

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122116\
 Data File : BF091856.D
 Acq On : 22 Dec 2016 1:40
 Operator : UM/SJ
 Sample : H6156-02
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-1101

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-BF122116.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	4.910	244	247	252	rBV	1074088	1309746	12.63%	1.630%
2	5.344	283	285	288	rBV	3956712	4268671	41.16%	5.313%
3	6.385	374	376	379	rBV	2788149	2754472	26.56%	3.428%
4	6.510	385	387	390	rBV	5231957	7597085	73.26%	9.456%
5	6.739	405	407	410	rVV	1523975	1907977	18.40%	2.375%
6	6.807	410	413	417	rVV2	98366	180413	1.74%	0.225%
7	6.899	418	421	424	rVV	5567666	6926457	66.79%	8.621%
8	7.002	426	430	433	rVV2	89924	110149	1.06%	0.137%
9	7.310	454	457	460	rBV	4265148	6329948	61.04%	7.878%
10	7.402	462	465	466	rVV	146199	191265	1.84%	0.238%
11	7.447	466	469	473	rVV	149213	228543	2.20%	0.284%
12	7.527	473	476	479	rVB	78028	136274	1.31%	0.170%
13	7.699	486	491	495	rVB3	58414	119469	1.15%	0.149%
14	7.779	495	498	500	rBV2	138561	269155	2.60%	0.335%
15	7.813	500	501	504	rVB	228119	385756	3.72%	0.480%
16	8.030	517	520	524	rBV	2692144	2816033	27.15%	3.505%
17	8.533	556	564	568	rBV3	92532	273902	2.64%	0.341%
18	8.842	588	591	593	rBV2	159218	185936	1.79%	0.231%
19	9.116	611	615	617	rBV	7458228	10370569	100.00%	12.908%
20	9.208	618	623	628	rVB6	45882	111357	1.07%	0.139%
21	9.311	628	632	634	rVB2	116726	166548	1.61%	0.207%
22	9.368	634	637	641	rBV5	47241	128700	1.24%	0.160%
23	9.505	648	649	651	rVV	141280	127187	1.23%	0.158%
24	9.562	651	654	657	rVV	105577	208369	2.01%	0.259%
25	9.642	657	661	663	rVB2	183030	228117	2.20%	0.284%
26	9.779	670	673	675	rBV	2856989	3207371	30.93%	3.992%
27	9.813	675	676	678	rVB	2041261	1556053	15.00%	1.937%
28	10.225	706	712	715	rVB5	99454	227145	2.19%	0.283%
29	10.293	715	718	720	rBV4	86012	198499	1.91%	0.247%
30	10.328	720	721	724	rVB	126822	115598	1.11%	0.144%
31	10.373	724	725	726	rBV	108748	129380	1.25%	0.161%
32	10.419	726	729	730	rVV2	164792	281164	2.71%	0.350%
33	10.442	730	731	734	rVB2	103818	179662	1.73%	0.224%
34	10.568	738	742	745	rBV2	3913693	6262507	60.39%	7.795%

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35	10.694	751	753	757	rVB	333769	346642	3.34%	0.431%
36	10.865	765	768	770	rBV2	60003	121744	1.17%	0.152%
37	10.911	770	772	774	rVV3	79952	122648	1.18%	0.153%
38	11.139	788	792	796	rBV4	141107	311209	3.00%	0.387%
39	11.219	796	799	800	rVV3	63709	109906	1.06%	0.137%
40	11.265	800	803	805	rVV	3038664	3025179	29.17%	3.765%
41	11.299	805	806	810	rVV4	64133	118909	1.15%	0.148%
42	11.482	820	822	824	rBV2	66091	110069	1.06%	0.137%
43	11.562	827	829	833	rVB3	63030	131890	1.27%	0.164%
44	11.654	833	837	838	rBV2	69265	148351	1.43%	0.185%
45	11.756	844	846	848	rBV2	96987	171496	1.65%	0.213%
46	11.802	848	850	852	rVB3	106407	153522	1.48%	0.191%
47	11.836	852	853	855	rBV	83415	114872	1.11%	0.143%
48	11.871	855	856	859	rVB	224500	245261	2.36%	0.305%
49	11.939	859	862	865	rBV2	92509	198758	1.92%	0.247%
50	12.282	891	892	897	rVB2	48445	120329	1.16%	0.150%
51	12.694	925	928	930	rBV	218500	275158	2.65%	0.342%
52	12.854	939	942	944	rBV	7203437	8006646	77.21%	9.965%
53	13.002	953	955	960	rVB	236157	247046	2.38%	0.307%
54	13.757	1019	1021	1027	rBV2	76190	144009	1.39%	0.179%
55	13.894	1030	1033	1036	rBV	2689031	2524330	24.34%	3.142%
56	14.980	1126	1128	1131	rBV	122608	142101	1.37%	0.177%
57	15.288	1152	1155	1157	rBV	1664407	1995082	19.24%	2.483%
58	15.334	1157	1159	1163	rVB	198769	292893	2.82%	0.365%
59	15.723	1190	1193	1197	rBV	225268	354923	3.42%	0.442%
60	16.180	1230	1233	1238	rVB	235538	404697	3.90%	0.504%
61	16.717	1276	1280	1287	rVB	157288	332060	3.20%	0.413%
62	17.346	1330	1335	1340	rBV	115106	290536	2.80%	0.362%
63	18.089	1395	1400	1406	rVB	57276	170850	1.65%	0.213%
64	18.969	1473	1477	1483	rVB2	37735	124425	1.20%	0.155%

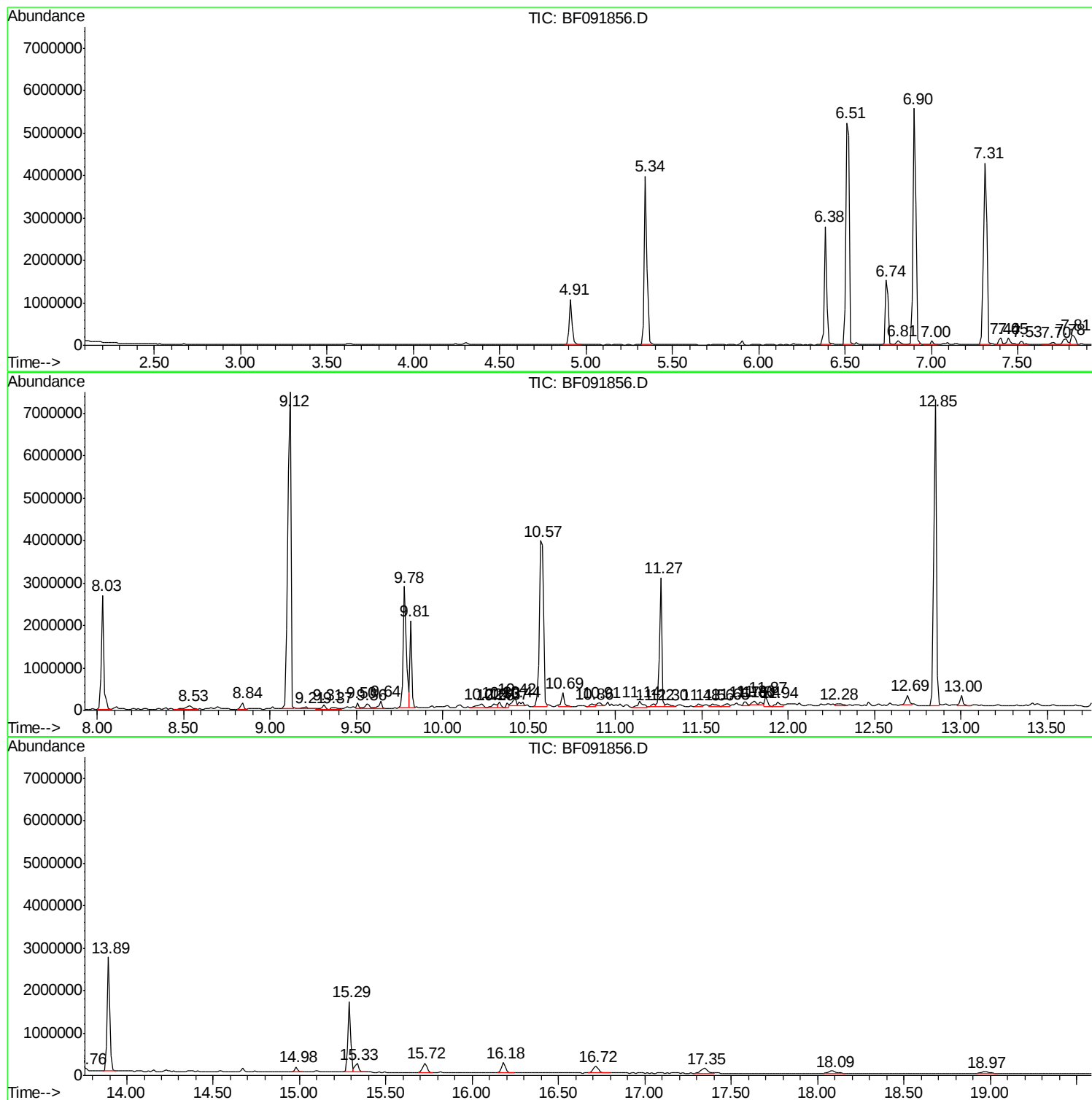
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ClientSampled :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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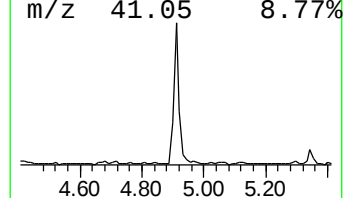
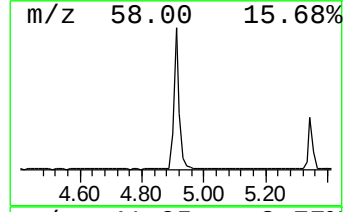
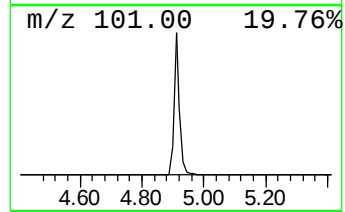
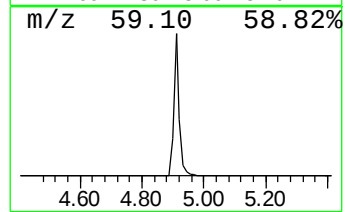
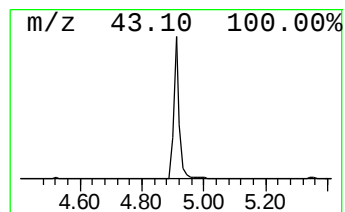
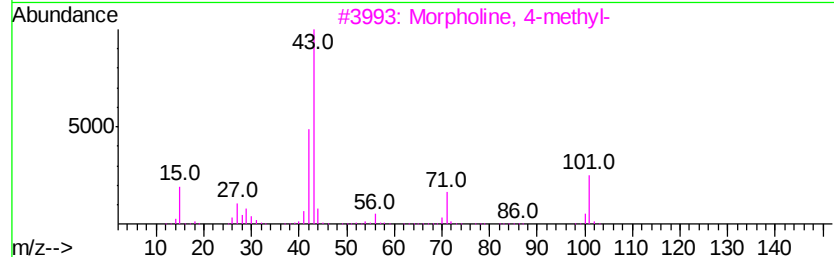
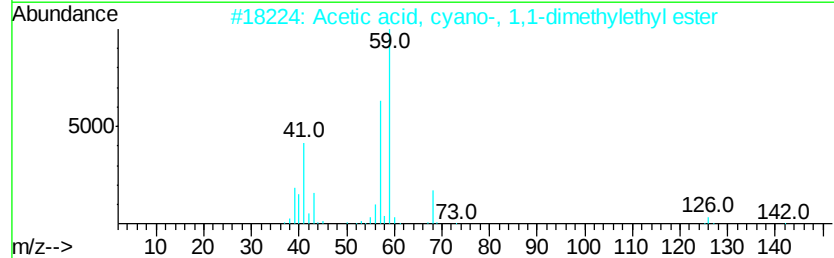
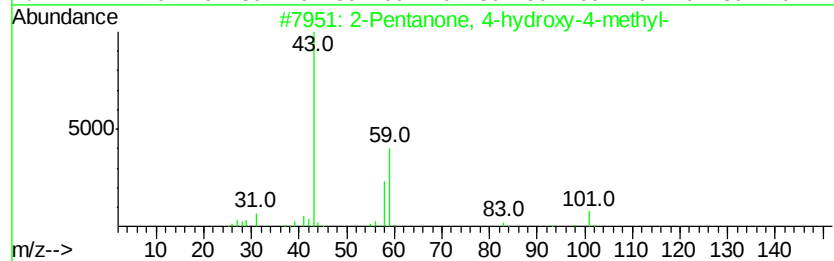
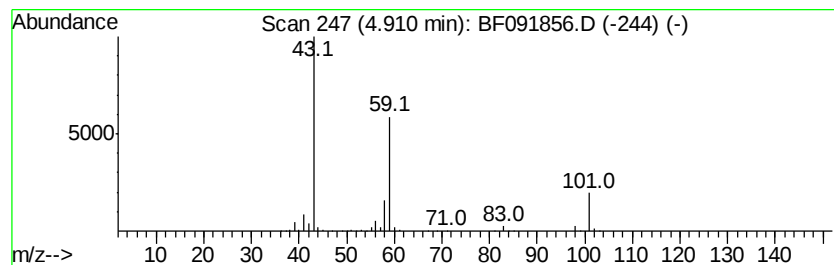
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.91	13.73 ng	1309750	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9



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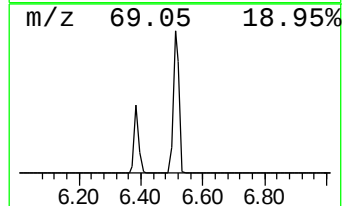
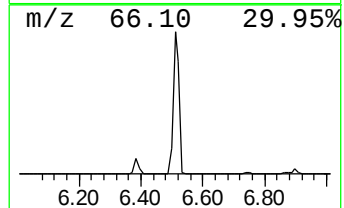
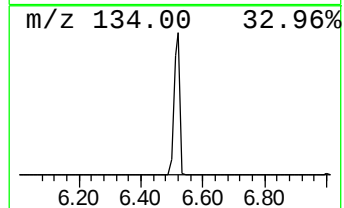
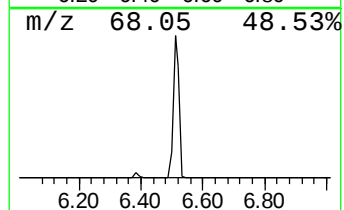
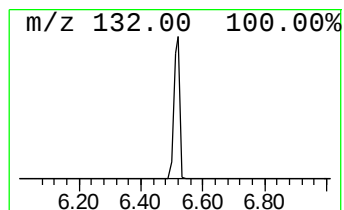
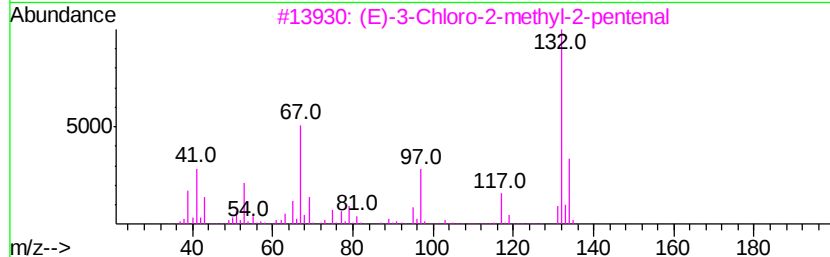
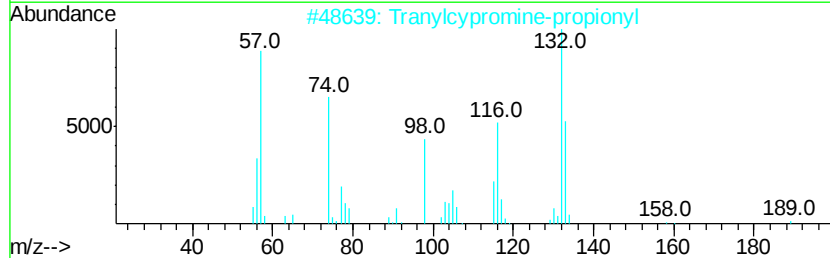
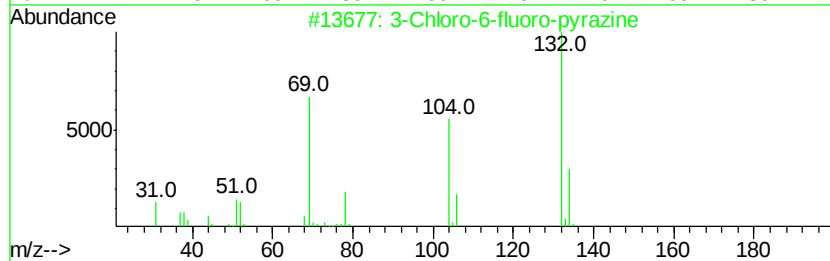
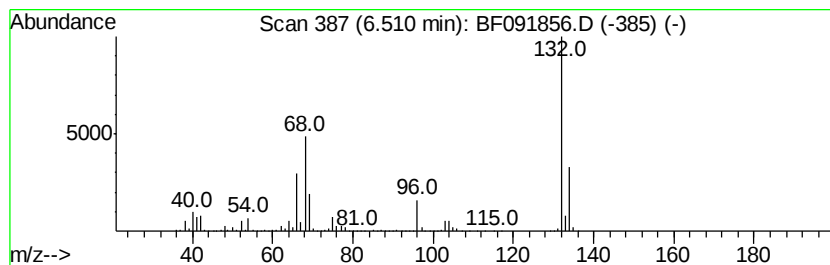
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.51 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	79.64 ng	7597090	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		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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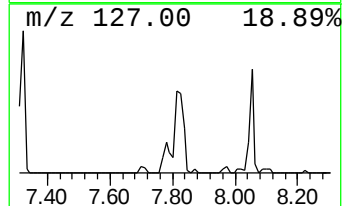
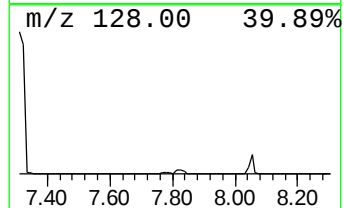
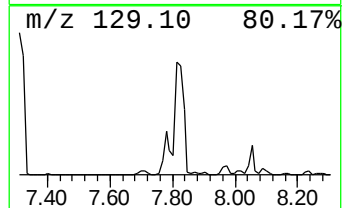
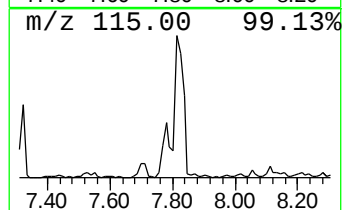
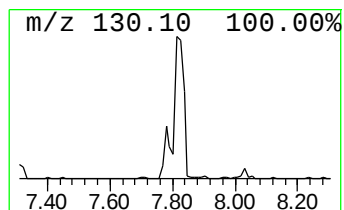
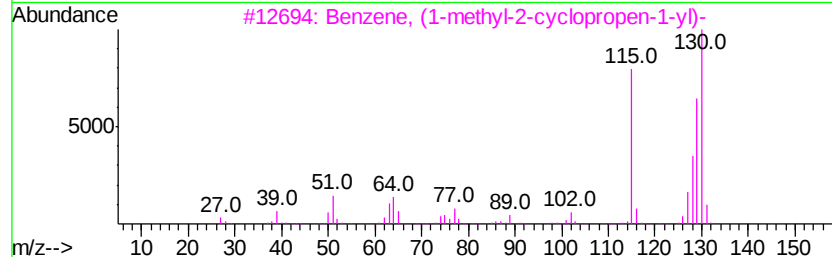
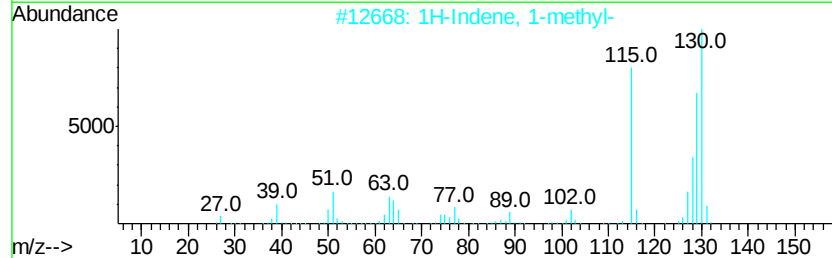
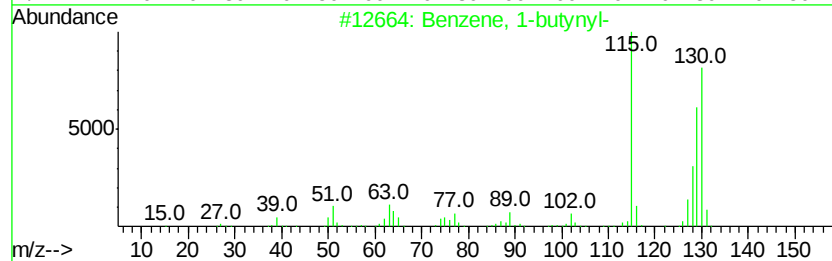
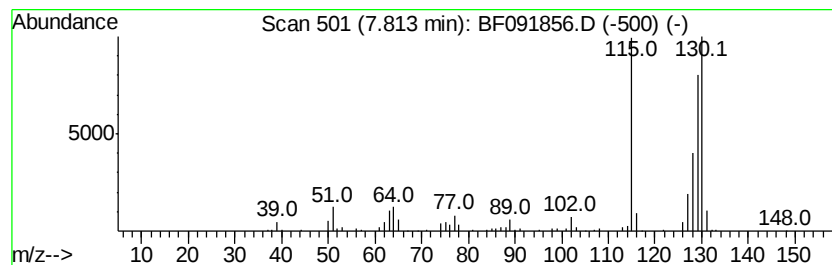
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Peak Number 4 Benzene, 1-butynyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.81	2.74 ng	385756	Naphthalene-d8	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-butynyl-	130	C10H10	000622-76-4	96
2		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	94
3		Benzene, (1-methyl-2-cyclopropen-1-yl)-	130	C10H10	065051-83-4	94
4		2-Methylindene	130	C10H10	002177-47-1	94
5		1,4-Dihydronaphthalene	130	C10H10	000612-17-9	92



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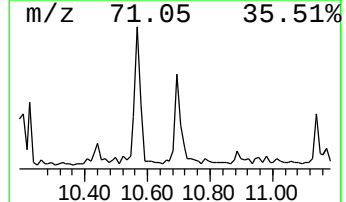
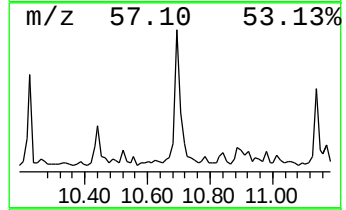
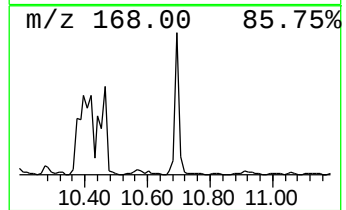
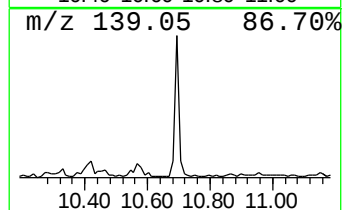
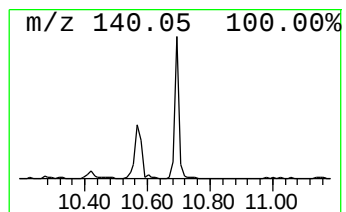
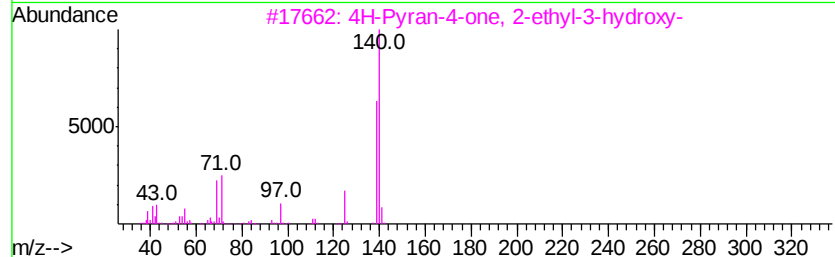
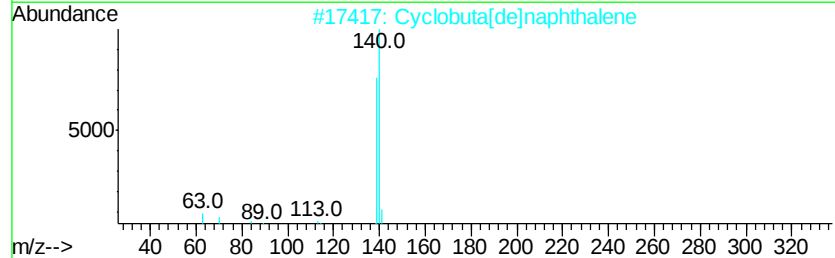
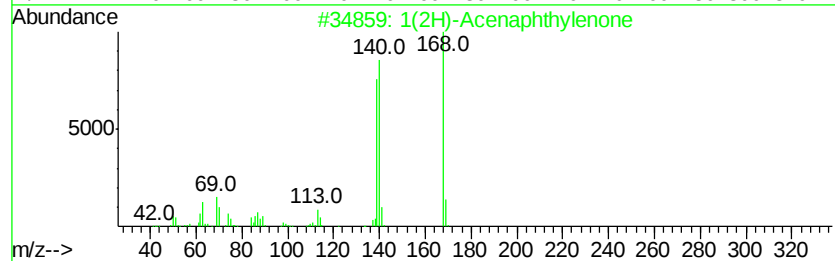
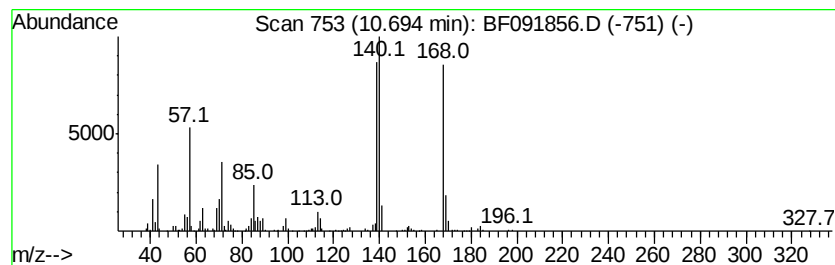
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 1(2H)-Acenaphthylenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.69	2.29 ng	346642	Phenanthrene-d10	11.27		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1(2H)-Acenaphthvlenone	168	C12H8O	002235-15-6	95
2		Cyclobuta[de]naphthalene	140	C11H8	024973-91-9	55
3		4H-Pyran-4-one, 2-ethyl-3-hydroxy-	140	C7H8O3	004940-11-8	47
4		7(4H)-Benzothiazolone, 2-amino-5...	168	C7H8N2OS	017583-10-7	47
5		1H-Inden-1-one, 4,7-difluoro-2,3...	168	C9H6F2O	130408-16-1	47



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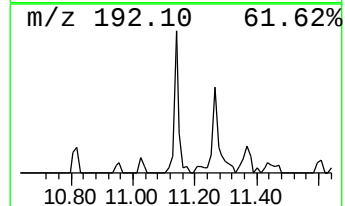
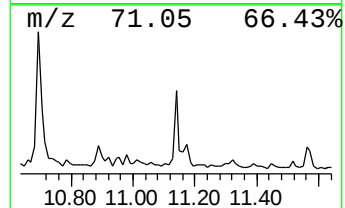
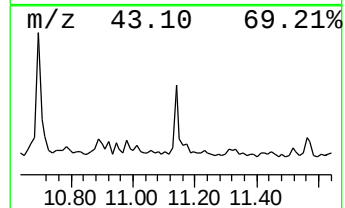
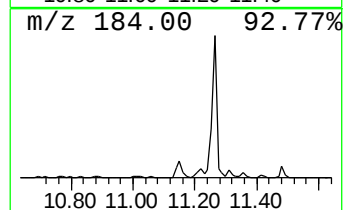
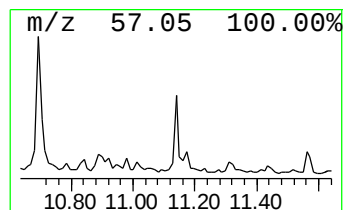
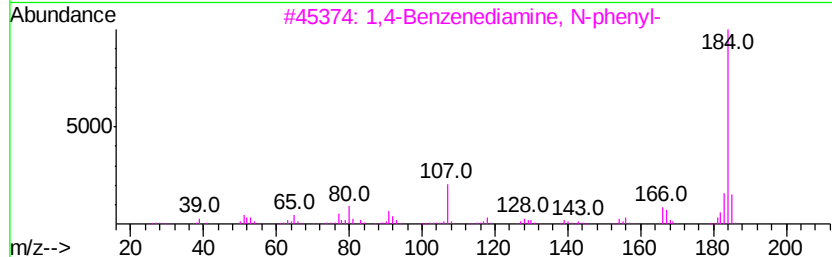
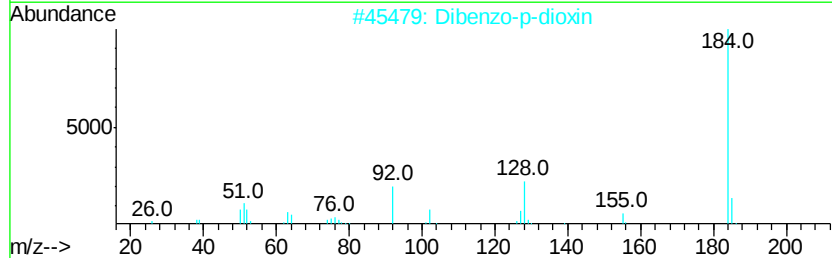
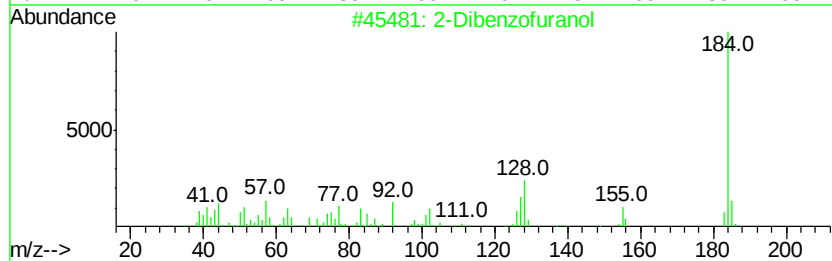
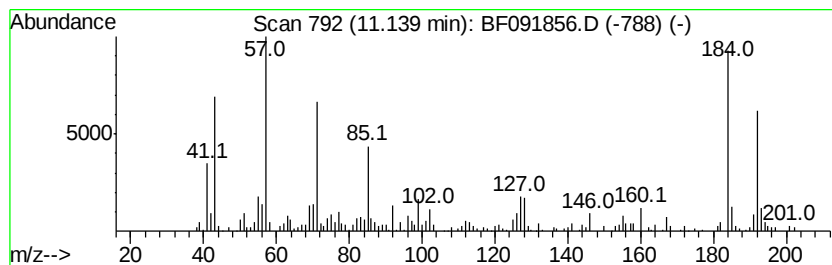
Instrument :
BNA_F
ClientSampleId :
MW-1101

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2-Dibenzofuranol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.14	2.06 ng	311209	Phenanthrene-d10	11.27	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Dibenzofuranol	184	C12H8O2	000086-77-1	86
2	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	38
3	1,4-Benzenediamine, N-phenyl-	184	C12H12N2	000101-54-2	25
4	2,2'-Bipyridine, 4,4'-dimethyl-	184	C12H12N2	001134-35-6	25
5	Benzidine	184	C12H12N2	000092-87-5	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122116\
Data File : BF091856.D
Acq On : 22 Dec 2016 1:40
Operator : UM/SJ
Sample : H6156-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-1101

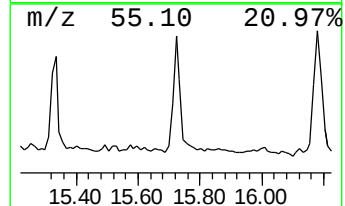
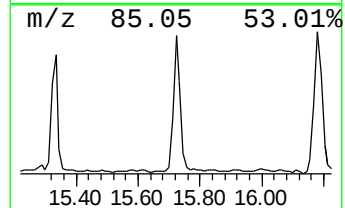
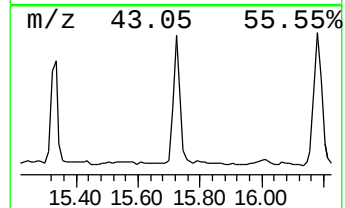
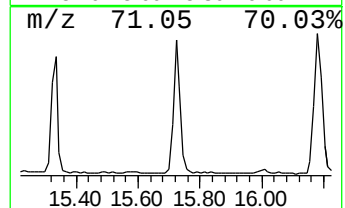
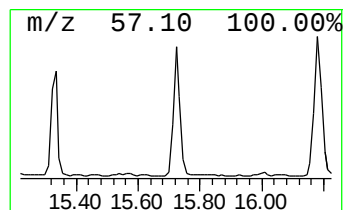
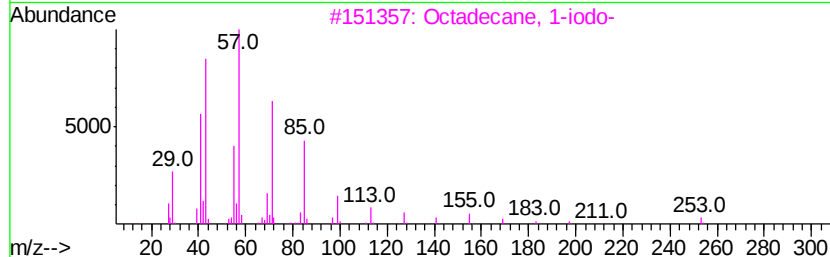
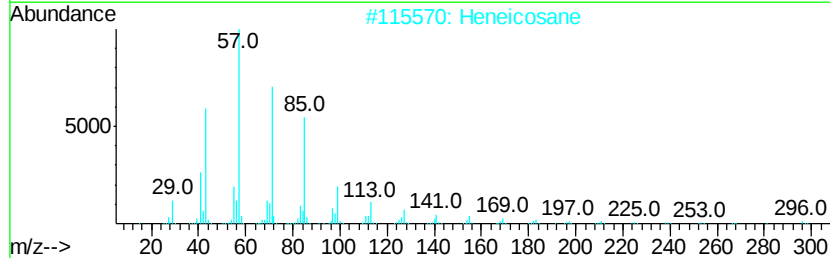
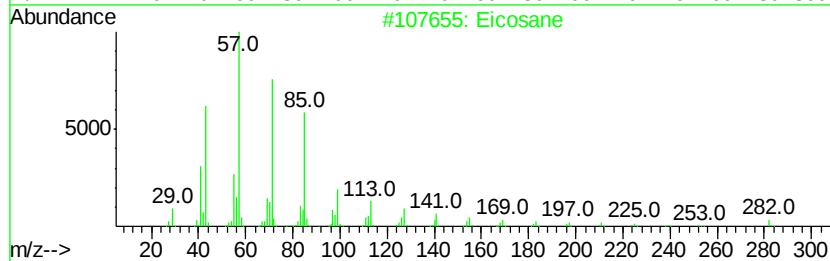
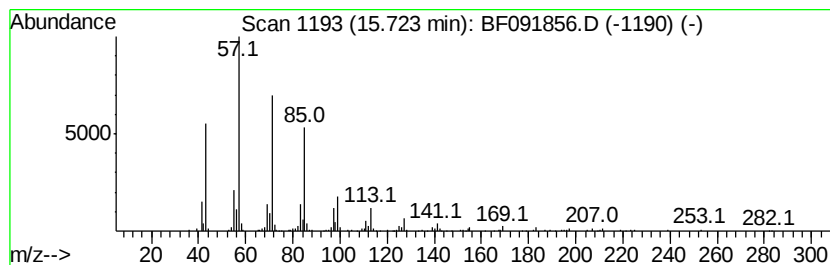
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.72	3.56 ng	354923	Perylene-d12	15.29

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	91
2	Heneicosane		296	C21H44	000629-94-7	90
3	Octadecane, 1-iodo-		380	C18H37I	000629-93-6	86
4	Hexacosane		366	C26H54	000630-01-3	86
5	Pentacosane		352	C25H52	000629-99-2	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122116\
Data File : BF091856.D
Acq On : 22 Dec 2016 1:40
Operator : UM/SJ
Sample : H6156-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-1101

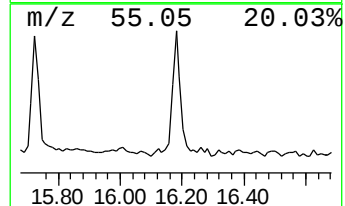
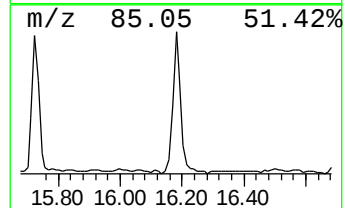
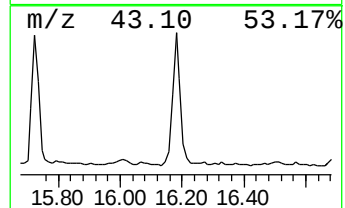
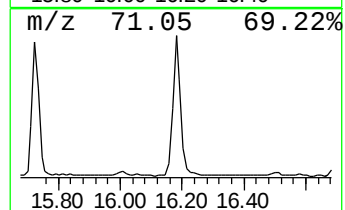
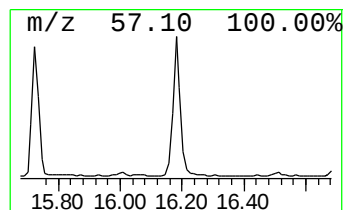
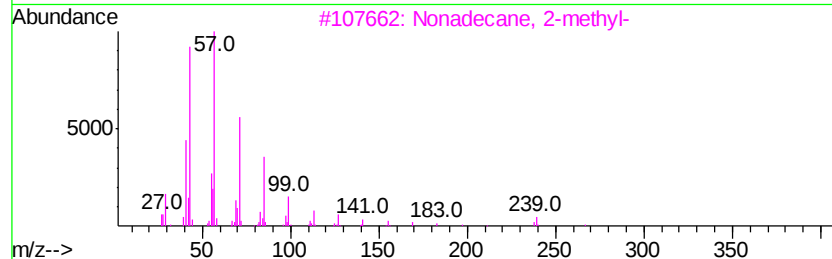
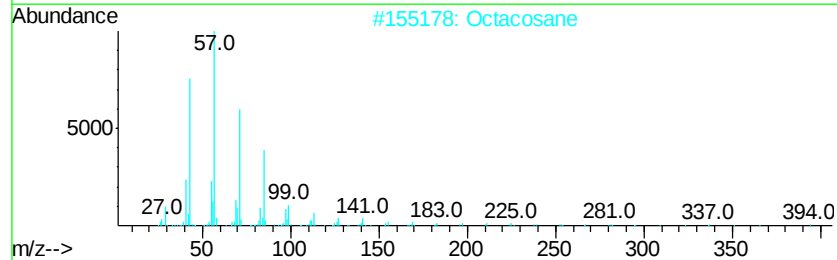
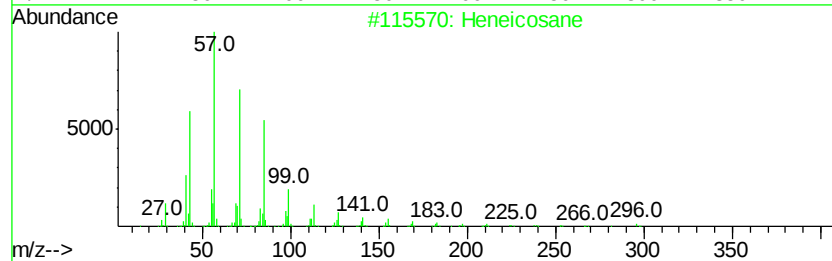
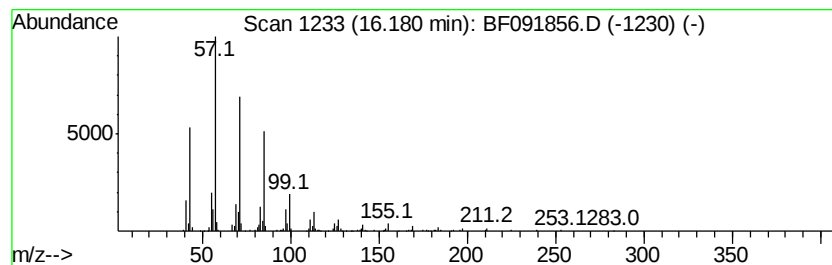
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heneicosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.18	4.06 ng	404697	Perylene-d12	15.29

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Octacosane		394	C28H58	000630-02-4	90
3	Nonadecane, 2-methyl-		282	C20H42	001560-86-7	90
4	Pentadecane, 8-heptyl-		310	C22H46	071005-15-7	87
5	Octadecane		254	C18H38	000593-45-3	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122116\
Data File : BF091856.D
Acq On : 22 Dec 2016 1:40
Operator : UM/SJ
Sample : H6156-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

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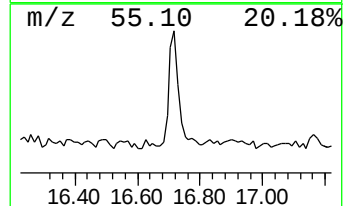
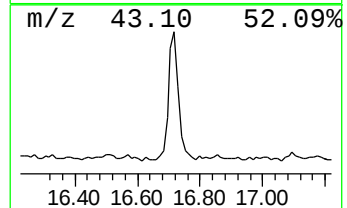
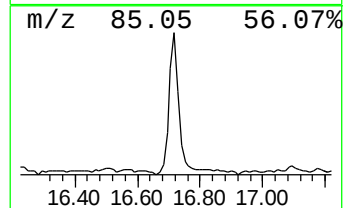
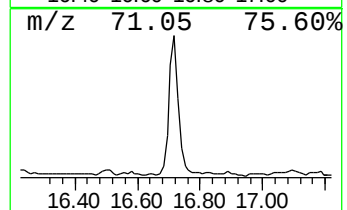
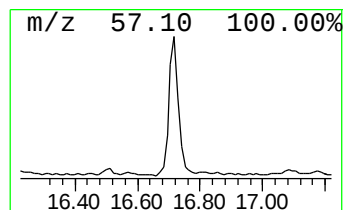
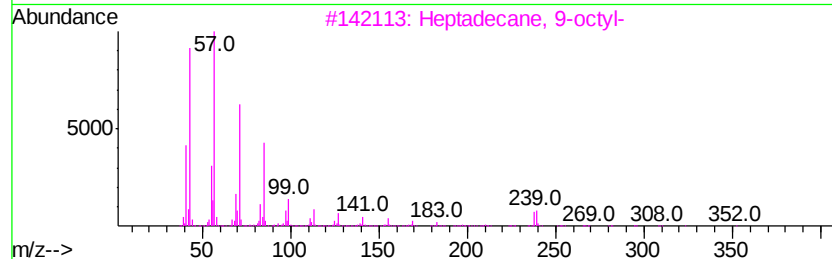
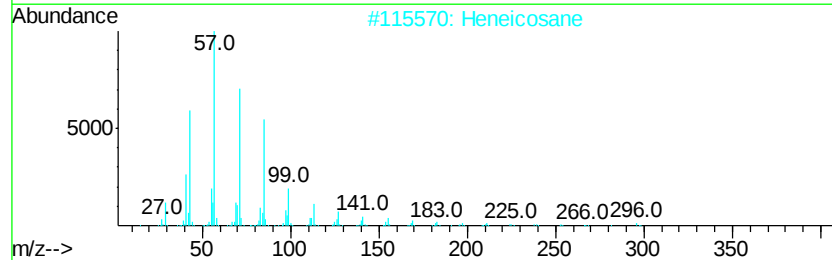
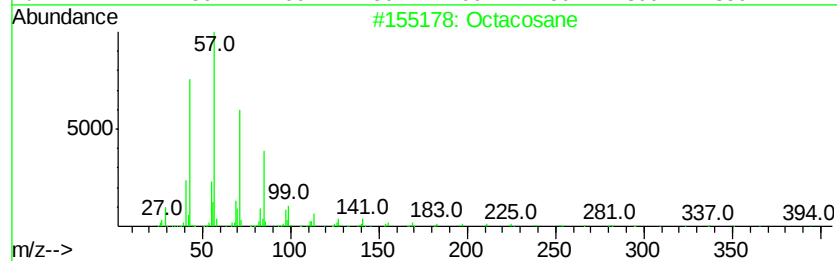
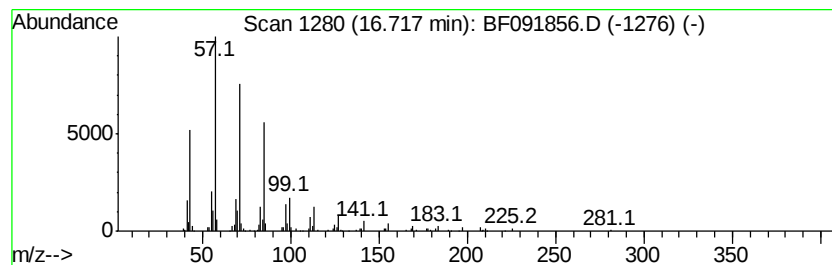
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Octacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.72	3.33 ng	332060	Perylene-d12	15.29

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	91
2			Heneicosane	296	C21H44	000629-94-7	91
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4			Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
5			Docosane	310	C22H46	000629-97-0	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122116\
Data File : BF091856.D
Acq On : 22 Dec 2016 1:40
Operator : UM/SJ
Sample : H6156-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-1101

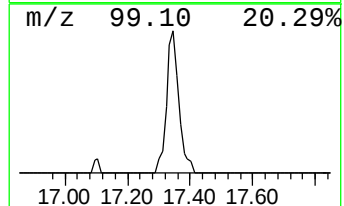
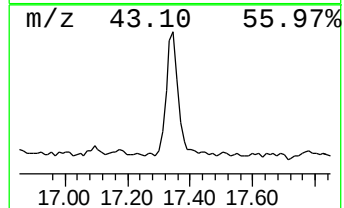
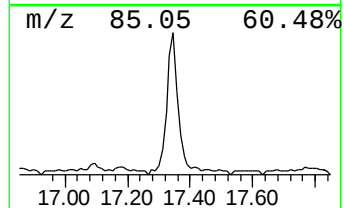
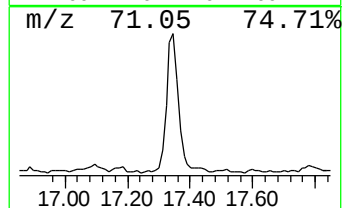
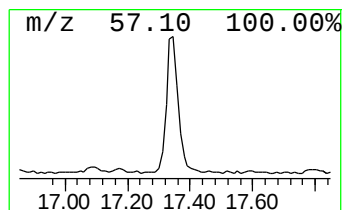
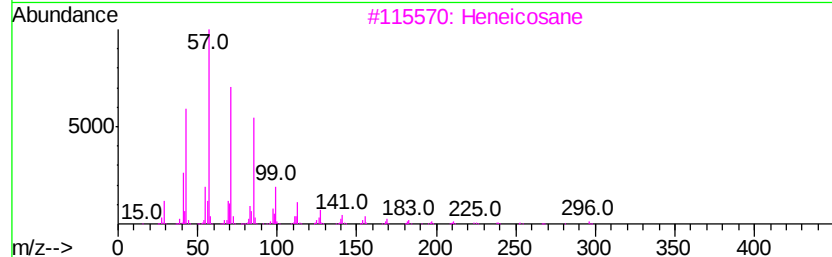
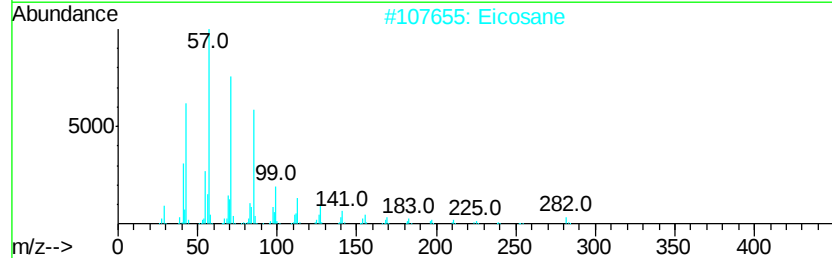
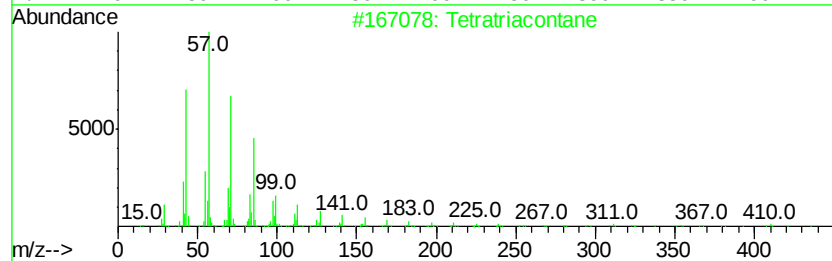
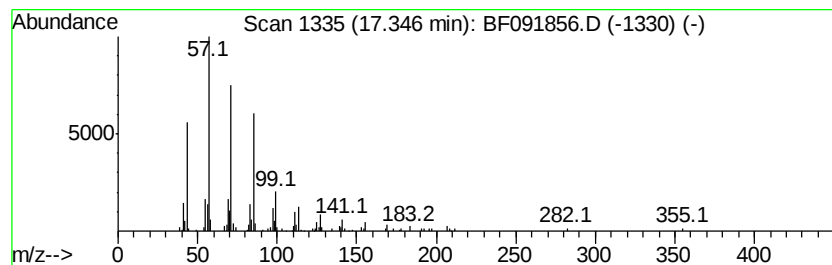
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetratriacontane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.35	2.91 ng	290536	Perylene-d12	15.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratriacontane	479	C34H70	014167-59-0	91
2		Eicosane	282	C20H42	000112-95-8	91
3		Heneicosane	296	C21H44	000629-94-7	90
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90
5		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122116\
Data File : BF091856.D
Acq On : 22 Dec 2016 1:40
Operator : UM/SJ
Sample : H6156-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.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
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unknown6.51	6.51	79.6	ng	7597090	1	6.74	1907980	20.0
Benzene, 1-butynyl-	7.81	2.7	ng	385756	2	8.03	2816030	20.0
1(2H)-Acenaphthyl...	10.69	2.3	ng	346642	4	11.27	3025180	20.0
2-Dibenzofuranol	11.14	2.1	ng	311209	4	11.27	3025180	20.0
Eicosane	15.72	3.6	ng	354923	6	15.29	1995080	20.0
Heneicosane	16.18	4.1	ng	404697	6	15.29	1995080	20.0
Octacosane	16.72	3.3	ng	332060	6	15.29	1995080	20.0
Tetratriacontane	17.35	2.9	ng	290536	6	15.29	1995080	20.0