

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052717\
 Data File : BF095517.D
 Acq On : 28 May 2017 2:06
 Operator : SJ/MA
 Sample : I3244-13
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM5-COMP-0-6

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	5.095	533	537	558	rBV	86500	221904	10.38%	1.052%
2	5.701	636	640	680	rBV	694388	1980375	92.66%	9.386%
3	7.283	905	909	925	rBV	658800	1569680	73.45%	7.439%
4	7.401	925	929	940	rVB	1225320	2137163	100.00%	10.129%
5	7.736	982	986	998	rBV	438048	544527	25.48%	2.581%
6	7.972	1021	1026	1045	rBV	1272519	1631083	76.32%	7.730%
7	8.642	1136	1140	1172	rBV	489362	1160170	54.29%	5.499%
8	9.766	1327	1331	1358	rBV	310659	722222	33.79%	3.423%
9	11.530	1626	1631	1656	rBV	1144965	2030167	94.99%	9.622%
10	12.242	1747	1752	1769	rBV	84107	170312	7.97%	0.807%
11	12.595	1807	1812	1819	rBV2	419035	698310	32.67%	3.310%
12	13.424	1947	1953	1958	rBV	20314	25041	1.17%	0.119%
13	13.518	1962	1969	1977	rVB2	13597	25101	1.17%	0.119%
14	13.895	2028	2033	2058	rBV	454190	934913	43.75%	4.431%
15	14.195	2081	2084	2088	rBV	17392	22350	1.05%	0.106%
16	14.677	2160	2166	2172	rBV	22606	31363	1.47%	0.149%
17	15.001	2216	2221	2225	rBV2	310257	467148	21.86%	2.214%
18	15.036	2225	2227	2240	rVV	142726	265507	12.42%	1.258%
19	15.142	2241	2245	2256	rVB2	21898	45058	2.11%	0.214%
20	15.483	2298	2303	2317	rVB5	36687	84200	3.94%	0.399%
21	15.842	2360	2364	2371	rVB3	17031	26553	1.24%	0.126%
22	15.954	2378	2383	2401	rVB3	70909	150613	7.05%	0.714%
23	16.783	2520	2524	2533	rBV	220281	299446	14.01%	1.419%
24	16.853	2533	2536	2542	rVB3	39933	48286	2.26%	0.229%
25	17.042	2563	2568	2571	rBV3	20275	29974	1.40%	0.142%
26	17.089	2571	2576	2583	rVV	222743	327269	15.31%	1.551%
27	17.312	2610	2614	2624	rVV	1029707	1239613	58.00%	5.875%
28	17.406	2627	2630	2636	rVB3	16884	27073	1.27%	0.128%
29	17.553	2652	2655	2663	rVB	27278	45792	2.14%	0.217%
30	17.642	2666	2670	2674	rBV	23668	31809	1.49%	0.151%
31	17.695	2674	2679	2684	rBV3	25138	43142	2.02%	0.204%
32	17.877	2707	2710	2717	rVB	60245	63716	2.98%	0.302%
33	17.936	2717	2720	2724	rBV5	16660	24973	1.17%	0.118%
34	18.012	2729	2733	2741	rVB2	25556	32164	1.50%	0.152%

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35	18.436	2800	2805	2807	rBV2	54558	67163	3.14%	0.318%
36	18.553	2816	2825	2835	rBV5	48685	134598	6.30%	0.638%
37	18.671	2840	2845	2849	rVV2	500164	741388	34.69%	3.514%
38	18.706	2849	2851	2863	rVB3	135312	236156	11.05%	1.119%
39	18.806	2863	2868	2870	rBV2	29722	40364	1.89%	0.191%
40	18.842	2870	2874	2878	rVB	85497	94228	4.41%	0.447%
41	18.895	2878	2883	2886	rBV	129361	140763	6.59%	0.667%
42	18.924	2886	2888	2891	rVV	44874	46198	2.16%	0.219%
43	18.965	2891	2895	2900	rVB3	62067	75377	3.53%	0.357%
44	19.189	2927	2933	2936	rBV3	22784	31901	1.49%	0.151%
45	19.336	2955	2958	2962	rBV4	44526	58218	2.72%	0.276%
46	19.365	2962	2963	2970	rVB7	20113	23892	1.12%	0.113%
47	19.783	3028	3034	3041	rBV	99319	136511	6.39%	0.647%
48	19.947	3058	3062	3065	rBV	126113	166732	7.80%	0.790%
49	19.977	3065	3067	3074	rVV2	51223	76029	3.56%	0.360%
50	20.065	3078	3082	3087	rBV	88195	99426	4.65%	0.471%
51	20.241	3108	3112	3118	rBV	82065	95190	4.45%	0.451%
52	20.300	3118	3122	3126	rBV	112981	138390	6.48%	0.656%
53	20.353	3126	3131	3140	rBV	332530	508745	23.80%	2.411%
54	20.583	3167	3170	3177	rVB6	34995	50738	2.37%	0.240%
55	20.788	3200	3205	3209	rBV	130782	180563	8.45%	0.856%
56	21.112	3256	3260	3267	rBV8	20289	40439	1.89%	0.192%
57	21.459	3315	3319	3325	rVB7	23391	40485	1.89%	0.192%
58	21.718	3353	3363	3369	rBV4	171729	352597	16.50%	1.671%
59	21.865	3383	3388	3393	rVB4	72936	129321	6.05%	0.613%
60	22.106	3421	3429	3440	rBV4	38973	114354	5.35%	0.542%
61	22.365	3467	3473	3474	rBV5	12661	24872	1.16%	0.118%
62	22.965	3568	3575	3584	rVB9	36653	98098	4.59%	0.465%

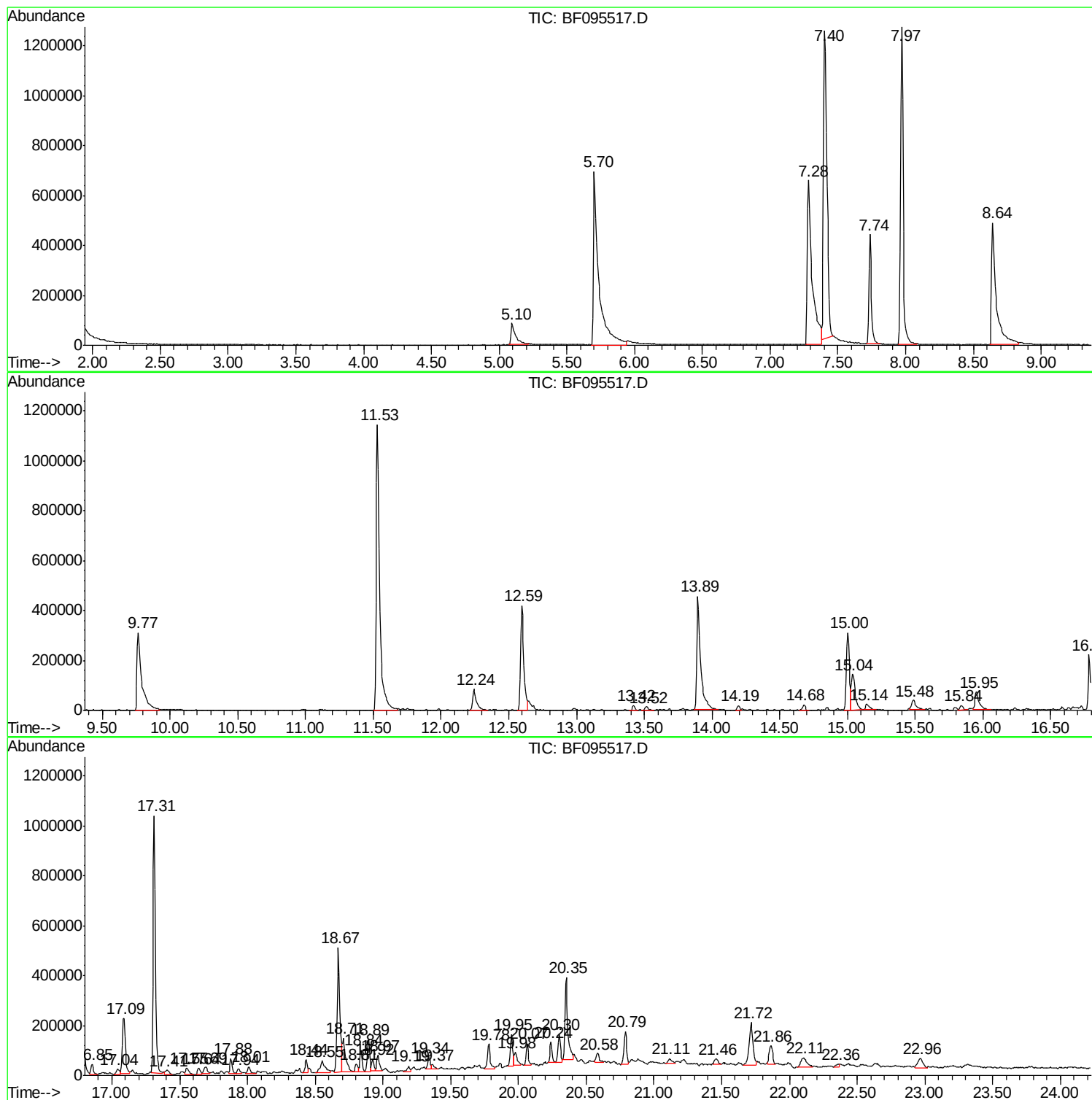
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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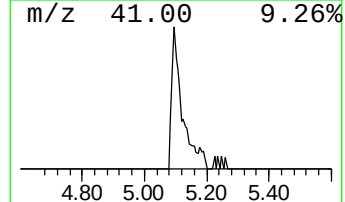
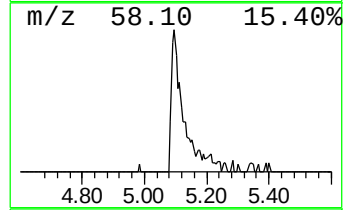
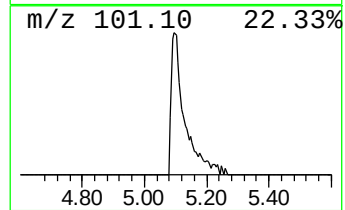
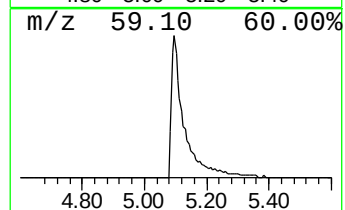
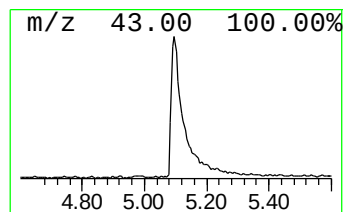
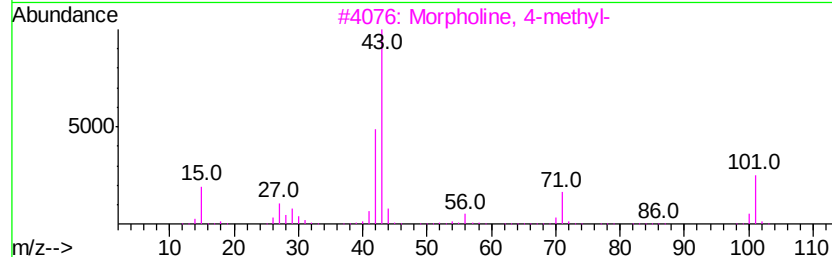
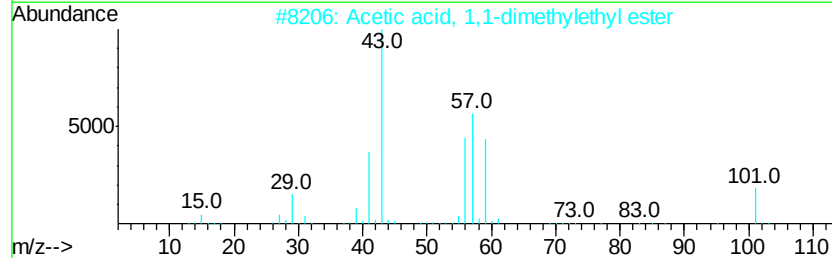
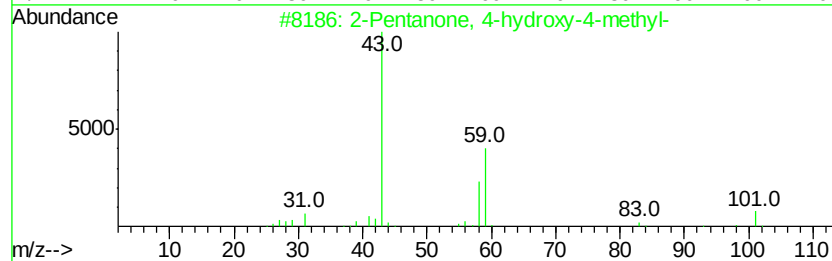
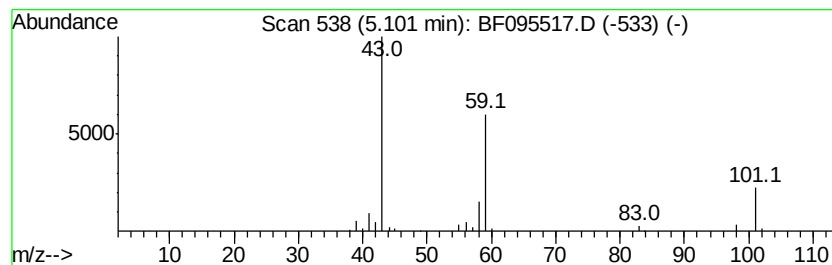
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	8.15 ng	221904	1,4-Dichlorobenzene-d4	7.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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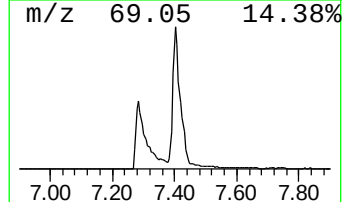
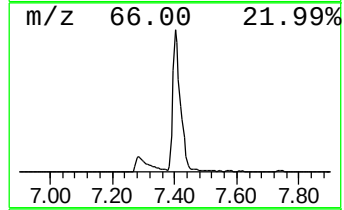
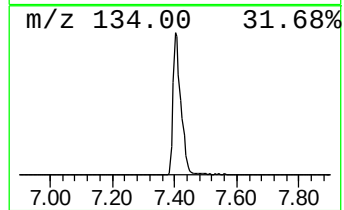
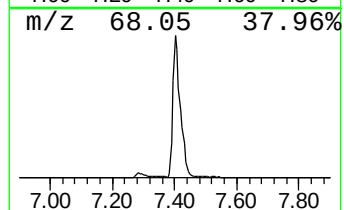
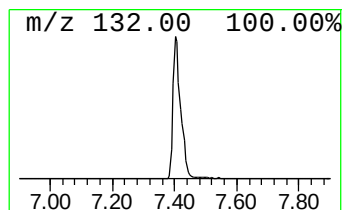
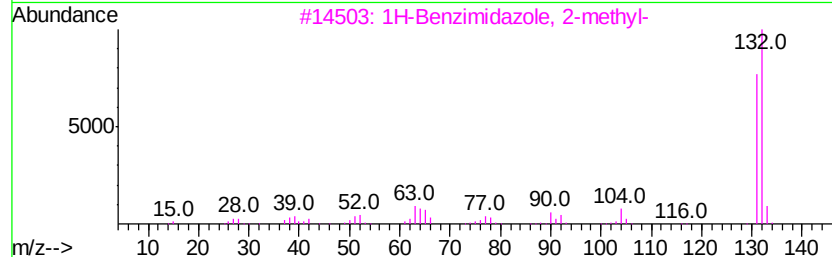
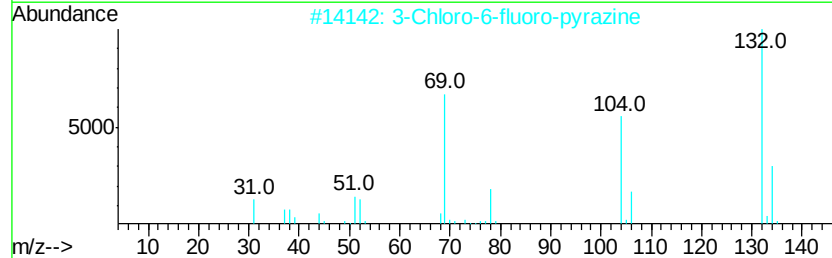
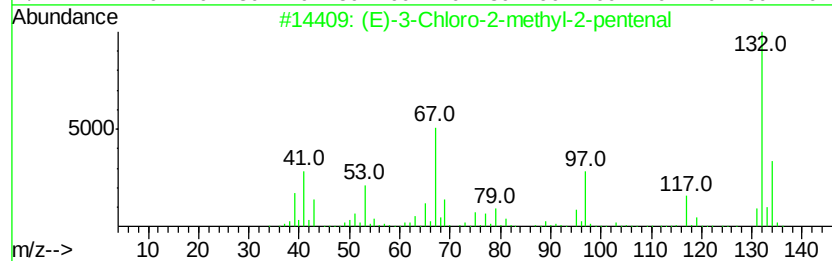
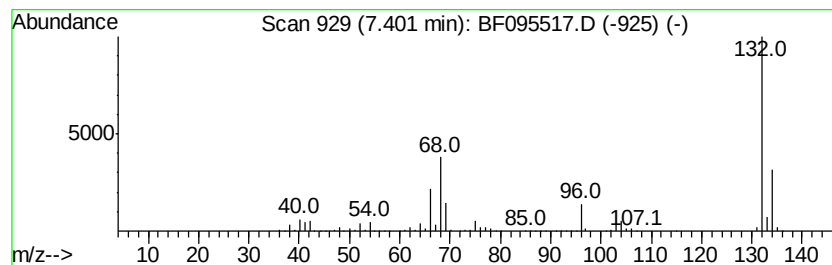
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown7.40 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.40	78.50 ng	2137160	1,4-Dichlorobenzene-d4	7.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
4		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	10
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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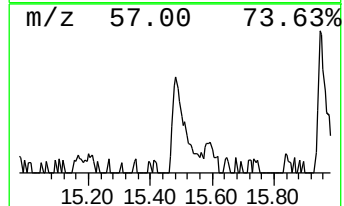
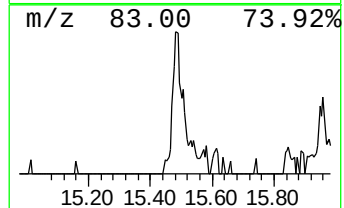
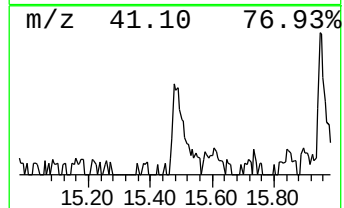
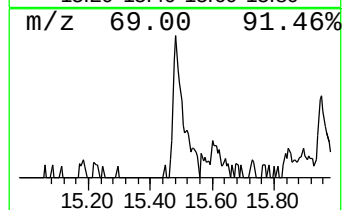
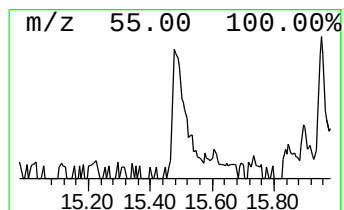
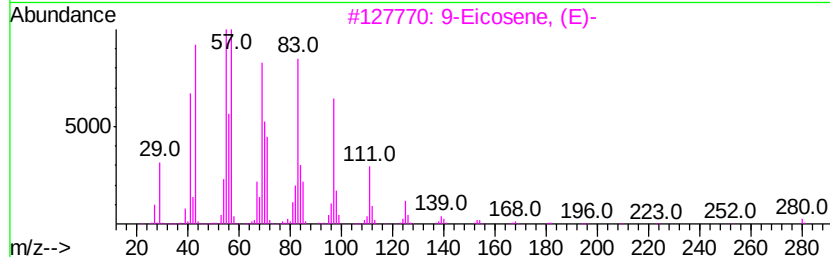
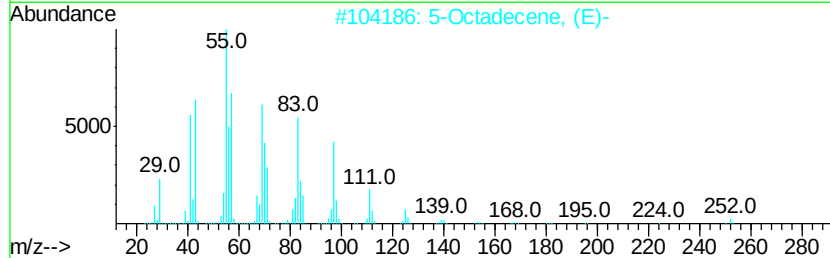
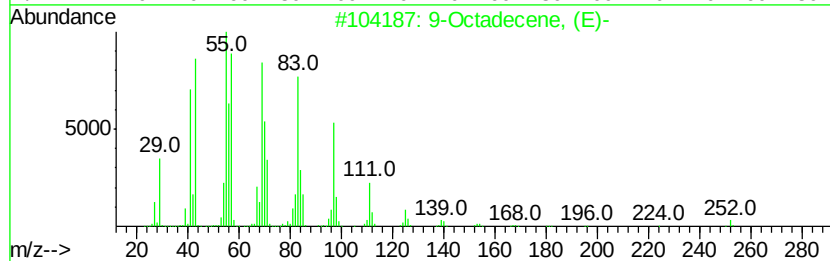
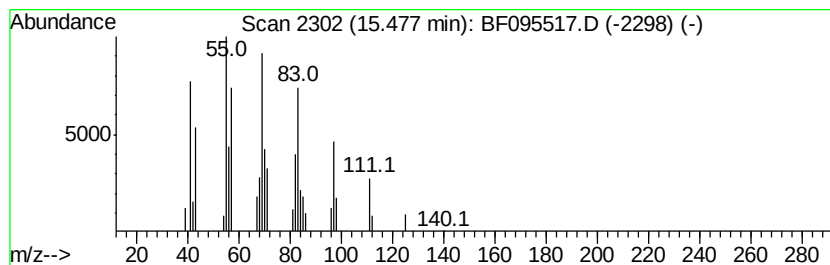
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 9-Octadecene, (E)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.48	3.60 ng	84200	Phenanthrene-d10	15.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecene, (E)-			252	C18H36	007206-25-9	87
2	5-Octadecene, (E)-			252	C18H36	007206-21-5	83
3	9-Eicosene, (E)-			280	C20H40	074685-29-3	70
4	Acetic acid, trifluoro-, undecyl...			268	C13H23F3O2	053800-01-4	68
5	4-Heptafluorobutyryloxyhexadecane			438	C20H33F7O2	1000282-97-2	62



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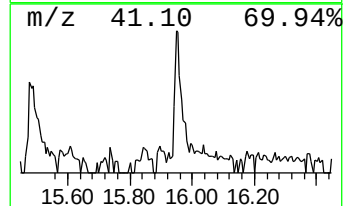
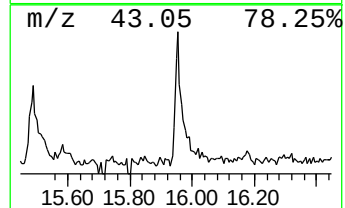
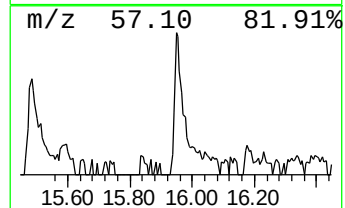
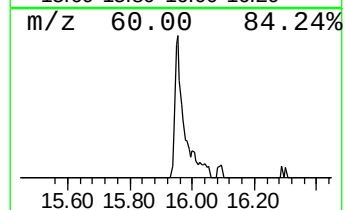
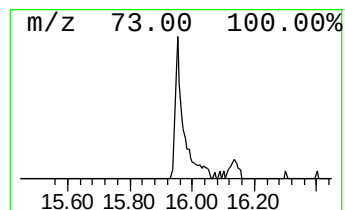
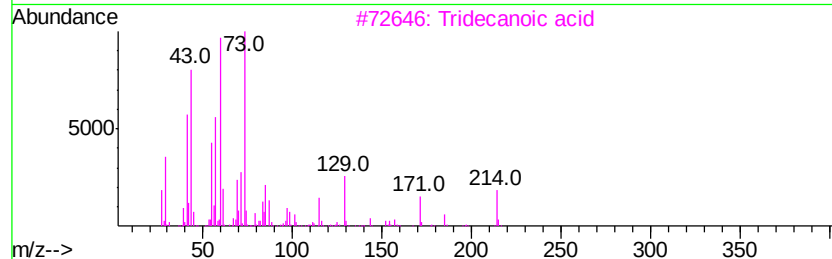
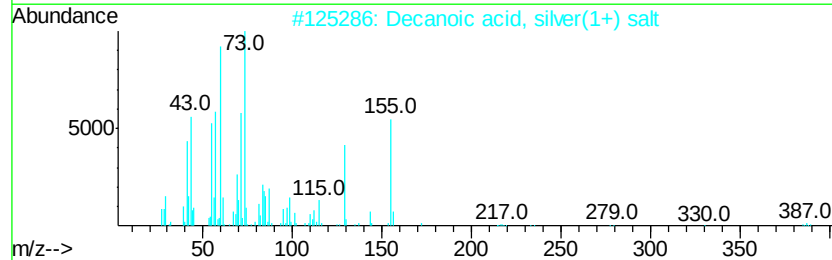
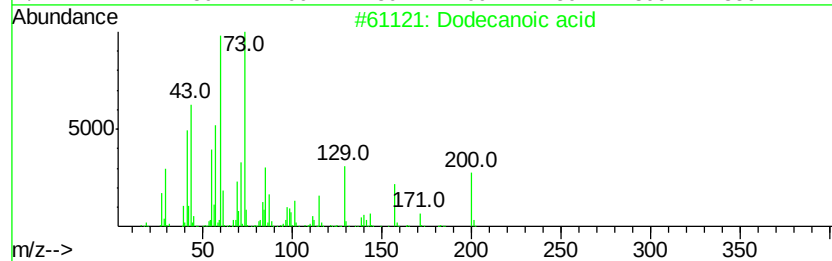
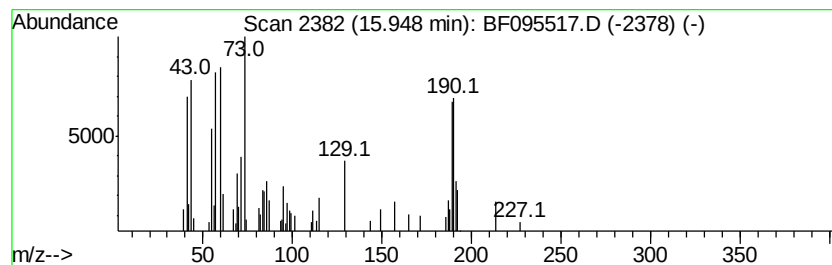
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown15.95 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.95	6.45 ng	150613	Phenanthrene-d10	15.00	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecanoic acid	200	C12H24O2	000143-07-7	49
2	Decanoic acid, silver(1+) salt	278	C10H19AgO2	013126-67-5	38
3	Tridecanoic acid	214	C13H26O2	000638-53-9	38
4	Trehalose	342	C12H22O11	000099-20-7	35
5	.alpha.-D-Glucopyranose, 4-O-.be...	342	C12H22O11	014641-93-1	35



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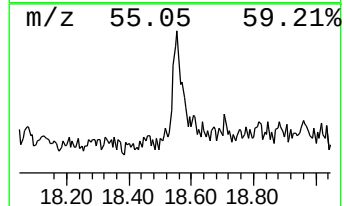
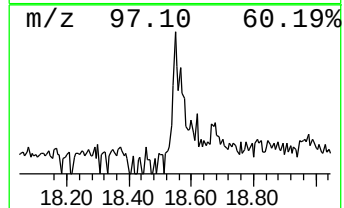
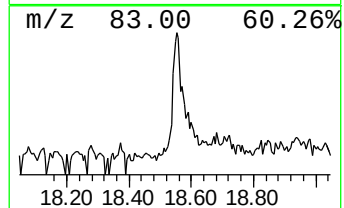
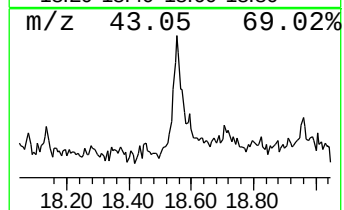
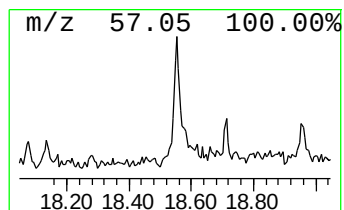
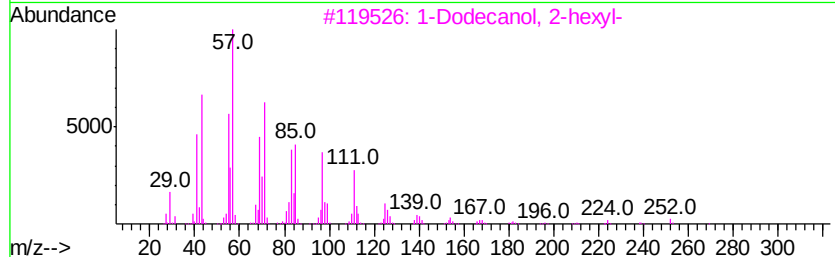
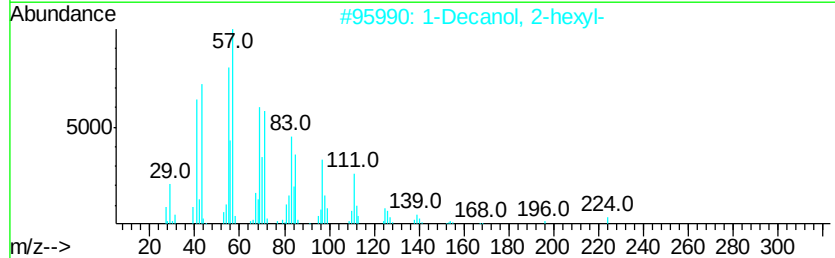
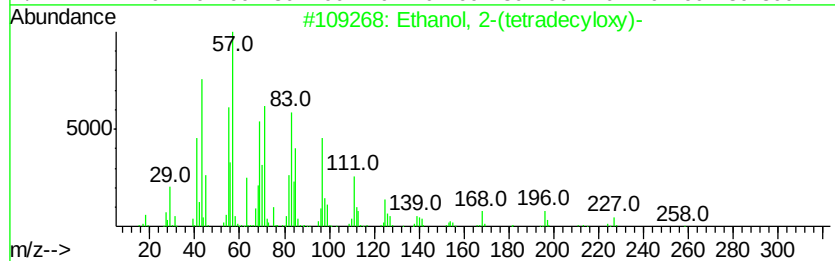
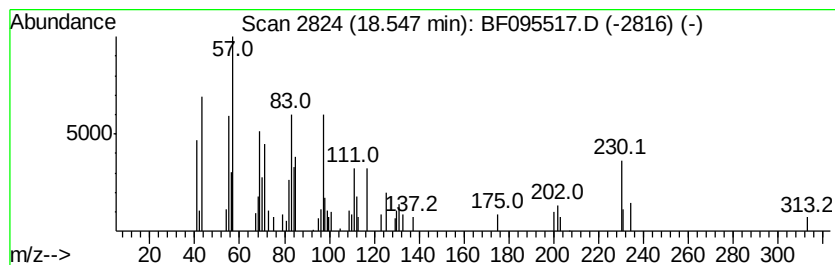
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Ethanol, 2-(tetradecyloxy)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.			
18.55	3.63 ng	134598	Chrysene-d12	18.67			
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	76
2			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	64
3			1-Dodecanol, 2-hexyl-	270	C18H38O	110225-00-8	49
4			1-Heptadecene	238	C17H34	006765-39-5	49
5			Carbonic acid, tridecyl 2,2,2-tr...	374	C16H29Cl3O3	1000314-55-9	49



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052717\
Data File : BF095517.D
Acq On : 28 May 2017 2:06
Operator : SJ/MA
Sample : I3244-13
Misc :
ALS Vial : 16 Sample Multiplier: 1

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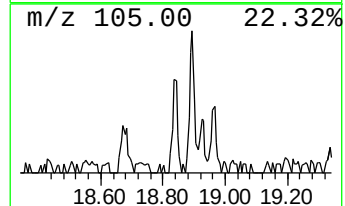
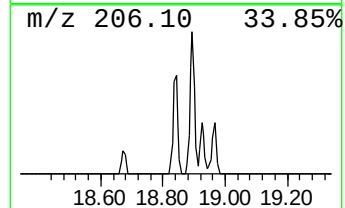
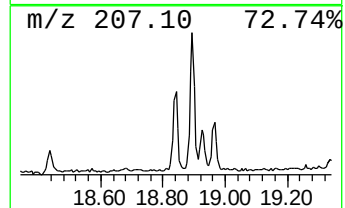
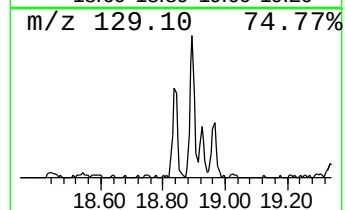
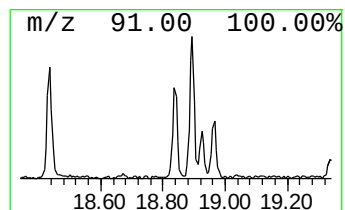
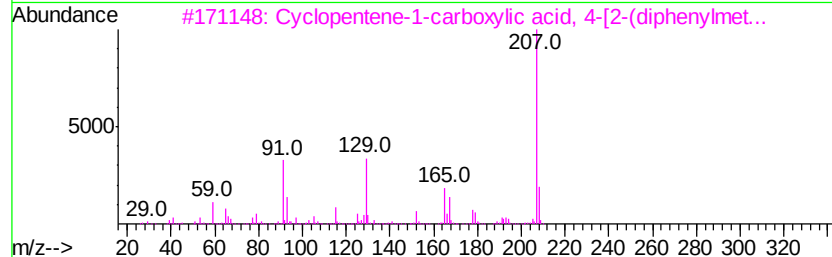
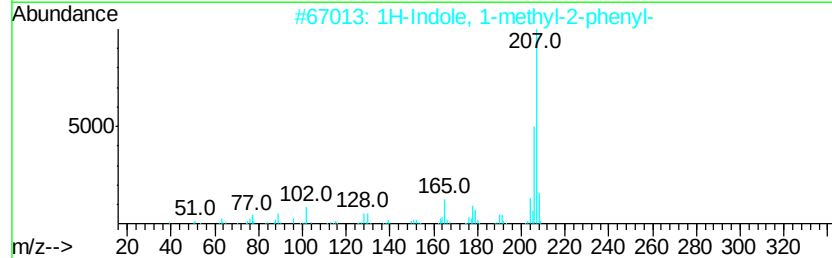
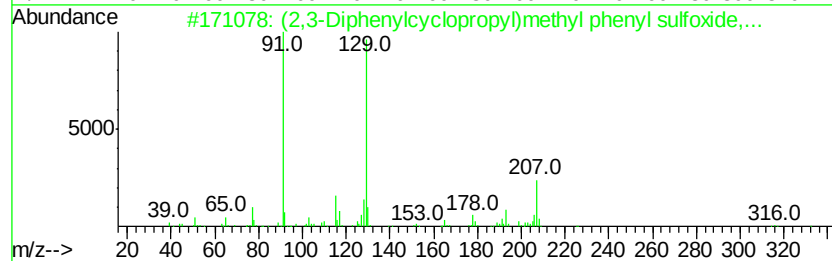
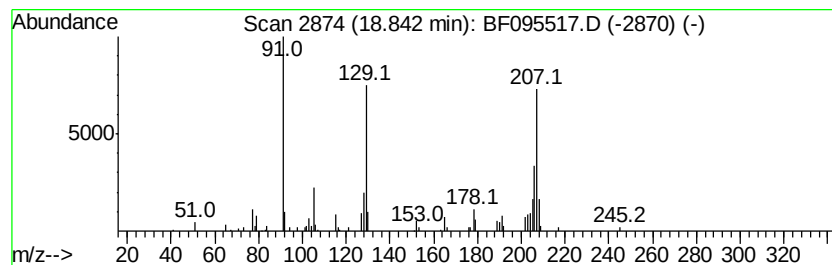
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 (2,3-Diphenylcyclopropyl)me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.84	2.54 ng	94228	Chrysene-d12	18.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(2,3-Diphenylcyclopropyl)methyl ...	332	C22H20OS	131758-71-9	56
2		1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	46
3		Cyclopentene-1-carboxylic acid, ...	332	C23H24O2	1000159-40-6	42
4		1H-Indole, 5-methyl-2-phenyl-	207	C15H13N	013228-36-9	41
5		Quinoline	129	C9H7N	000091-22-5	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052717\
Data File : BF095517.D
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Instrument :
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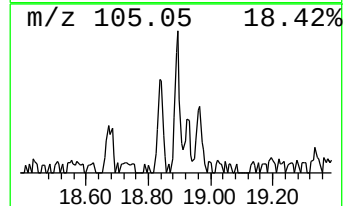
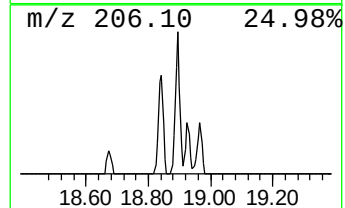
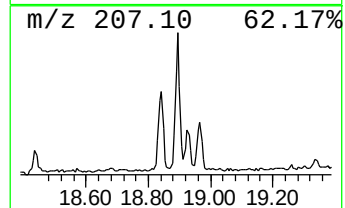
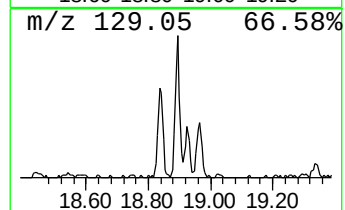
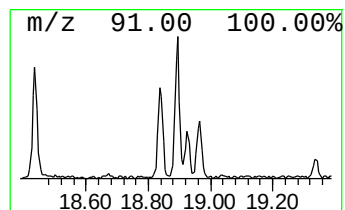
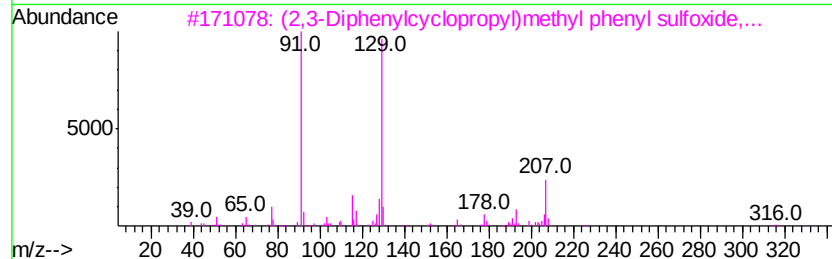
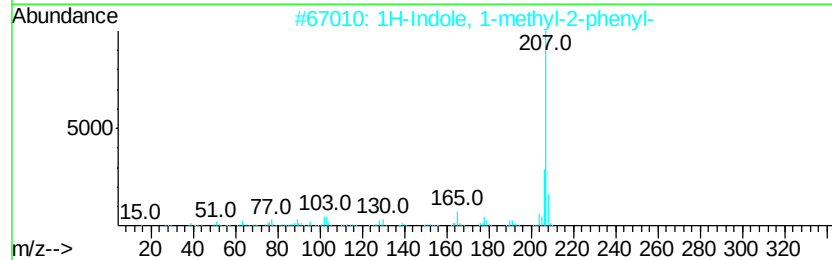
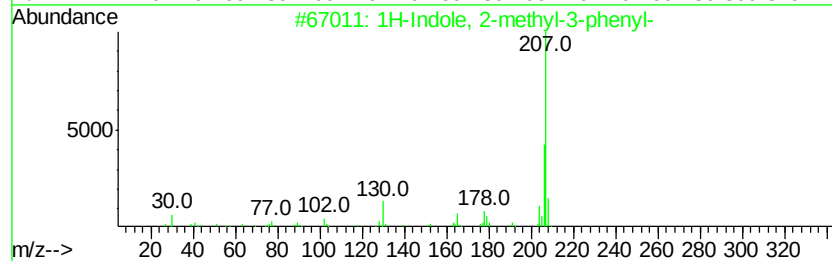
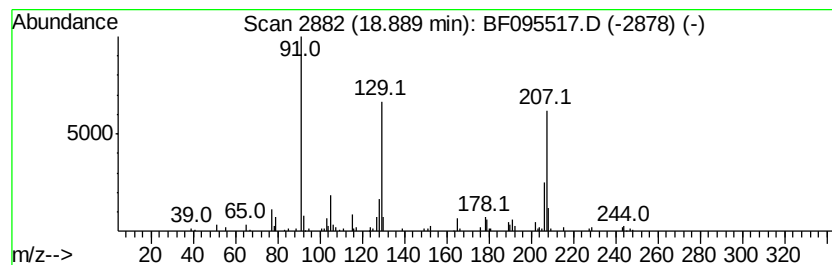
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown18.89 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.89	3.80 ng	140763	Chrysene-d12	18.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	45
2		1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	41
3		(2,3-Diphenylcyclopropyl)methyl ...	332	C22H20OS	131758-71-9	37
4		1-benzylindole	207	C15H13N	1000334-31-3	35
5		Indolizine, 2-(4-methylphenyl)-	207	C15H13N	007496-81-3	30



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052717\
Data File : BF095517.D
Acq On : 28 May 2017 2:06
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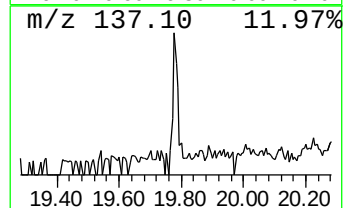
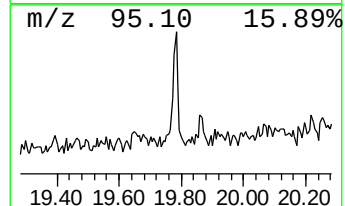
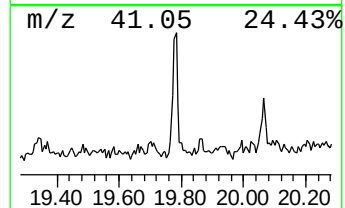
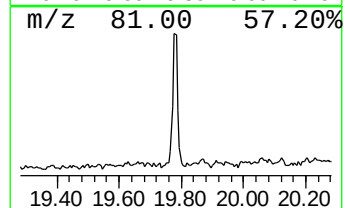
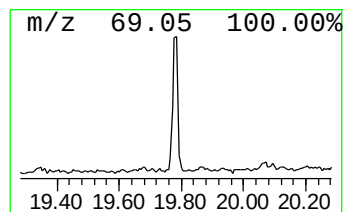
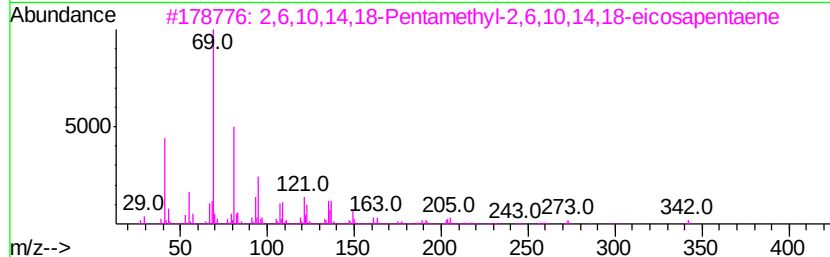
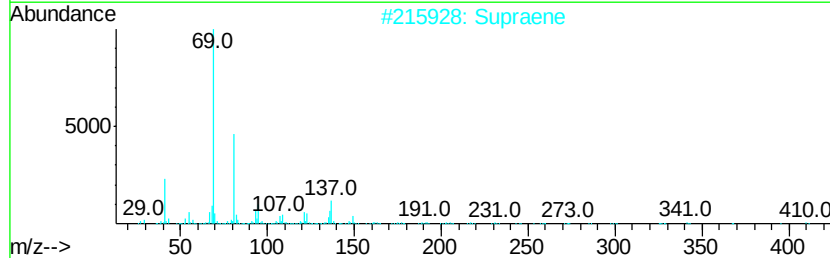
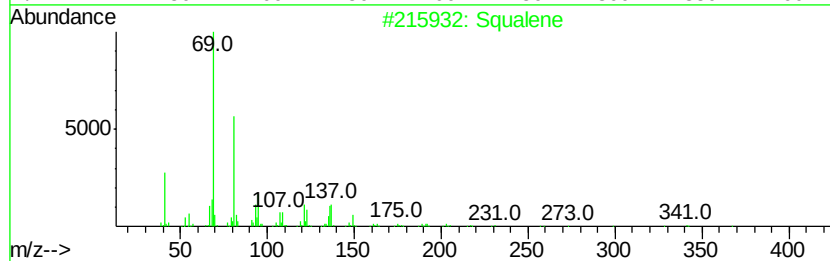
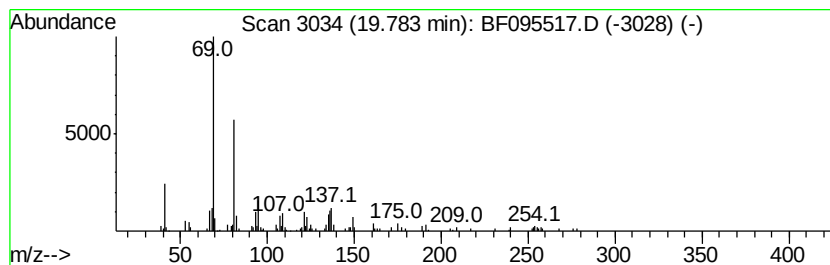
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Squalene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.78	5.37 ng	136511	Perylene-d12	20.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	87
2		Supraene	410	C30H50	007683-64-9	87
3		2,6,10,14,18-Pentamethyl-2,6,10,...	342	C25H42	075581-03-2	86
4		Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	86
5		Hexadeca-2,6,10,14-tetraen-1-ol,...	290	C20H34O	007614-21-3	78



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052717\
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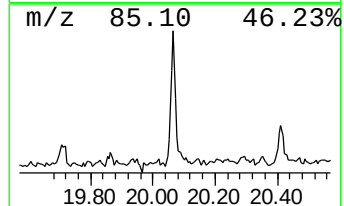
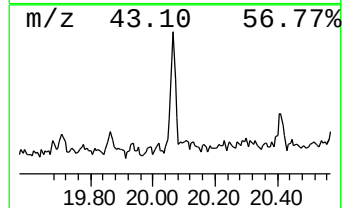
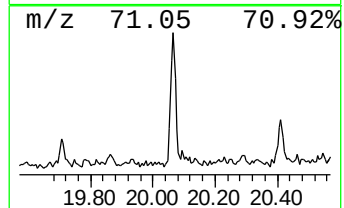
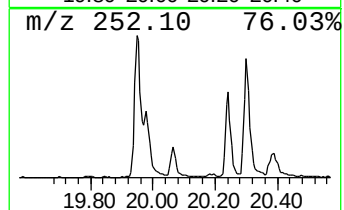
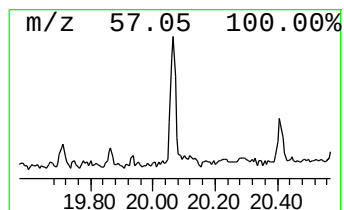
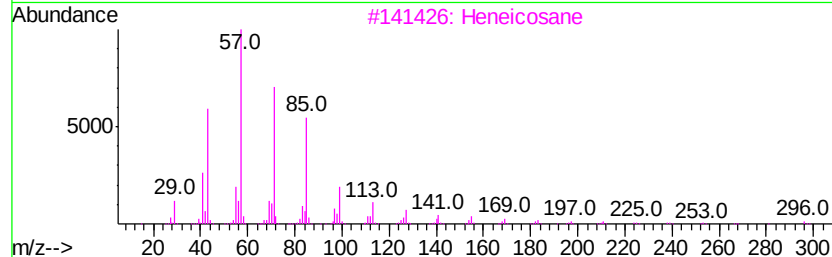
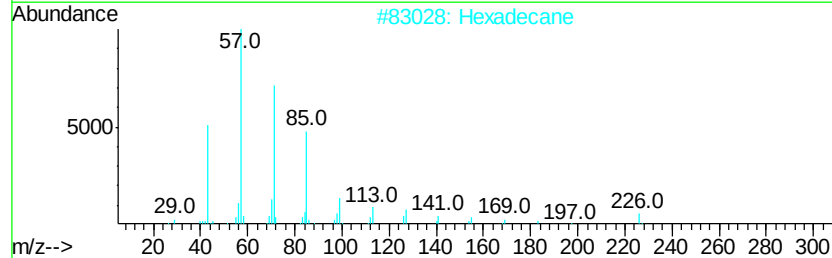
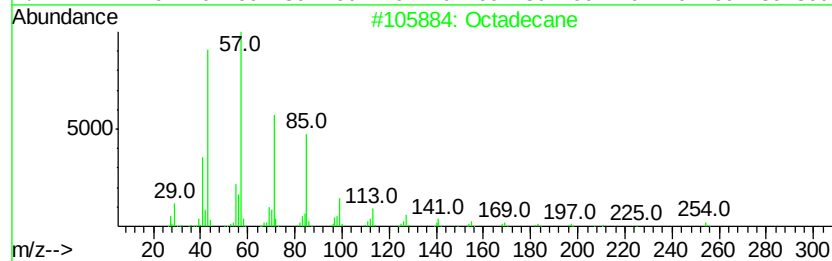
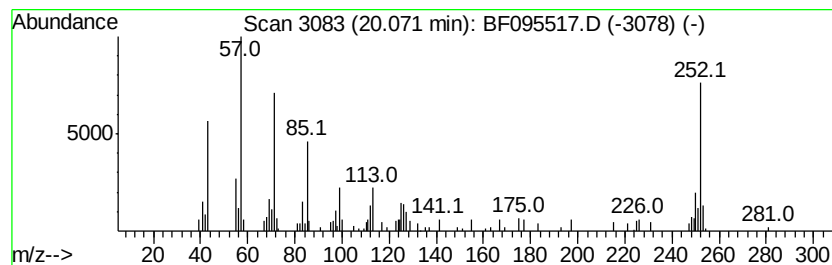
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 unknown20.07 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.07	3.91 ng	99426	Perylene-d12	20.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	45
2		Hexadecane	226	C16H34	000544-76-3	45
3		Heneicosane	296	C21H44	000629-94-7	45
4		Benzo[k]fluoranthene	252	C20H12	000207-08-9	45
5		Triacotane	422	C30H62	000638-68-6	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052717\
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Acq On : 28 May 2017 2:06
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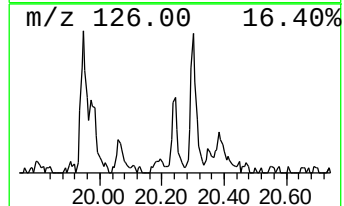
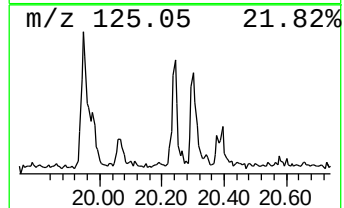
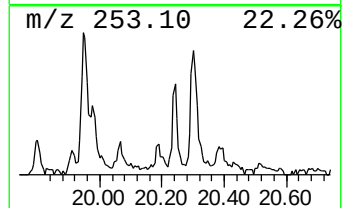
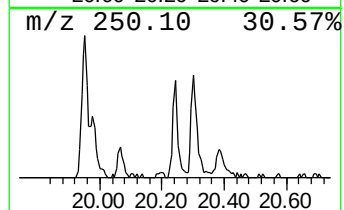
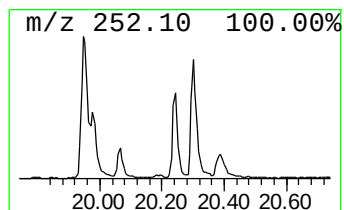
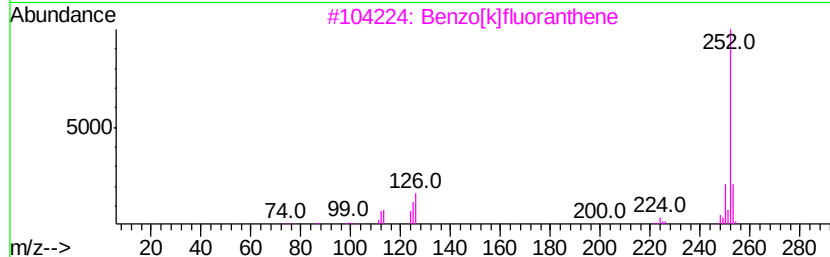
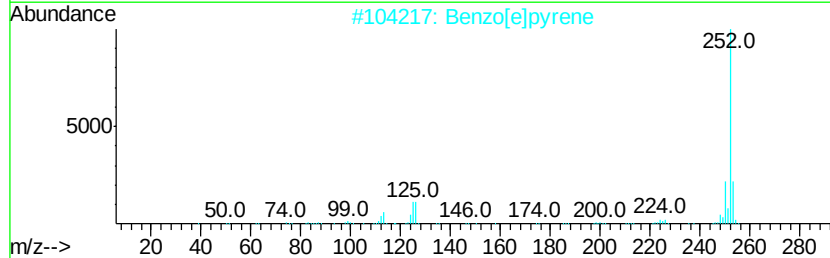
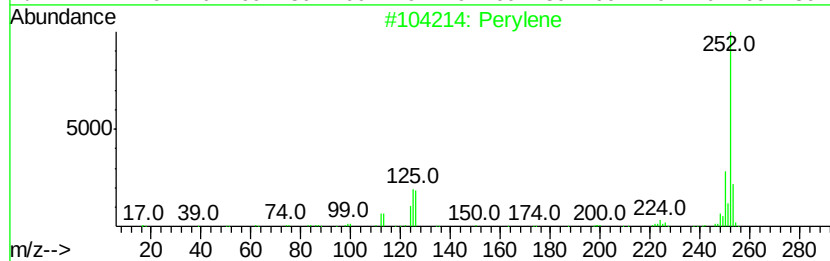
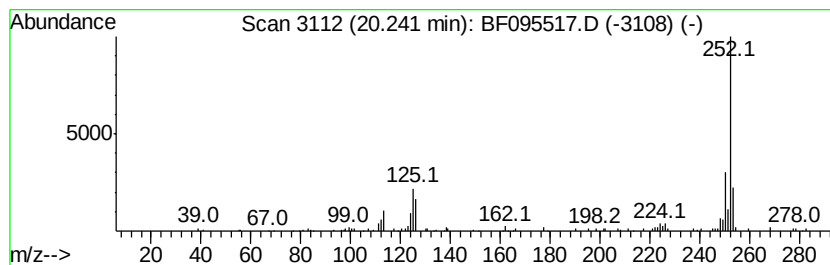
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Perylene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.24	3.74 ng	95190	Perylene-d12	20.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Perylene	252	C20H12	000198-55-0	99
2			Benzo[e]pyrene	252	C20H12	000192-97-2	97
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
4			Benzo[i]fluoranthene	252	C20H12	000205-82-3	94
5			Benz[e]acephenanthrylene	252	C20H12	000205-99-2	94



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052717\
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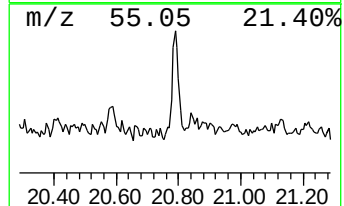
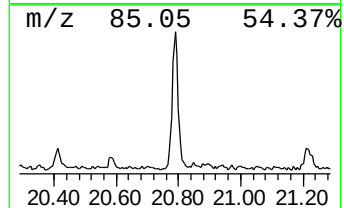
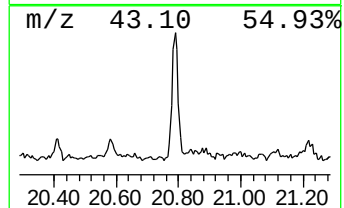
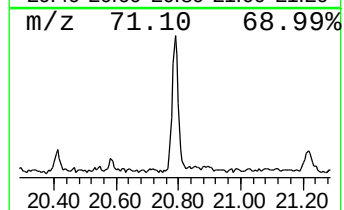
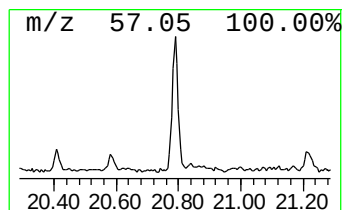
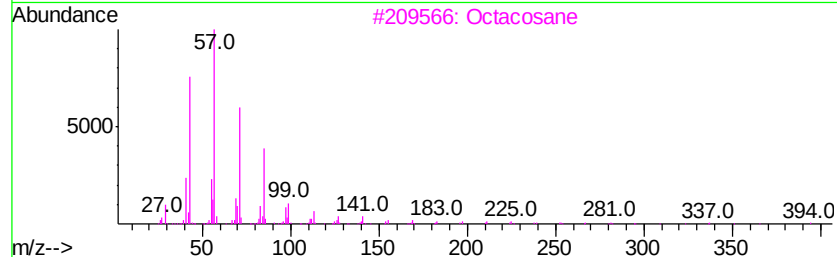
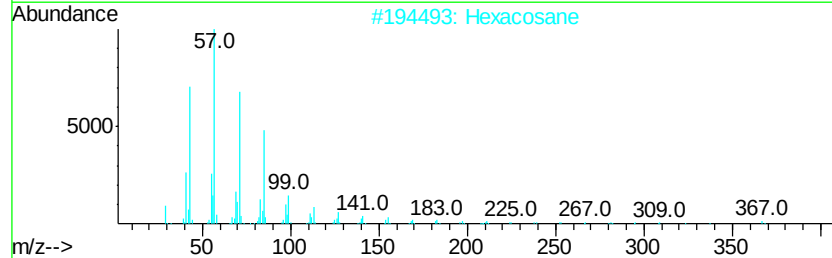
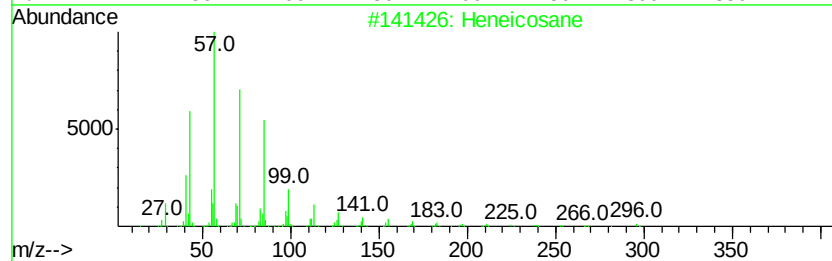
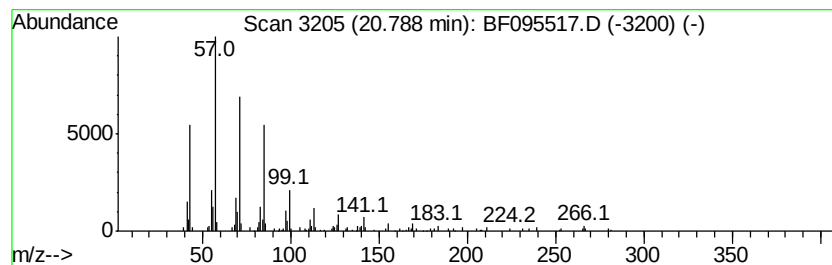
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heneicosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.79	7.10 ng	180563	Perylene-d12	20.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	90
2		Hexacosane	366	C26H54	000630-01-3	87
3		Octacosane	394	C28H58	000630-02-4	87
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	86
5		Tetracosane	338	C24H50	000646-31-1	83



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Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
NM5-COMP-0-6

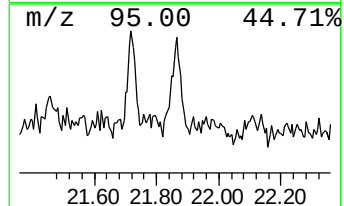
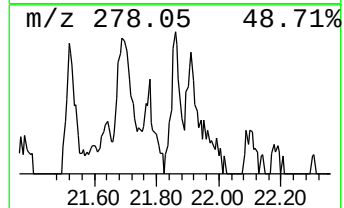
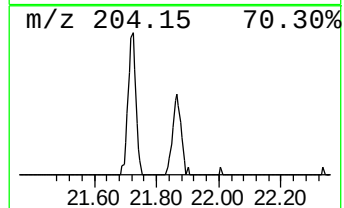
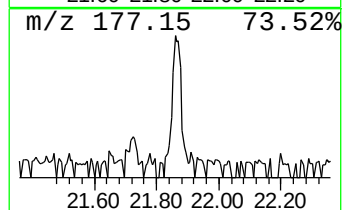
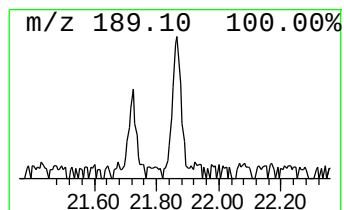
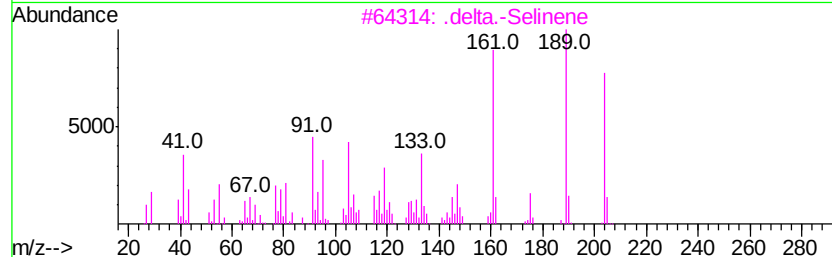
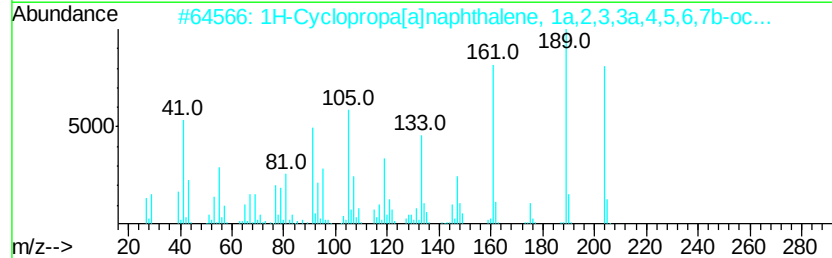
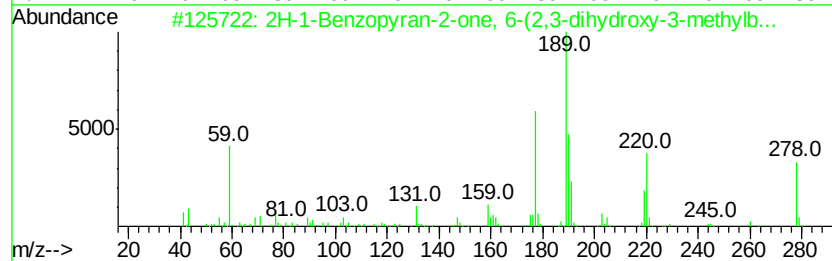
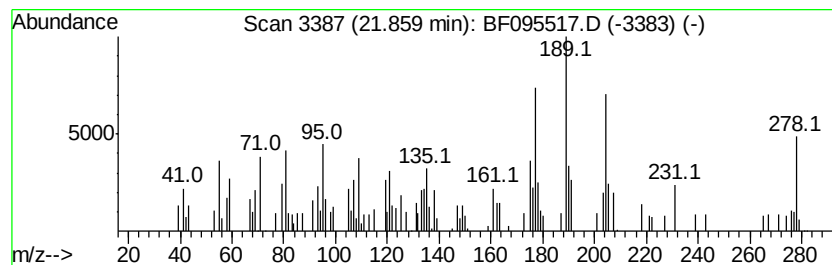
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown21.86 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.86	5.08 ng	129321	Perylene-d12	20.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H-1-Benzopyran-2-one, 6-(2,3-di...	278	C15H18O5	036149-96-9	46
2		1H-Cyclopropa[<i>a</i>]naphthalene, 1a,...	204	C15H24	000489-29-2	35
3		.delta.-Selinene	204	C15H24	028624-23-9	35
4		1,3a-Ethano-3aH-indene, 1,2,3,6,...	204	C15H24	004545-68-0	27
5		2-Acetyl-3,5-dimethylbenzo(b)thi...	204	C12H12OS	006179-05-1	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052717\
Data File : BF095517.D
Acq On : 28 May 2017 2:06
Operator : SJ/MA
Sample : I3244-13
Misc :
ALS Vial : 16 Sample Multiplier: 1

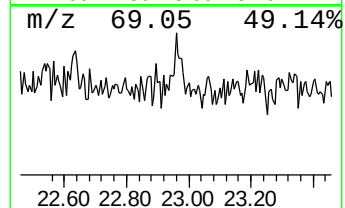
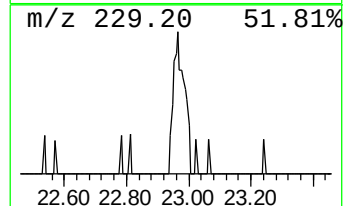
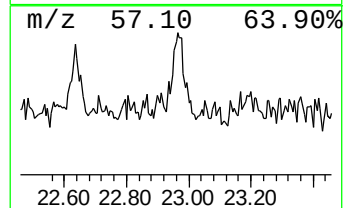
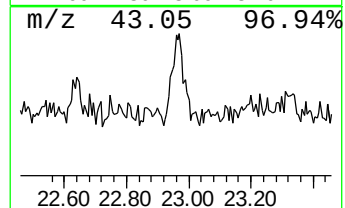
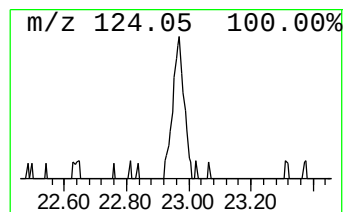
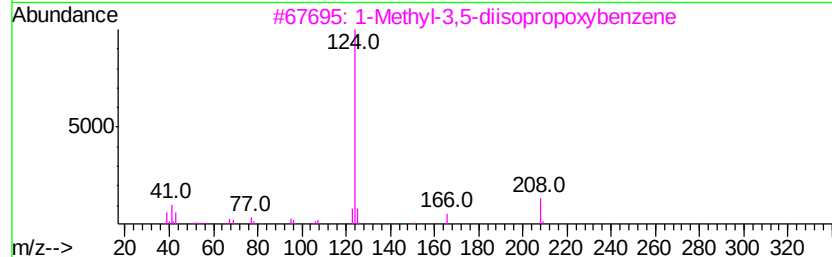
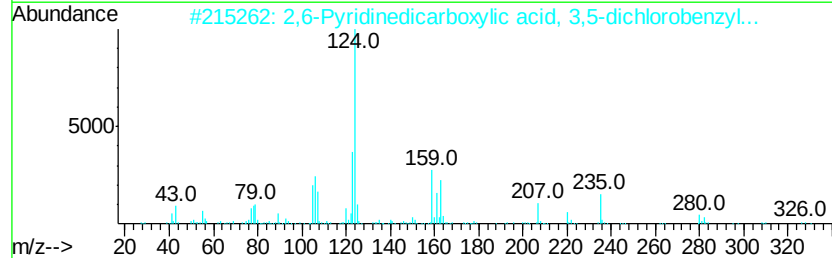
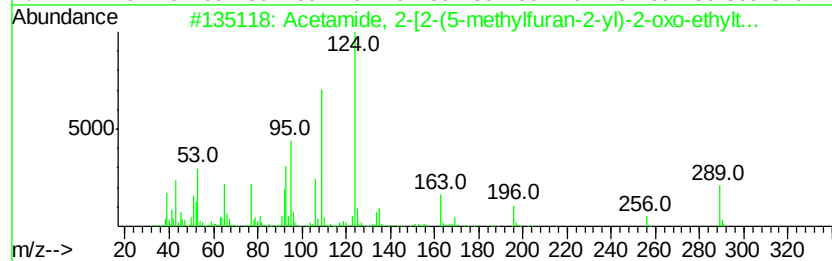
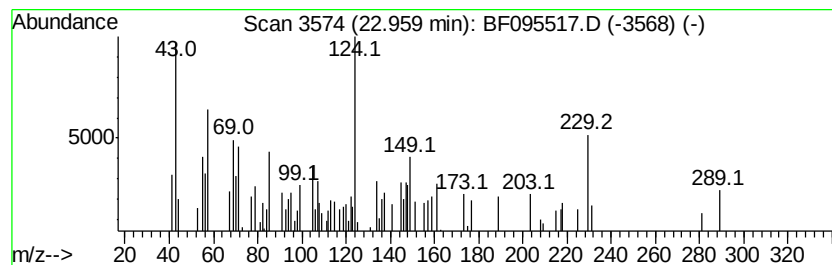
Instrument :
BNA_F
ClientSampleId :
NM5-COMP-0-6

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 15 unknown22.96 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.96	3.86 ng	98098	Perylene-d12	20.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Acetamide, 2-[2-(5-methylfuran-2-yl)-2-oxo-ethyl]...	289 C15H15N03S	1000259-79-3 27
2		2,6-Pyridinedicarboxylic acid, 3,5-dichlorobenzyl...	409 C20H21Cl2N04	1000369-10-5 22
3		1-Methyl-3,5-diisopropoxybenzene	208 C13H20O2	1000281-22-2 22
4		2,6-Pyridinedicarboxylic acid, 3,5-dichlorobenzyl...	423 C21H23Cl2N04	1000369-10-6 16
5		Trisiloxane, 1,3,5-triethenyl-1,...	272 C11H24O2Si3	001529-65-3 14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052717\
 Data File : BF095517.D
 Acq On : 28 May 2017 2:06
 Operator : SJ/MA
 Sample : I3244-13
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM5-COMP-0-6

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...	5.10	8.2	ng	221904	1	7.74	544527	20.0
unknown7.40	7.40	78.5	ng	2137160	1	7.74	544527	20.0
9-Octadecene, (E)-	15.48	3.6	ng	84200	4	15.00	467148	20.0
unknown15.95	15.95	6.5	ng	150613	4	15.00	467148	20.0
Ethanol, 2-(tetra...	18.55	3.6	ng	134598	5	18.67	741388	20.0
(2,3-Diphenylcycl...	18.84	2.5	ng	94228	5	18.67	741388	20.0
unknown18.89	18.89	3.8	ng	140763	5	18.67	741388	20.0
Squalene	19.78	5.4	ng	136511	6	20.35	508745	20.0
unknown20.07	20.07	3.9	ng	99426	6	20.35	508745	20.0
Perylene	20.24	3.7	ng	95190	6	20.35	508745	20.0
Heneicosane	20.79	7.1	ng	180563	6	20.35	508745	20.0
unknown21.86	21.86	5.1	ng	129321	6	20.35	508745	20.0
unknown22.96	22.96	3.9	ng	98098	6	20.35	508745	20.0