

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050416\
 Data File : BF086694.D
 Acq On : 4 May 2016 23:31
 Operator : UM/SJ
 Sample : H2715-12
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 26WARD-4-CS-1(0-8FTBGS)

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-BF050416.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	4.921	246	248	254	rBV	118977	225757	4.28%	0.349%
2	5.333	281	284	287	rBV	4484490	4854400	92.00%	7.509%
3	6.384	373	376	379	rBV	4147690	5043832	95.60%	7.803%
4	6.510	384	387	389	rVV	5259191	5276237	100.00%	8.162%
5	6.578	391	393	398	rVB2	67259	97316	1.84%	0.151%
6	6.738	405	407	409	rBV	1185688	1113339	21.10%	1.722%
7	6.819	411	414	416	rVB2	70645	82330	1.56%	0.127%
8	6.899	418	421	423	rBV	3814791	3991772	75.66%	6.175%
9	7.310	454	457	459	rBV	3235596	3534628	66.99%	5.468%
10	7.344	459	460	463	rVB	60671	74336	1.41%	0.115%
11	7.539	472	477	484	rVB6	18680	90043	1.71%	0.139%
12	8.019	517	519	525	rBV2	870893	1379568	26.15%	2.134%
13	8.739	579	582	584	rBV	122184	114268	2.17%	0.177%
14	8.830	588	590	593	rBV	73342	91145	1.73%	0.141%
15	9.104	611	614	616	rBV	4772302	4465041	84.63%	6.907%
16	9.196	619	622	626	rVB2	41693	98535	1.87%	0.152%
17	9.367	634	637	641	rBV2	64887	111045	2.10%	0.172%
18	9.436	641	643	644	rBV	120882	120921	2.29%	0.187%
19	9.493	647	648	651	rVV	303151	383457	7.27%	0.593%
20	9.550	651	653	656	rVB	52297	76987	1.46%	0.119%
21	9.630	658	660	662	rBV	61687	88015	1.67%	0.136%
22	9.722	666	668	669	rBV	81215	83067	1.57%	0.128%
23	9.779	670	673	674	rBV	1119118	1215809	23.04%	1.881%
24	9.802	674	675	678	rVB	245504	325446	6.17%	0.503%
25	9.927	684	686	689	rBV2	38715	79045	1.50%	0.122%
26	9.985	689	691	692	rVV	154516	225762	4.28%	0.349%
27	10.019	692	694	698	rVB2	47956	79671	1.51%	0.123%
28	10.213	709	711	713	rVB	128625	150229	2.85%	0.232%
29	10.316	719	720	723	rVB	295354	333851	6.33%	0.516%
30	10.385	723	726	727	rBV	54389	74440	1.41%	0.115%
31	10.430	729	730	733	rVV2	65559	120036	2.28%	0.186%
32	10.476	733	734	737	rVB	105109	118999	2.26%	0.184%
33	10.567	739	742	744	rBV	2173108	2640290	50.04%	4.084%
34	10.602	744	745	748	rVV2	49897	71722	1.36%	0.111%

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

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050416.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	10.693	751	753	756	rVB	167697	248610	4.71%	0.385%
36	10.865	763	768	774	rBV4	74295	251103	4.76%	0.388%
37	10.945	774	775	777	rVB2	48157	54942	1.04%	0.085%
38	11.013	777	781	784	rBV4	59852	178123	3.38%	0.276%
39	11.116	788	790	791	rBV	119393	167630	3.18%	0.259%
40	11.150	791	793	800	rVB2	163341	263911	5.00%	0.408%
41	11.288	800	805	807	rBV2	1680663	3359550	63.67%	5.197%
42	11.333	807	809	811	rVB	645212	603437	11.44%	0.933%
43	11.493	819	823	827	rBV3	86067	191060	3.62%	0.296%
44	11.562	827	829	830	rVB2	50498	68881	1.31%	0.107%
45	11.756	844	846	847	rBV	317058	297749	5.64%	0.461%
46	11.790	847	849	851	rVV2	362704	570456	10.81%	0.882%
47	11.836	851	853	854	rVV	162041	205000	3.89%	0.317%
48	11.870	854	856	860	rVV2	468155	714197	13.54%	1.105%
49	11.962	860	864	866	rVV2	34771	79690	1.51%	0.123%
50	12.053	870	872	876	rVB	194014	226300	4.29%	0.350%
51	12.202	883	885	886	rBV	74658	68178	1.29%	0.105%
52	12.305	891	894	896	rBV	157492	217584	4.12%	0.337%
53	12.339	896	897	900	rVB2	85087	159787	3.03%	0.247%
54	12.396	900	902	906	rVB2	68022	111310	2.11%	0.172%
55	12.465	906	908	911	rBV	2180339	2263907	42.91%	3.502%
56	12.510	911	912	917	rVB5	55145	108606	2.06%	0.168%
57	12.613	919	921	925	rVB2	95699	134122	2.54%	0.207%
58	12.693	925	928	930	rBV	2227309	2554571	48.42%	3.952%
59	12.842	936	941	944	rBV	3121124	3318191	62.89%	5.133%
60	12.911	944	947	951	rVV3	170288	357556	6.78%	0.553%
61	13.025	952	957	959	rVV	351670	584131	11.07%	0.904%
62	13.082	961	962	964	rVV	271912	356741	6.76%	0.552%
63	13.128	964	966	970	rVV2	244057	420471	7.97%	0.650%
64	13.219	972	974	975	rVV	118116	128797	2.44%	0.199%
65	13.242	975	976	981	rVV2	127097	188398	3.57%	0.291%
66	13.425	988	992	997	rBV2	110451	305076	5.78%	0.472%
67	13.528	997	1001	1003	rBV2	87684	220202	4.17%	0.341%
68	13.642	1008	1011	1012	rBV2	110632	225135	4.27%	0.348%
69	13.665	1012	1013	1014	rVV	199684	185700	3.52%	0.287%
70	13.745	1018	1020	1024	rVV2	649923	749045	14.20%	1.159%
71	13.882	1029	1032	1034	rBV2	1394374	2421221	45.89%	3.745%

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72	13.996	1039	1042	1043	rBV2	153763	190347	3.61%	0.294%
73	14.065	1046	1048	1050	rBV3	76450	126677	2.40%	0.196%
74	14.271	1064	1066	1068	rBV	170243	217399	4.12%	0.336%
75	14.374	1072	1075	1076	rBV3	110891	185021	3.51%	0.286%
76	14.728	1104	1106	1111	rBV2	70442	161707	3.06%	0.250%
77	14.888	1118	1120	1125	rBV	637349	1378103	26.12%	2.132%
78	14.991	1127	1129	1131	rVB	122037	149520	2.83%	0.231%
79	15.162	1142	1144	1147	rVB	345108	471885	8.94%	0.730%
80	15.219	1147	1149	1152	rVV	522156	687603	13.03%	1.064%
81	15.276	1152	1154	1160	rVB2	495031	926341	17.56%	1.433%
82	16.602	1267	1270	1276	rBV2	155680	415260	7.87%	0.642%
83	16.991	1301	1304	1312	rVB	138415	343108	6.50%	0.531%

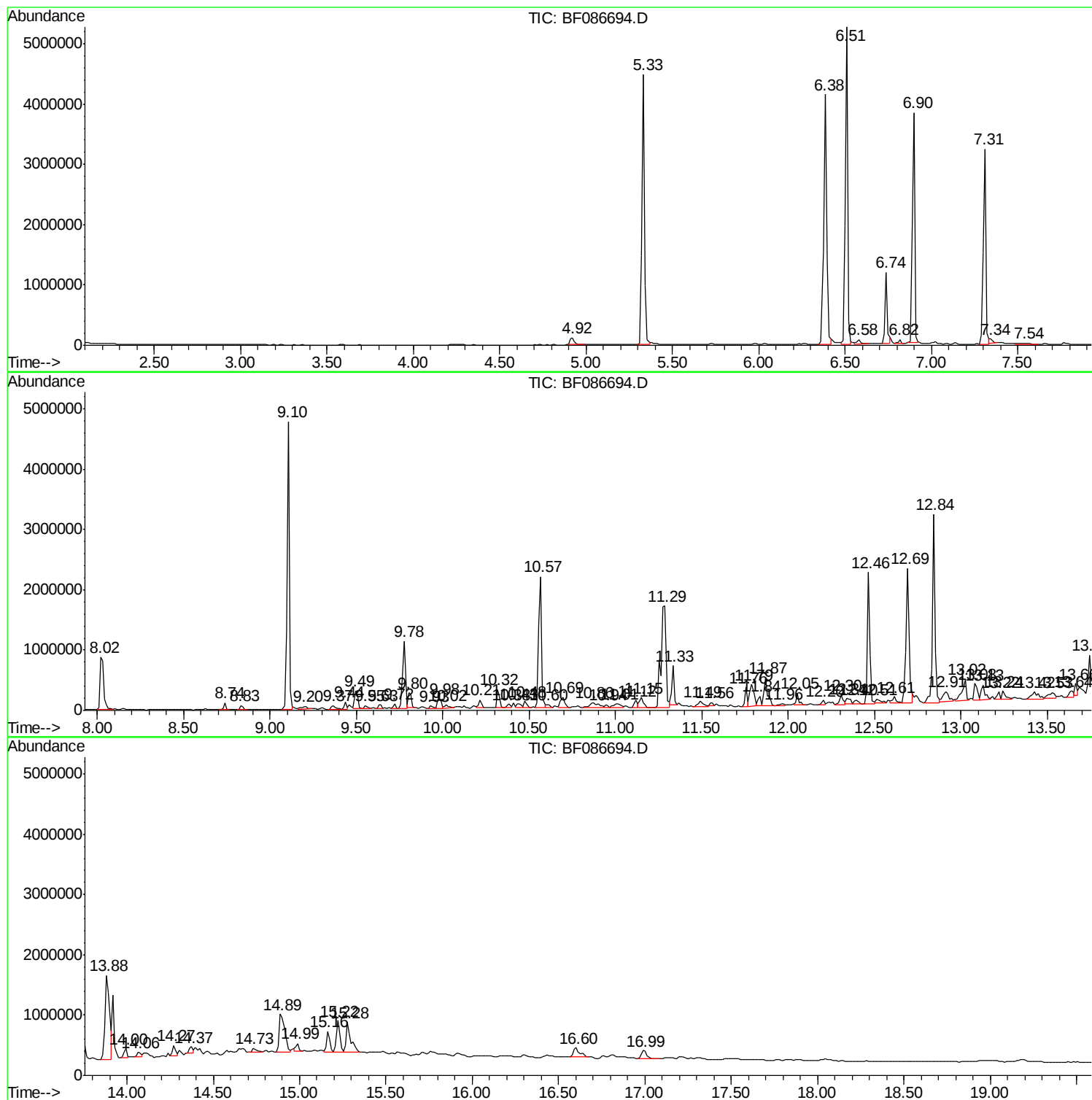
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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Acq On : 4 May 2016 23:31
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Sample : H2715-12
Misc :
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Instrument :
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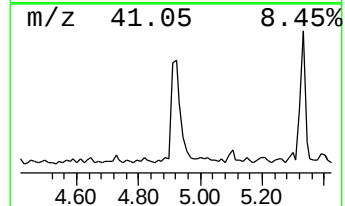
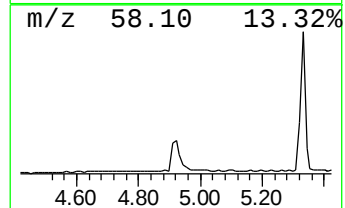
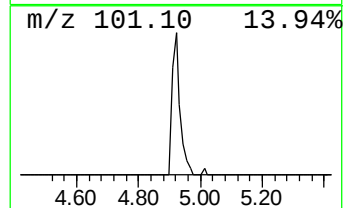
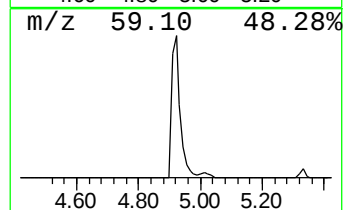
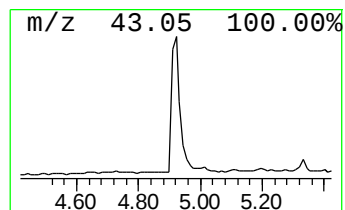
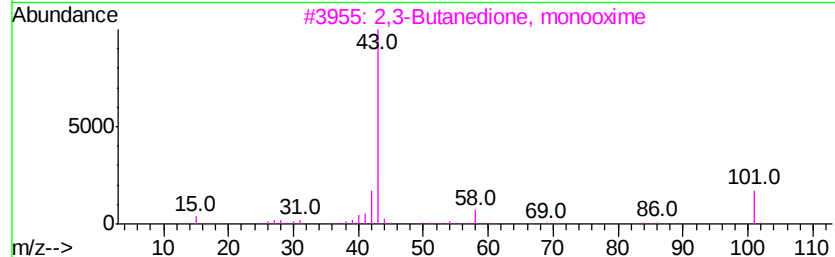
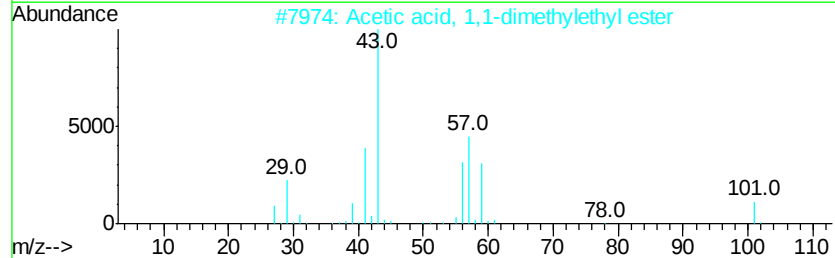
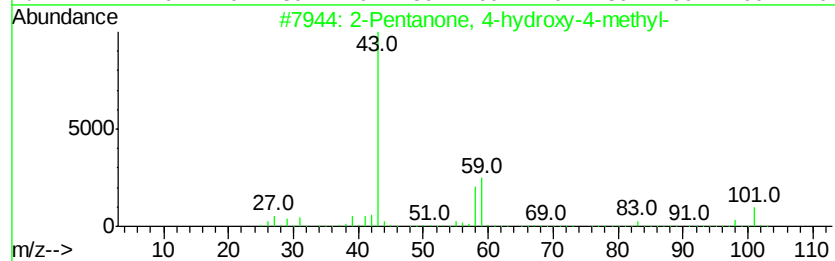
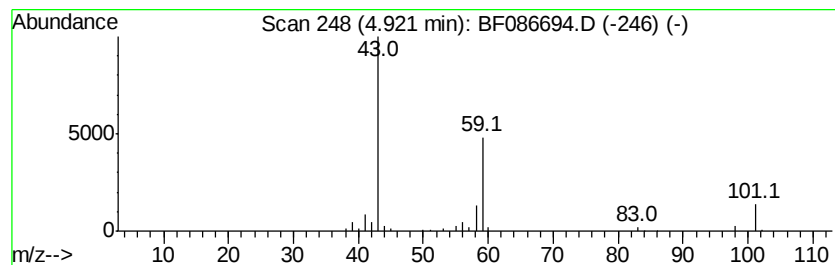
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.92	8.11 ng	225757	1,4-Dichlorobenzene-d4	6.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28		
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16		
4	Diacetamide	101	C4H7NO2	000625-77-4	9		
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9		



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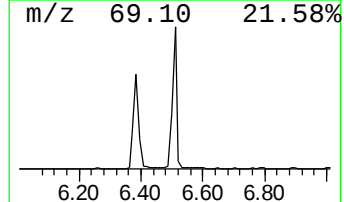
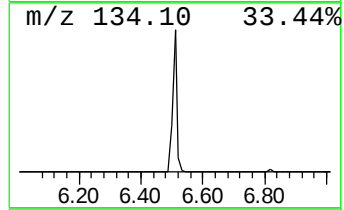
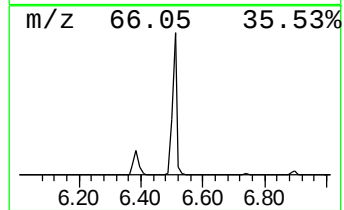
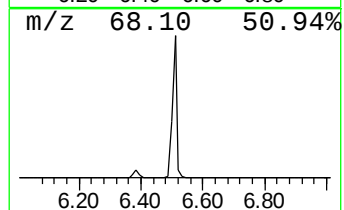
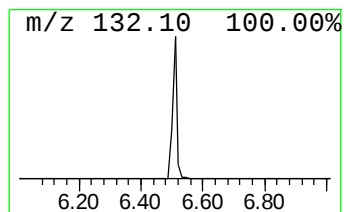
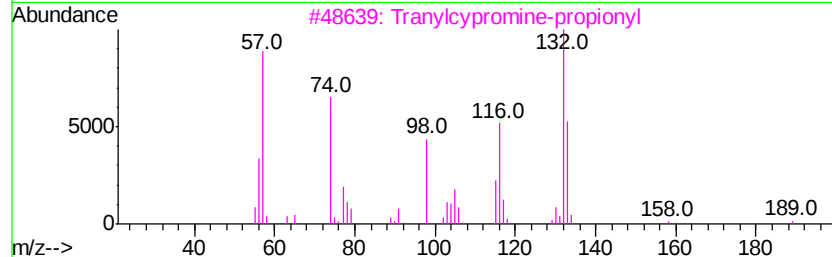
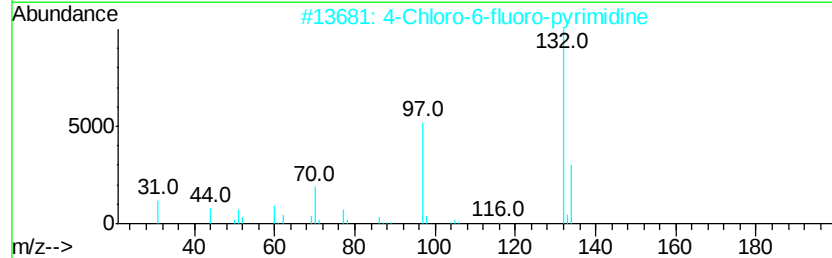
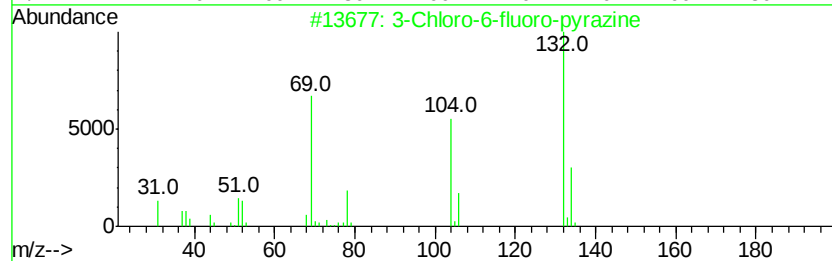
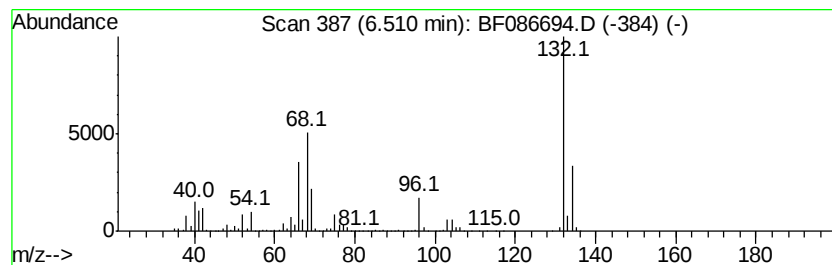
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.51 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	189.56 ng	5276240	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	22
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16



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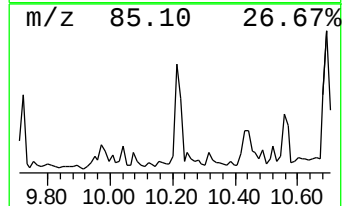
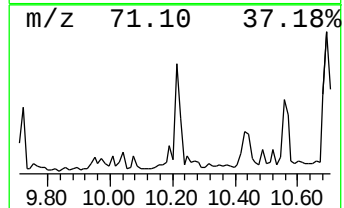
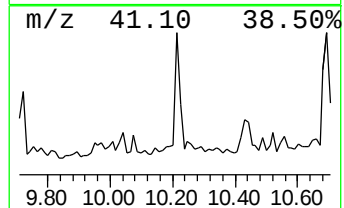
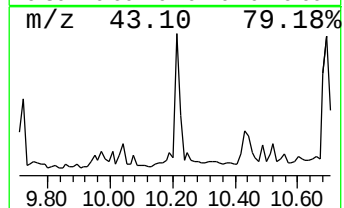
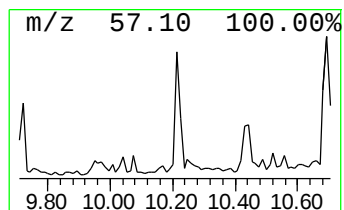
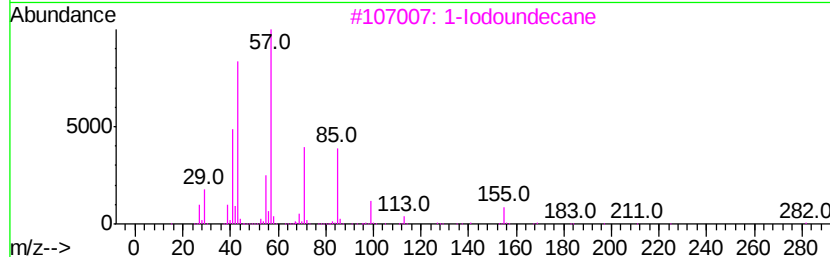
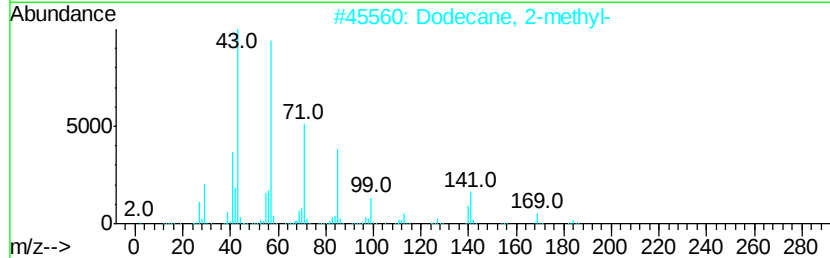
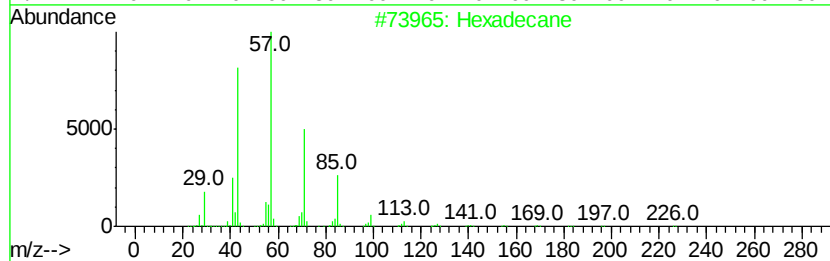
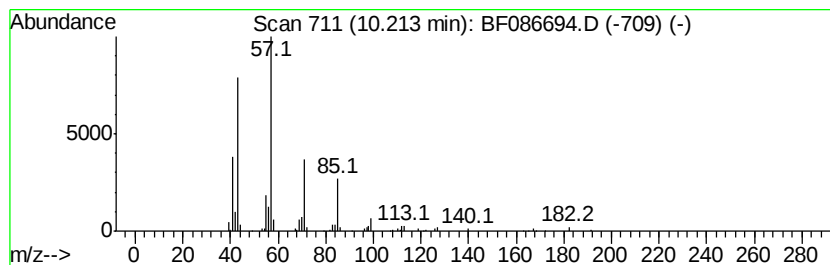
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Hexadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.21	4.94 ng	150229	Acenaphthene-d10	9.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	87
2			Dodecane, 2-methyl-	184	C13H28	001560-97-0	72
3			1-Iodoundecane	282	C11H23I	004282-44-4	72
4			Tridecane, 7-propyl-	226	C16H34	055045-09-5	72
5			Tridecane	184	C13H28	000629-50-5	64



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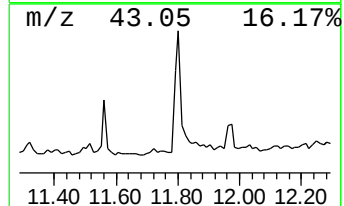
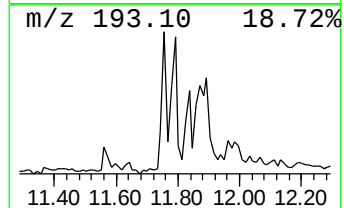
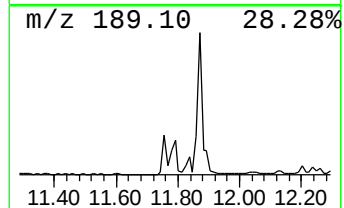
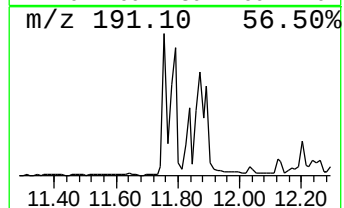
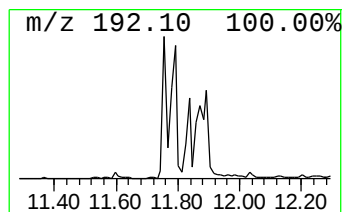
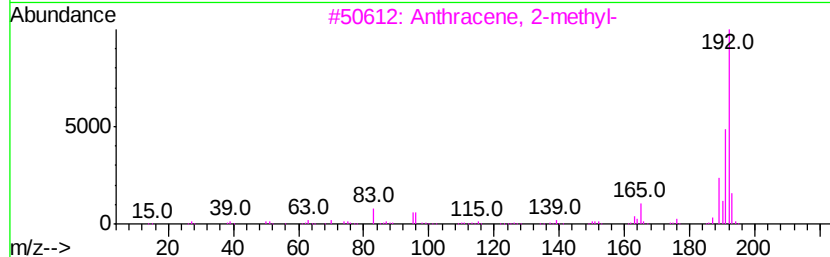
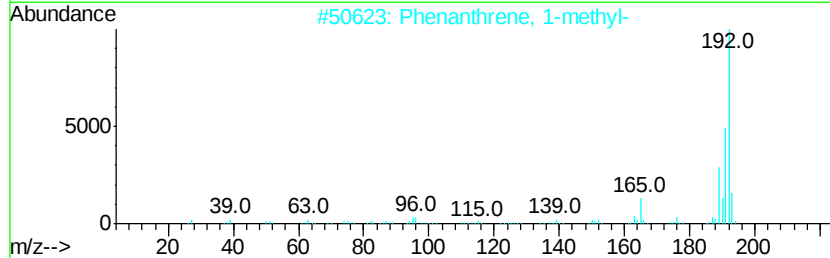
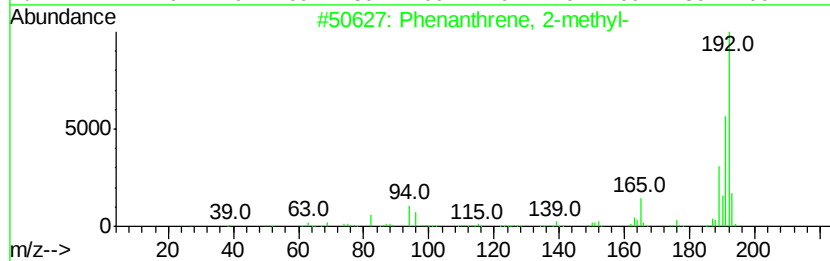
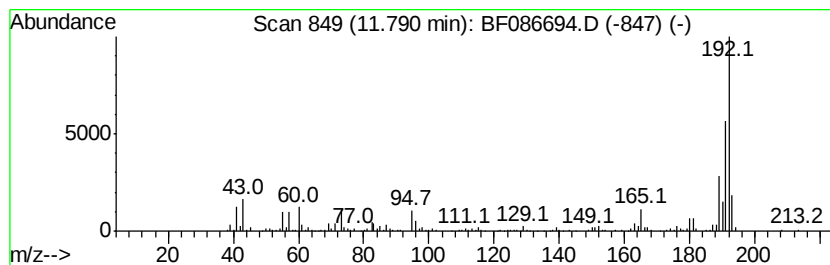
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	6.79 ng	570456	Phenanthrene-d10	11.25

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	94
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050416\
Data File : BF086694.D
Acq On : 4 May 2016 23:31
Operator : UM/SJ
Sample : H2715-12
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
26WARD-4-CS-1(0-8FTBGS)

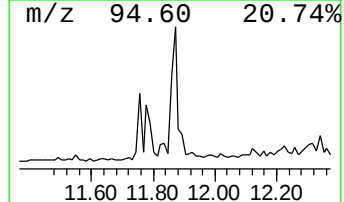
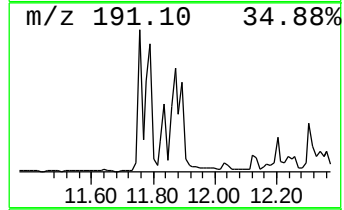
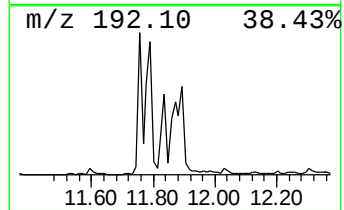
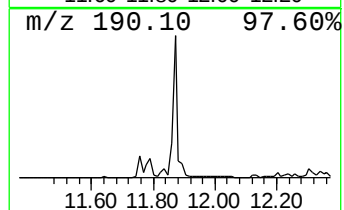
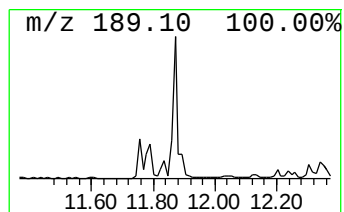
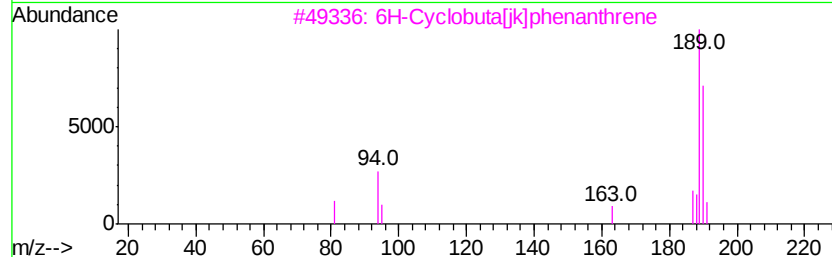
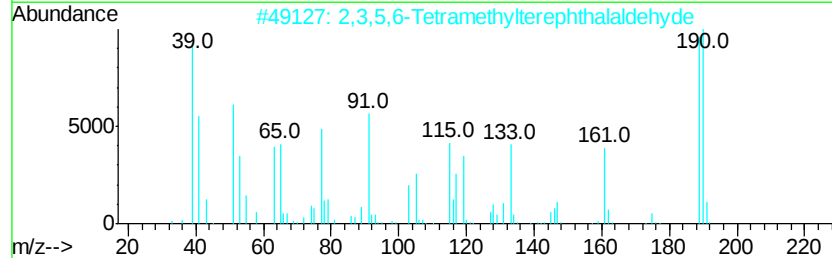
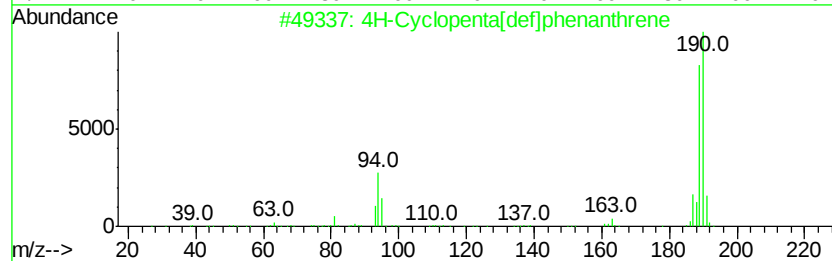
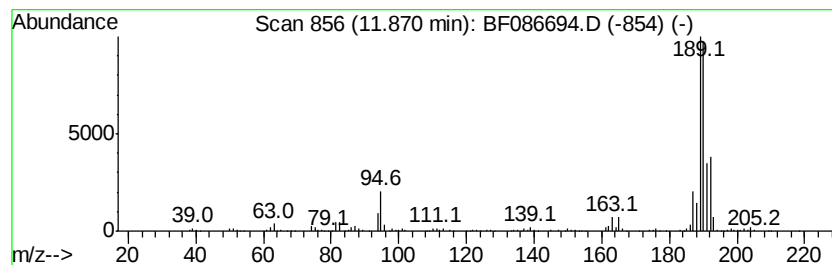
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.87	8.50 ng	714197	Phenanthrene-d10	11.25

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	94
2		2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	47
3		6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	45
4		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	42
5		4,4'-Difluorobiphenyl	190	C12H8F2	000398-23-2	17



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050416\
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Acq On : 4 May 2016 23:31
Operator : UM/SJ
Sample : H2715-12
Misc :
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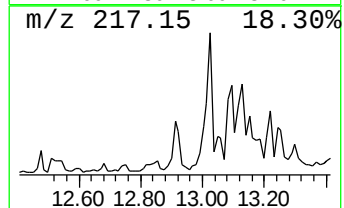
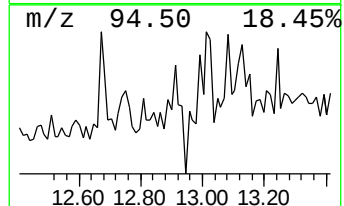
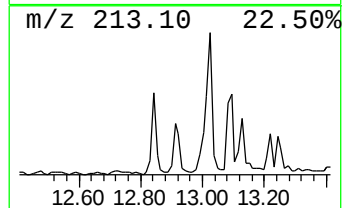
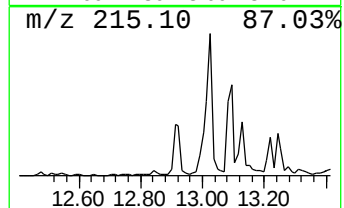
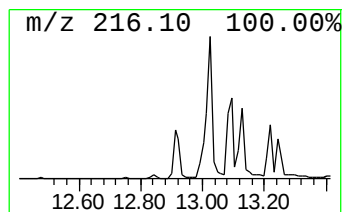
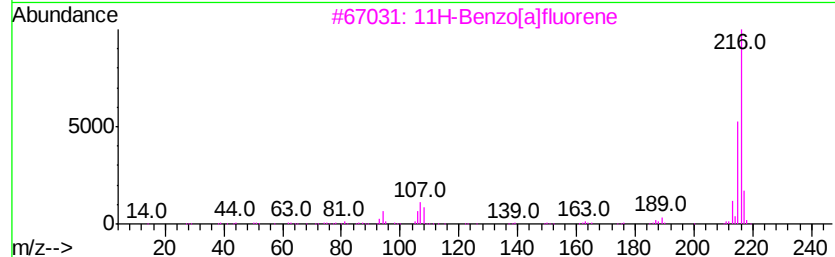
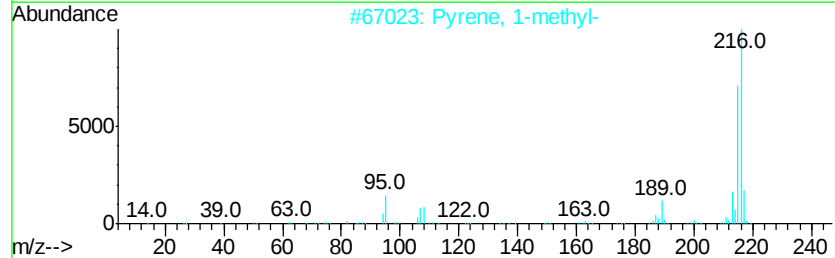
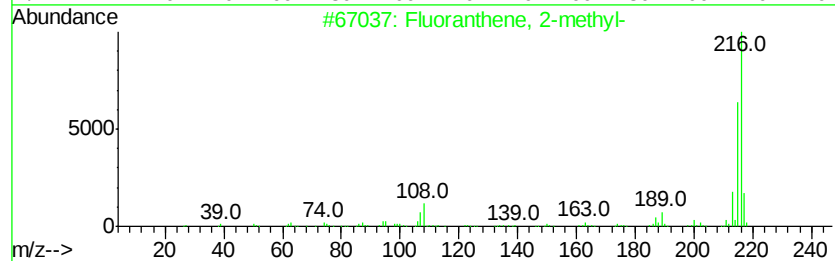
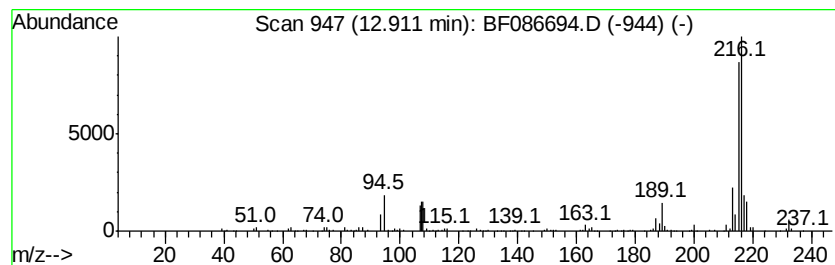
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Fluoranthene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.91	5.91 ng	357556	Chrysene-d12	13.89	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	95
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
4	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050416\
Data File : BF086694.D
Acq On : 4 May 2016 23:31
Operator : UM/SJ
Sample : H2715-12
Misc :
ALS Vial : 16 Sample Multiplier: 1

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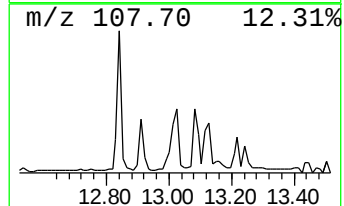
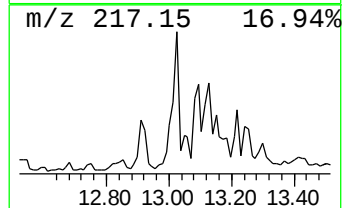
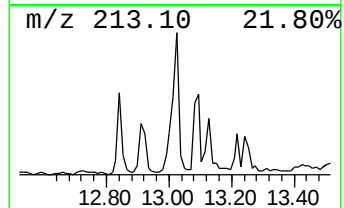
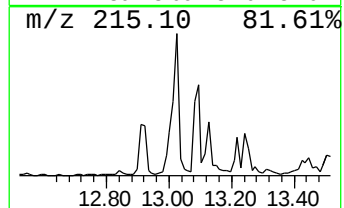
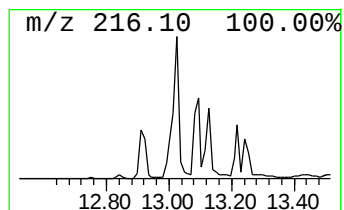
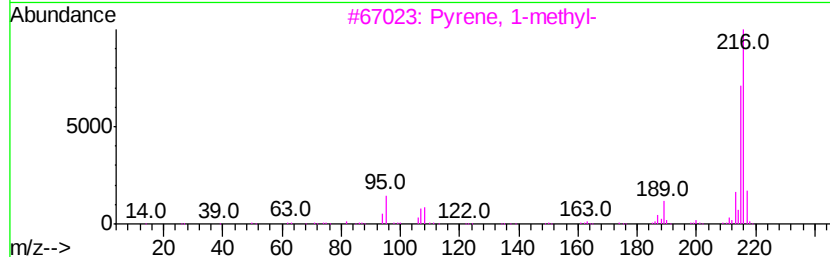
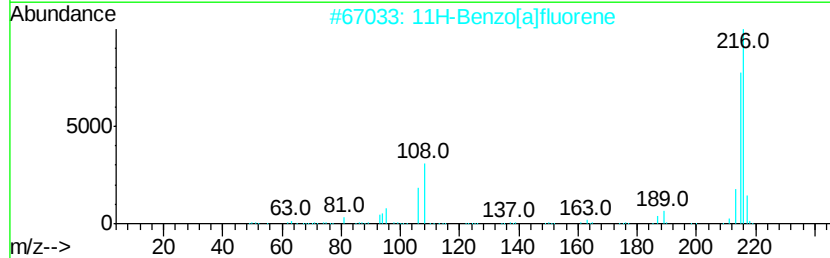
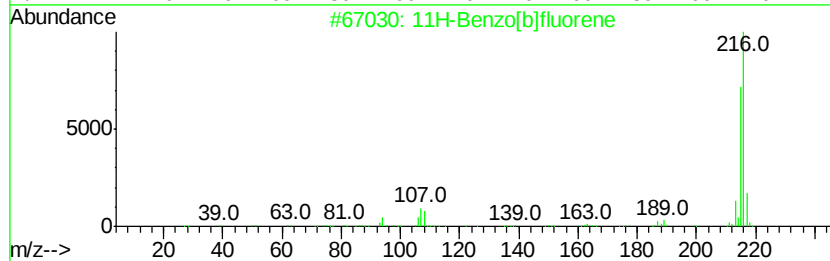
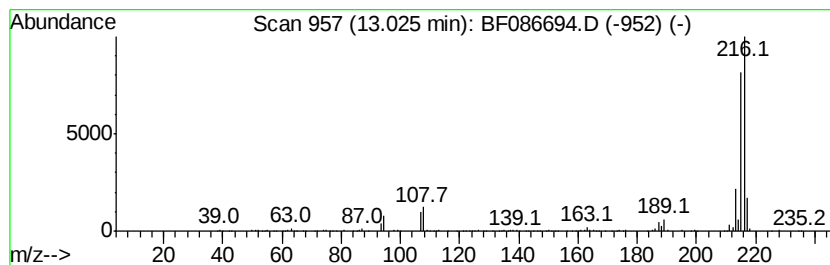
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 11H-Benzo[b]fluorene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	9.65 ng	584131	Chrysene-d12	13.89

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	96
2		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	94
3		Pvrene, 1-methyl-	216	C17H12	002381-21-7	93
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
5		7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050416\
Data File : BF086694.D
Acq On : 4 May 2016 23:31
Operator : UM/SJ
Sample : H2715-12
Misc :
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Instrument :
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26WARD-4-CS-1(0-8FTBGS)

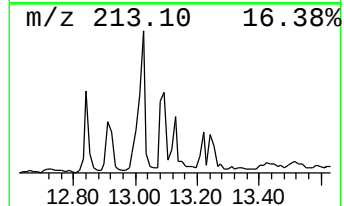
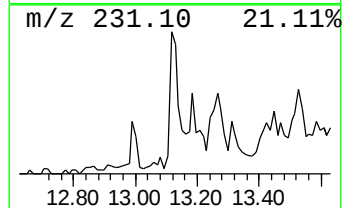
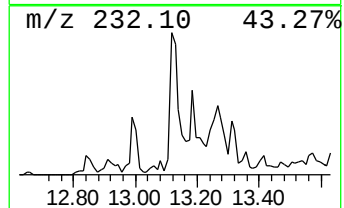
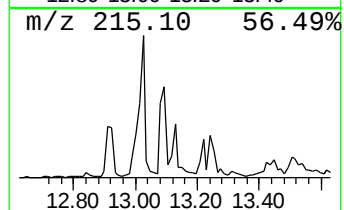
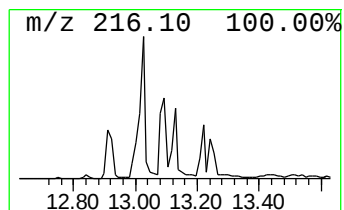
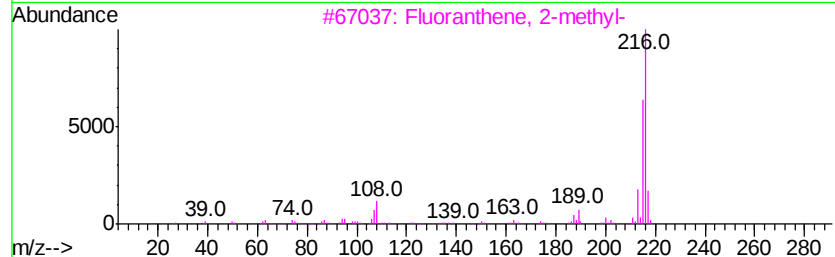
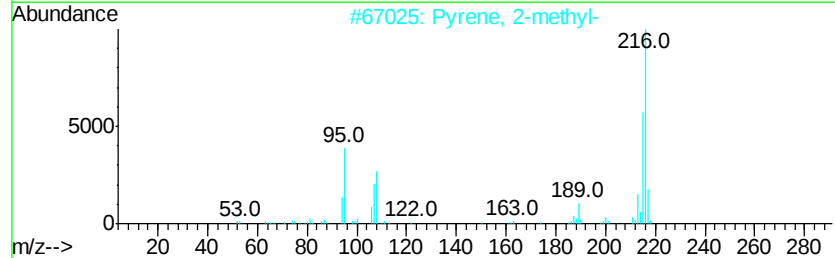
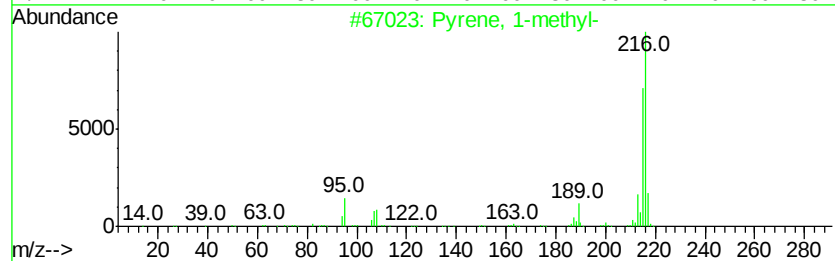
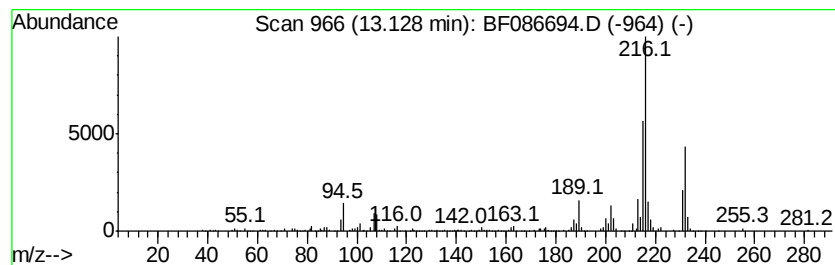
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Pyrene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.13	6.95 ng	420471	Chrysene-d12	13.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
2		Pyrene, 2-methyl-	216	C17H12	003442-78-2	64
3		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	55
4		11H-Benzo[alfluorene	216	C17H12	000238-84-6	53
5		Pyrene, 4-methyl-	216	C17H12	003353-12-6	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050416\
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Acq On : 4 May 2016 23:31
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Sample : H2715-12
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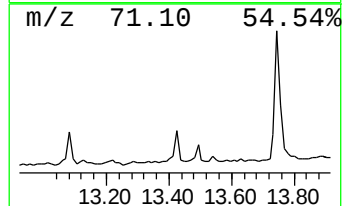
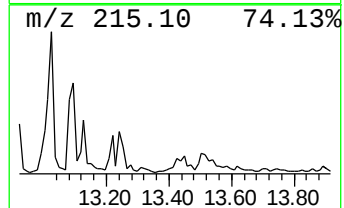
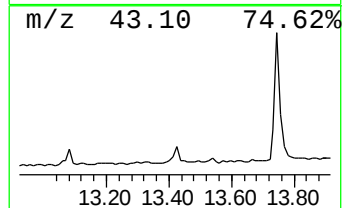
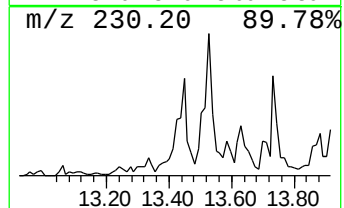
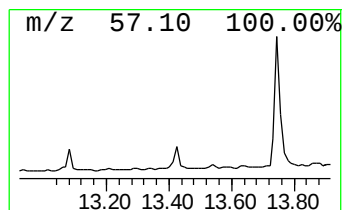
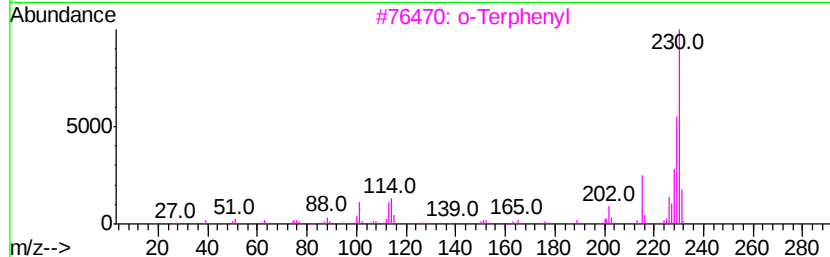
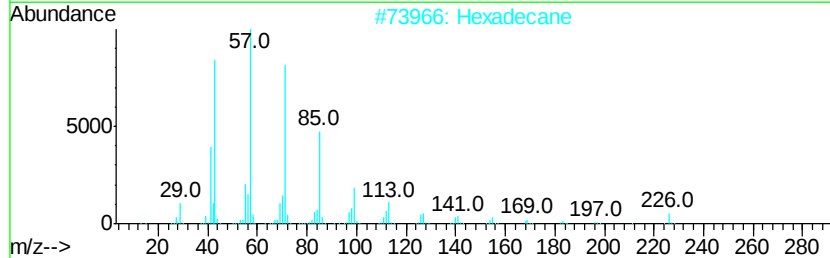
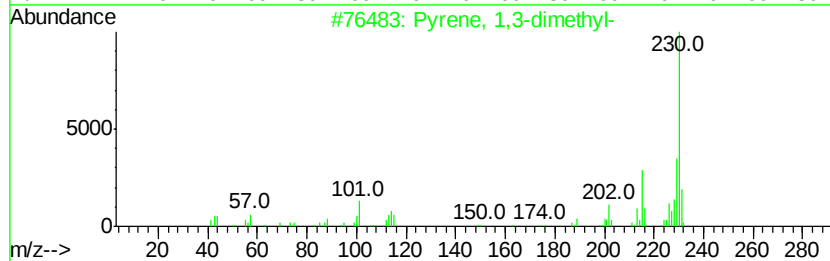
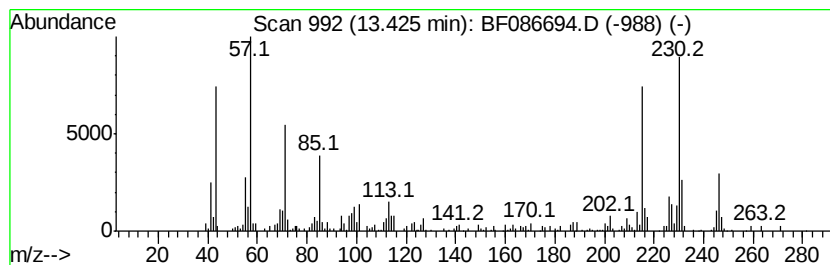
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Pyrene, 1,3-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.42	5.04 ng	305076	Chrysene-d12	13.89	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1,3-dimethyl-	230	C18H14	064401-21-4	55
2	Hexadecane	226	C16H34	000544-76-3	55
3	o-Terphenyl	230	C18H14	000084-15-1	35
4	5,6-Dihydrochrysene	230	C18H14	002091-92-1	30
5	Pyrene, 1-(methoxymethyl)-	246	C18H14O	091385-15-8	30



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050416\
Data File : BF086694.D
Acq On : 4 May 2016 23:31
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ALS Vial : 16 Sample Multiplier: 1

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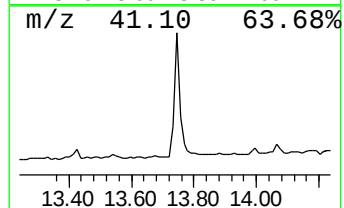
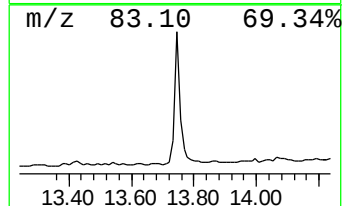
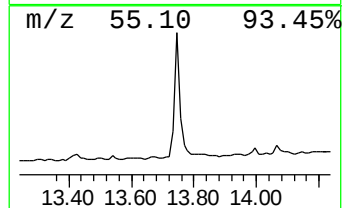
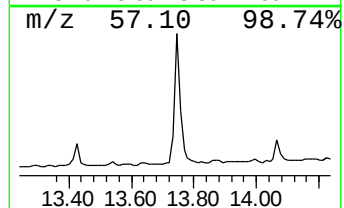
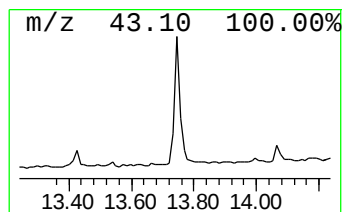
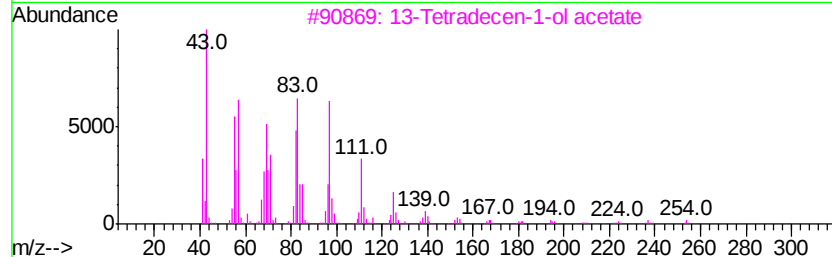
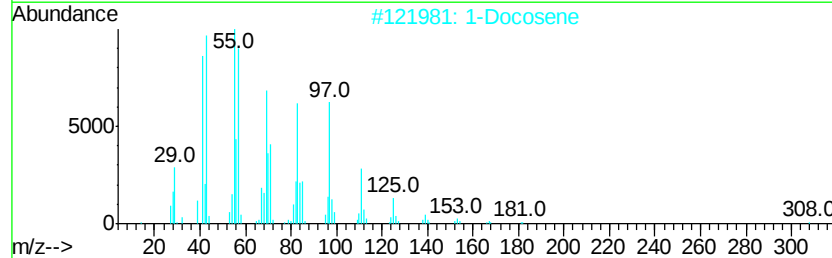
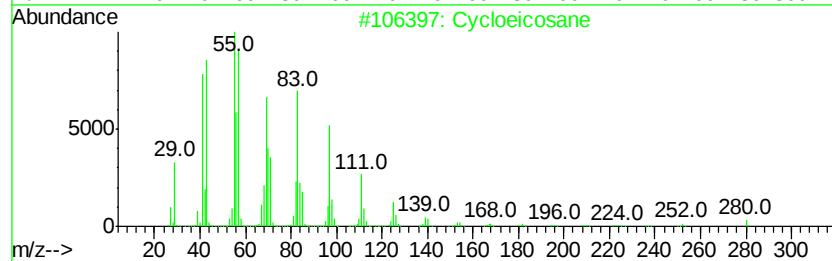
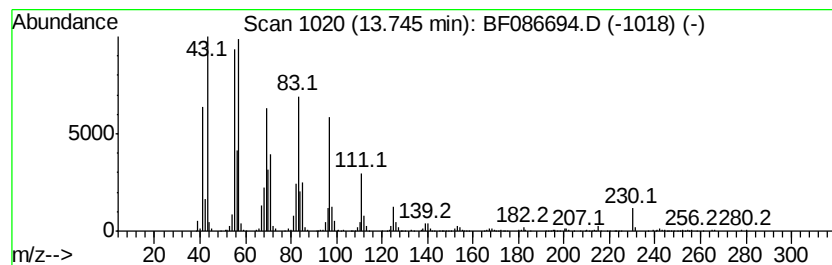
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Cycloeicosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.74	12.37 ng	749045	Chrysene-d12	13.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cycloeicosane	280	C20H40	000296-56-0	98
2		1-Docosene	308	C22H44	001599-67-3	95
3		13-Tetradecen-1-ol acetate	254	C16H30O2	056221-91-1	93
4		Trifluoroacetic acid, n-octadecyl...	366	C20H37F3O2	079392-43-1	91
5		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050416\
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ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
26WARD-4-CS-1(0-8FTBGS)

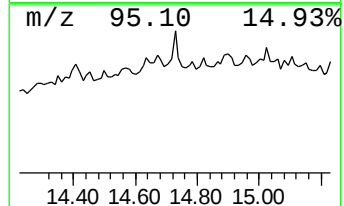
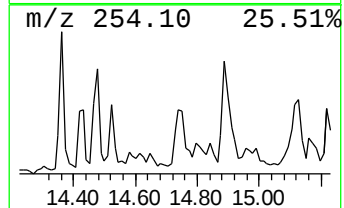
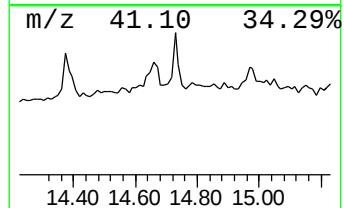
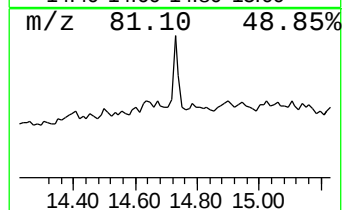
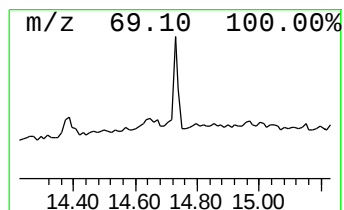
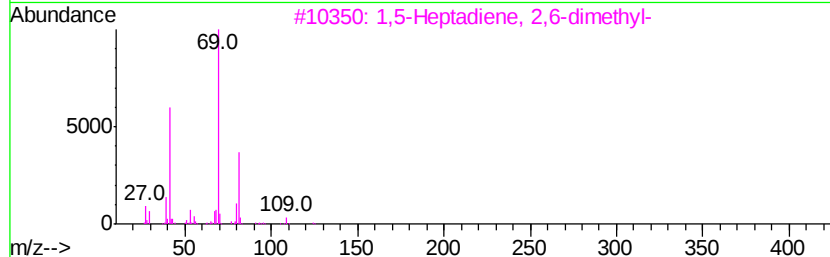
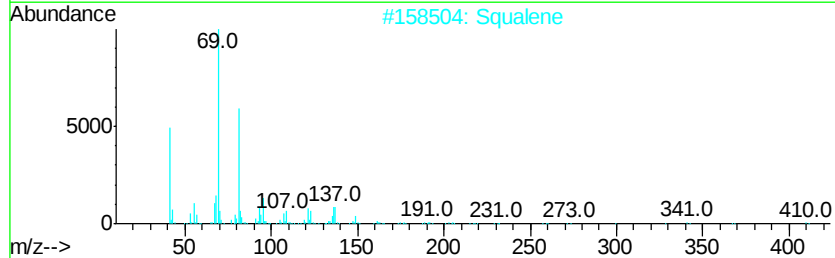
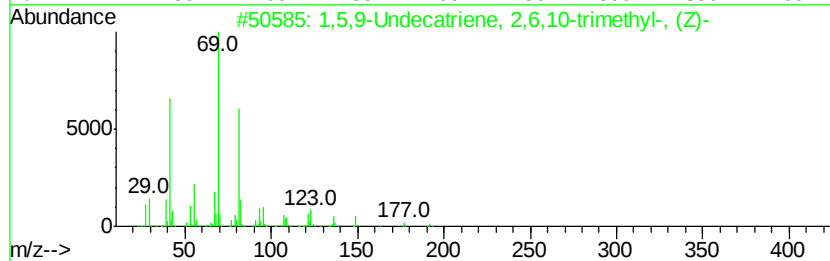
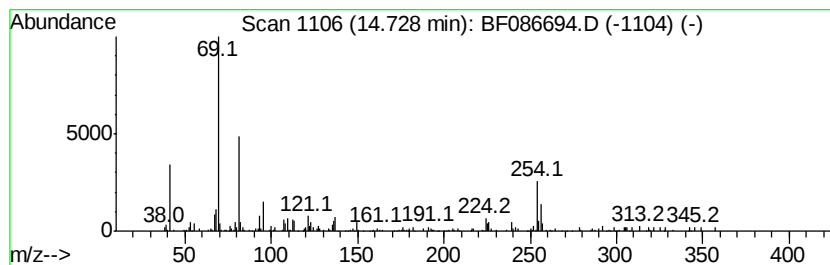
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 1,5,9-Undecatriene, 2,6,10-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.73	6.98 ng	161707	Perylene-d12	15.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,5,9-Undecatriene, 2,6,10-trime...	192	C14H24	062951-96-6	64
2		Squalene	410	C30H50	007683-64-9	50
3		1,5-Heptadiene, 2,6-dimethyl-	124	C9H16	006709-39-3	43
4		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	40
5		2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	40



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050416\
 Data File : BF086694.D
 Acq On : 4 May 2016 23:31
 Operator : UM/SJ
 Sample : H2715-12
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 26WARD-4-CS-1(0-8FTBGS)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.92	8.1 ng		225757	1	6.74	1113340	40.0
unknown6.51	6.51	189.6 ng		5276240	1	6.74	1113340	40.0
Hexadecane	10.21	4.9 ng		150229	3	9.78	1215810	40.0
Phenanthrene, 2-m...	11.79	6.8 ng		570456	4	11.25	3359550	40.0
4H-Cyclopenta[def...	11.87	8.5 ng		714197	4	11.25	3359550	40.0
Fluoranthene, 2-m...	12.91	5.9 ng		357556	5	13.89	2421220	40.0
11H-Benzo[b]fluorene	13.02	9.7 ng		584131	5	13.89	2421220	40.0
Pyrene, 1-methyl-	13.13	7.0 ng		420471	5	13.89	2421220	40.0
Pyrene, 1,3-dimet...	13.42	5.0 ng		305076	5	13.89	2421220	40.0
Cycloeicosane	13.74	12.4 ng		749045	5	13.89	2421220	40.0
1,5,9-Undecatrien...	14.73	7.0 ng		161707	6	15.28	926341	40.0