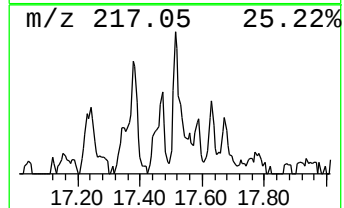
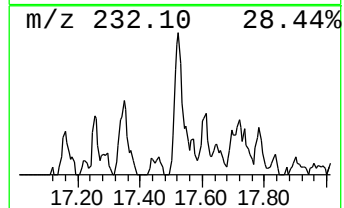
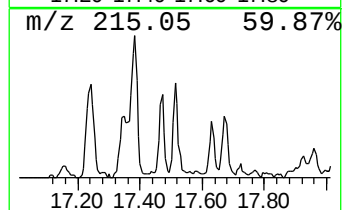
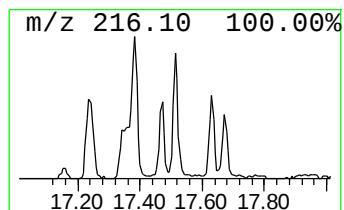
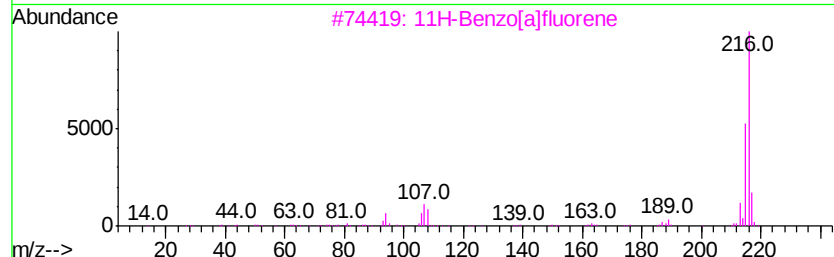
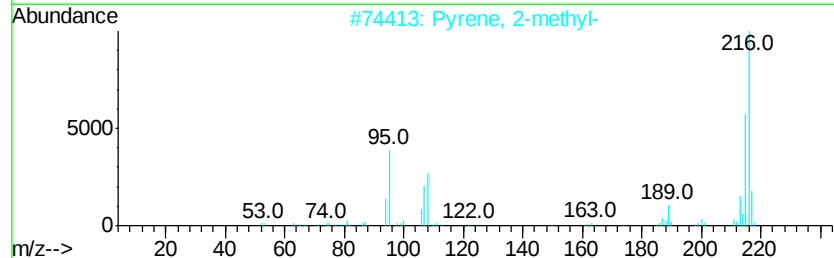
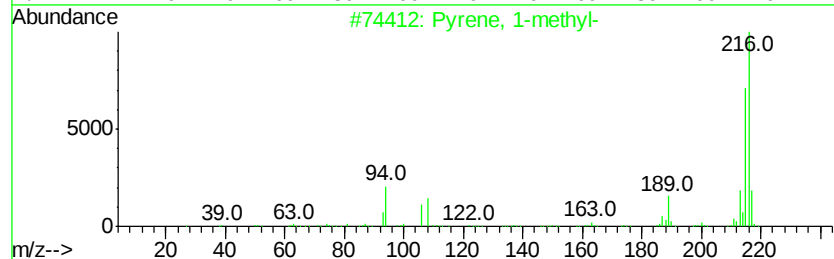
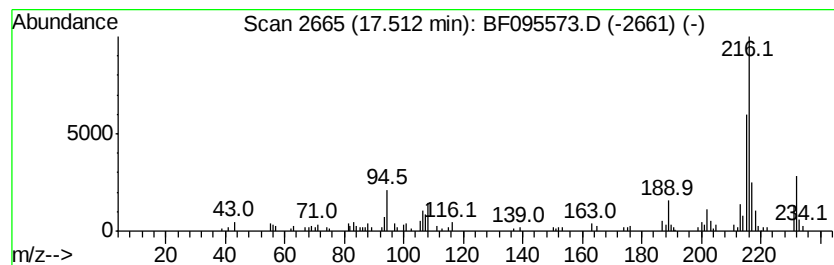
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Pyrene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.51	2.93 ng	158689	Chrysene-d12	18.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	95
2		Pyrene, 2-methyl-	216	C17H12	003442-78-2	83
3		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	76
4		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	64
5		Pyrene, 4-methyl-	216	C17H12	003353-12-6	60



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF053117\
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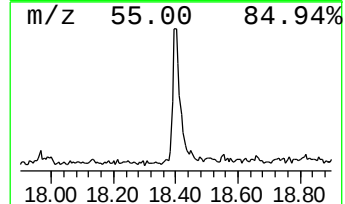
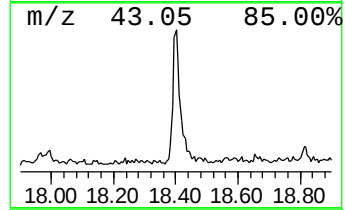
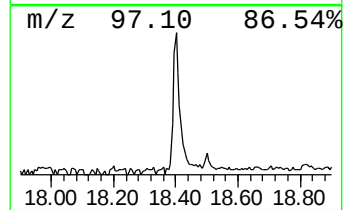
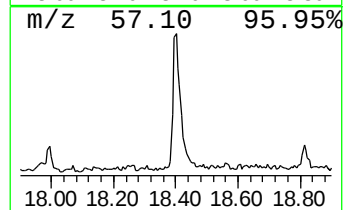
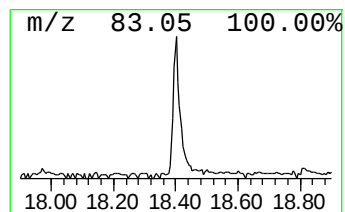
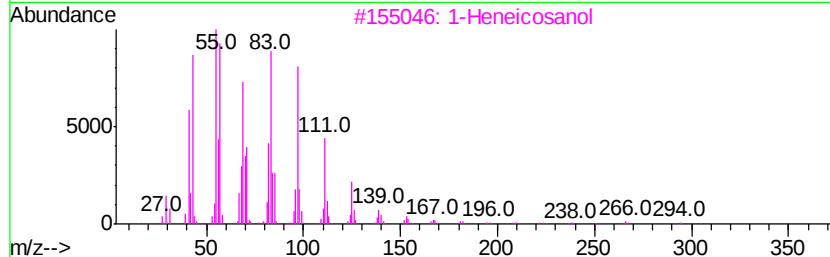
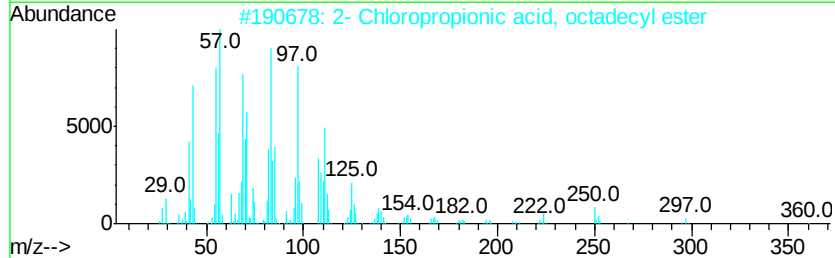
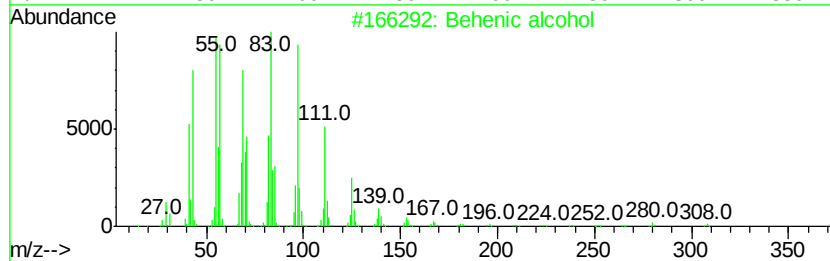
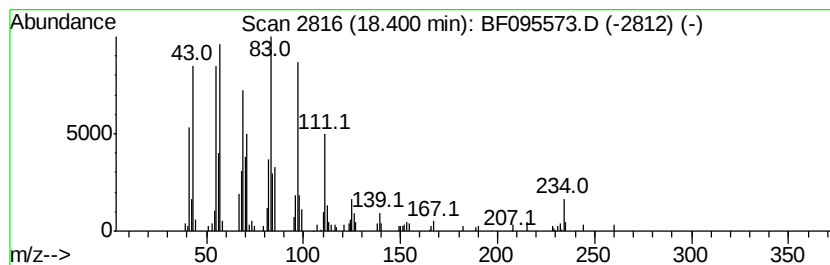
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Behenic alcohol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.40	4.59 ng	248348	Chrysene-d12	18.51	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Behenic alcohol	326	C22H46O	000661-19-8	94
2	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94
3	1-Heneicosanol	312	C21H44O	015594-90-8	93
4	n-Nonadecanol-1	284	C19H40O	001454-84-8	93
5	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF053117\
Data File : BF095573.D
Acq On : 31 May 2017 20:08
Operator : SJ/MA
Sample : I3329-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B-8(0-3)

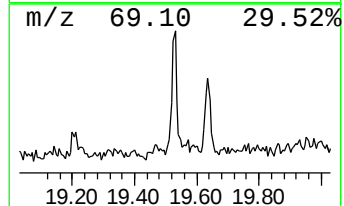
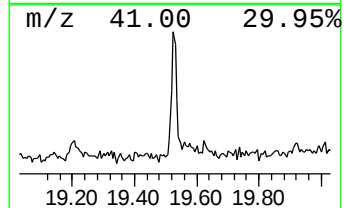
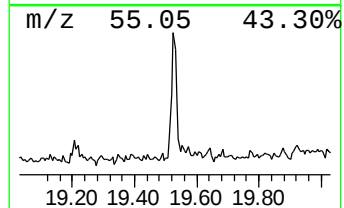
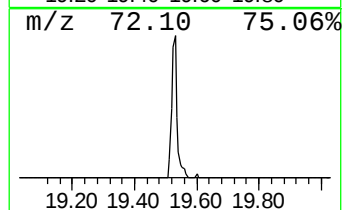
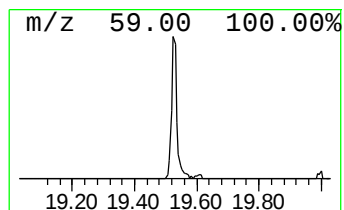
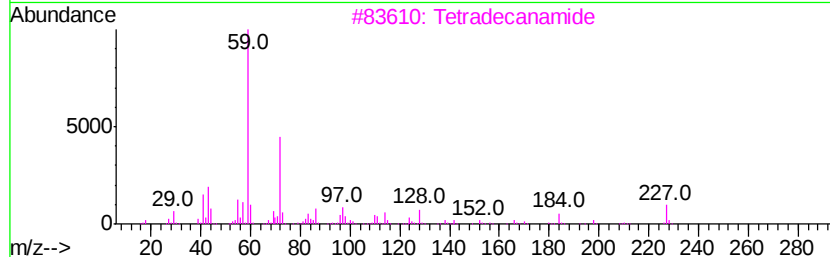
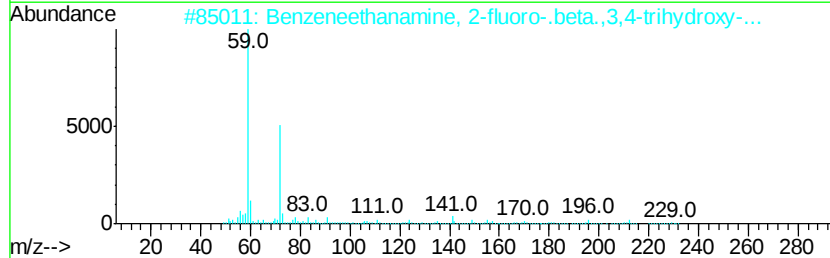
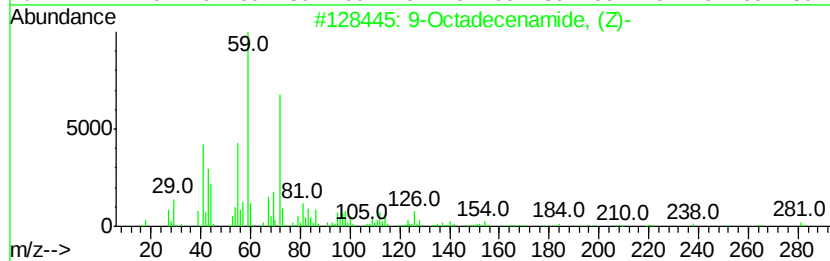
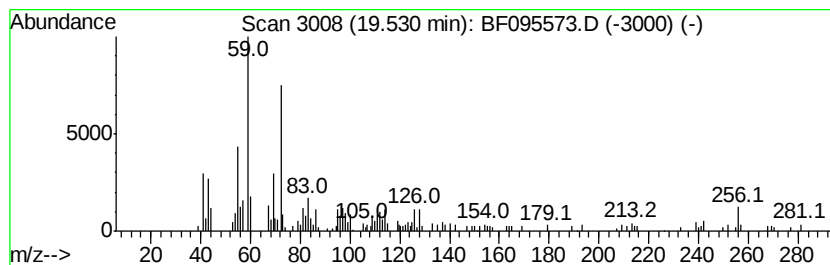
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 9-Octadecenamide, (Z)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.53	5.59 ng	158486	Perylene-d12	20.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	93
2		Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FN03	061338-98-5	50
3		Tetradecanamide	227	C14H29NO	000638-58-4	50
4		Hexadecanamide	255	C16H33NO	000629-54-9	47
5		1-(2-tert-Butoxycarbonylamino-3-...	314	C15H26N2O5	1000296-16-1	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF053117\
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Sample : I3329-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B-8(0-3)

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
2-Pentanone, 4-hy...	4.81	6.8 ng		215554	1	7.55	633498	20.0
unknown7.21	7.21	82.5 ng		2613850	1	7.55	633498	20.0
Anthracene, 2-met...	15.61	2.5 ng		124610	4	14.81	991354	20.0
Phenanthrene, 2-m...	15.65	3.3 ng		164658	4	14.81	991354	20.0
4H-Cyclopenta[def...	15.75	3.5 ng		174044	4	14.81	991354	20.0
n-Hexadecanoic acid	15.79	4.7 ng		231465	4	14.81	991354	20.0
Phenanthrene, 2,5...	16.41	2.5 ng		121850	4	14.81	991354	20.0
11H-Benzo[b]fluorene	17.24	2.0 ng		109475	5	18.51	1081880	20.0
Pyrene, 1-methyl-	17.51	2.9 ng		158689	5	18.51	1081880	20.0
Behenic alcohol	18.40	4.6 ng		248348	5	18.51	1081880	20.0
9-Octadecenamide,...	19.53	5.6 ng		158486	6	20.18	566731	20.0