

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032516\
 Data File : BF085789.D
 Acq On : 26 Mar 2016 4:21
 Operator : UM/SJ
 Sample : H1923-02
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-PG-B(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-BF032416.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	3.784	129	131	133	rVB2	36250	31066	1.04%	0.108%
2	3.864	135	138	140	rVB2	38955	51982	1.73%	0.181%
3	5.007	235	238	246	rBV	760684	1089797	36.34%	3.794%
4	5.430	272	275	278	rBV	2553838	2762041	92.11%	9.615%
5	5.830	307	310	314	rVB	97524	100407	3.35%	0.350%
6	6.470	363	366	368	rBV	2579486	2998495	100.00%	10.438%
7	6.596	374	377	379	rBV	3024570	2887823	96.31%	10.053%
8	6.825	394	397	399	rBV	761159	646924	21.57%	2.252%
9	6.985	408	411	413	rVB	2354716	2429804	81.03%	8.458%
10	7.396	444	447	449	rBV	1836726	2072239	69.11%	7.214%
11	7.499	452	456	460	rVB	30697	51535	1.72%	0.179%
12	7.853	483	487	492	rBV2	20639	38875	1.30%	0.135%
13	8.105	507	509	512	rBV	633932	749074	24.98%	2.608%
14	9.191	601	604	606	rBV	2718888	2847114	94.95%	9.911%
15	9.579	636	638	642	rVB	433883	364780	12.17%	1.270%
16	9.796	655	657	659	rBV	75071	58132	1.94%	0.202%
17	9.865	660	663	665	rVV	737345	711559	23.73%	2.477%
18	10.299	697	701	702	rBV	70307	110539	3.69%	0.385%
19	10.516	717	720	723	rBV	34372	60651	2.02%	0.211%
20	10.653	729	732	734	rBV	1626303	1608316	53.64%	5.599%
21	10.768	740	742	747	rBV	106157	124493	4.15%	0.433%
22	10.962	757	759	763	rBV3	15088	36573	1.22%	0.127%
23	11.214	779	781	782	rBV	65999	54196	1.81%	0.189%
24	11.339	790	792	796	rBV2	554585	723888	24.14%	2.520%
25	11.419	796	799	801	rVB2	23552	39265	1.31%	0.137%
26	11.579	809	813	816	rBV4	21406	31661	1.06%	0.110%
27	11.876	837	839	841	rBV	99344	99665	3.32%	0.347%
28	11.956	844	846	850	rVB2	40239	59043	1.97%	0.206%
29	12.551	896	898	900	rBV	233101	230202	7.68%	0.801%
30	12.654	905	907	910	rBV2	26005	41984	1.40%	0.146%
31	12.745	913	915	916	rBV	60896	77472	2.58%	0.270%
32	12.779	916	918	921	rVB	228359	244023	8.14%	0.849%
33	12.928	928	931	933	rBV	2123076	1910183	63.70%	6.650%
34	13.111	944	947	949	rVB	28871	39131	1.31%	0.136%

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35	13.157	949	951	954	rBV2	60308	84619	2.82%	0.295%
36	13.454	975	977	979	rBV	52923	52133	1.74%	0.181%
37	13.499	979	981	982	rVV	36476	39481	1.32%	0.137%
38	13.534	982	984	985	rVB	24994	31761	1.06%	0.111%
39	13.614	990	991	996	rVB	22160	35159	1.17%	0.122%
40	13.728	998	1001	1002	rBV	36506	47778	1.59%	0.166%
41	13.774	1002	1005	1007	rVV3	34239	74679	2.49%	0.260%
42	13.831	1007	1010	1014	rVB2	117163	185873	6.20%	0.647%
43	13.980	1019	1023	1024	rBV2	706524	832771	27.77%	2.899%
44	14.082	1029	1032	1033	rBV	33586	44984	1.50%	0.157%
45	14.140	1035	1037	1039	rBV2	40024	60906	2.03%	0.212%
46	14.448	1062	1064	1066	rBV3	58698	88286	2.94%	0.307%
47	14.608	1076	1078	1081	rBV	29378	37388	1.25%	0.130%
48	14.745	1088	1090	1091	rVB	32789	30103	1.00%	0.105%
49	14.997	1109	1112	1115	rBV	111251	206547	6.89%	0.719%
50	15.065	1115	1118	1119	rBV	139493	173631	5.79%	0.604%
51	15.283	1135	1137	1140	rVB	59197	71828	2.40%	0.250%
52	15.340	1140	1142	1145	rVB	75697	91330	3.05%	0.318%
53	15.397	1145	1147	1156	rBV	402293	573593	19.13%	1.997%
54	15.614	1163	1166	1167	rBV2	38567	66719	2.23%	0.232%
55	15.831	1182	1185	1187	rBV	72418	108411	3.62%	0.377%
56	16.048	1202	1204	1206	rBV2	23519	42005	1.40%	0.146%
57	16.300	1223	1226	1230	rBV2	30241	65055	2.17%	0.226%
58	16.574	1248	1250	1252	rBV2	22443	38007	1.27%	0.132%
59	16.780	1266	1268	1271	rVB2	35588	64364	2.15%	0.224%
60	16.848	1271	1274	1279	rVB	38691	79953	2.67%	0.278%
61	17.008	1286	1288	1289	rBV2	21818	31326	1.04%	0.109%
62	17.191	1302	1304	1310	rVB	39507	84903	2.83%	0.296%

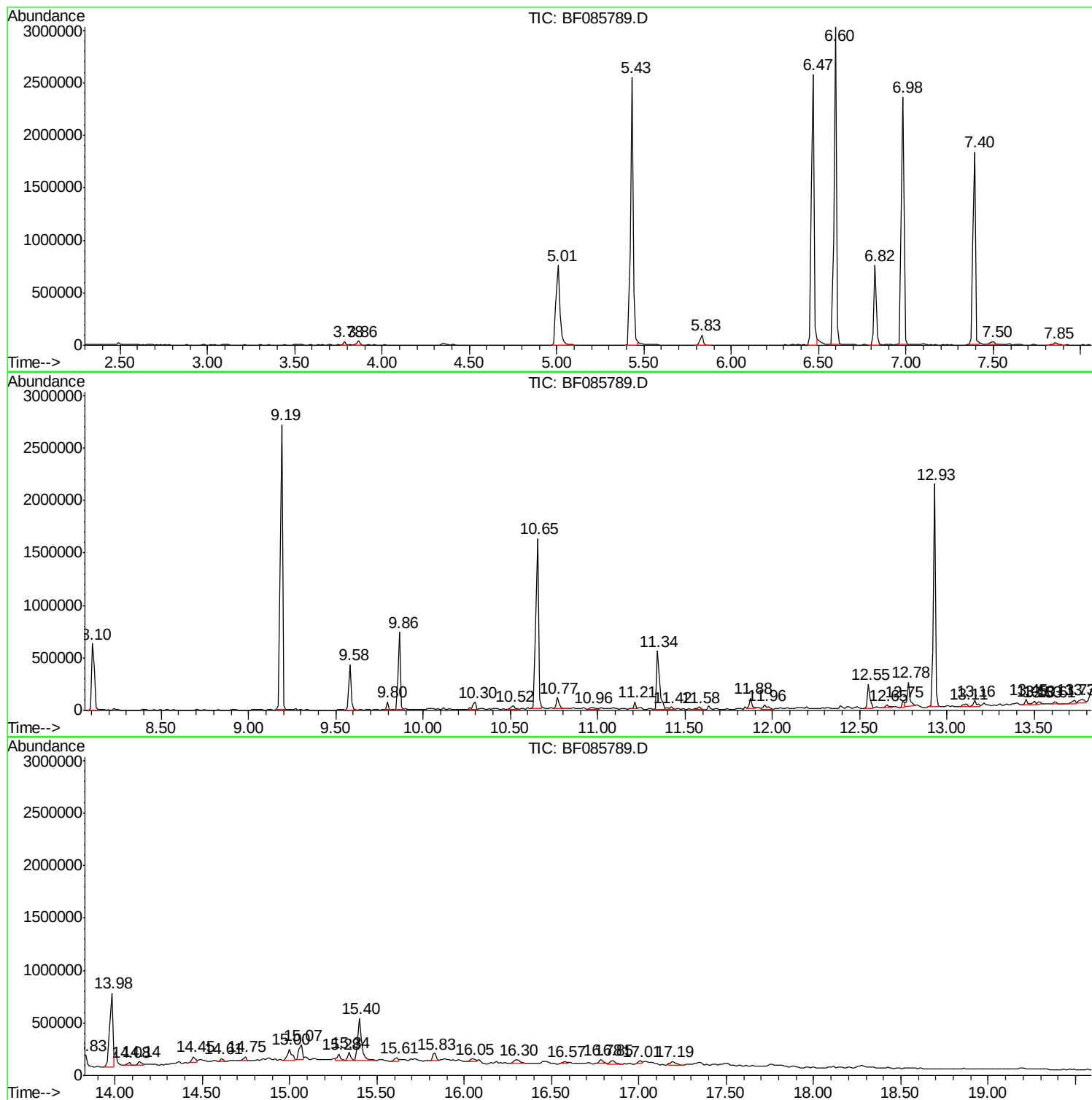
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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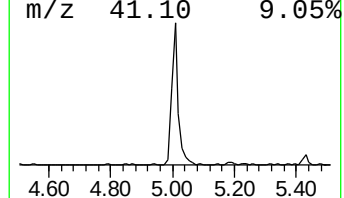
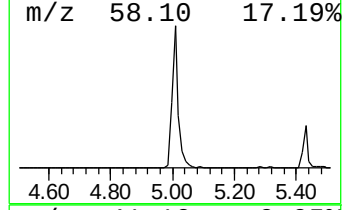
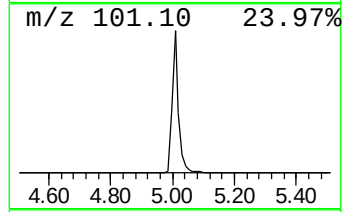
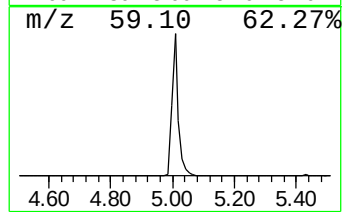
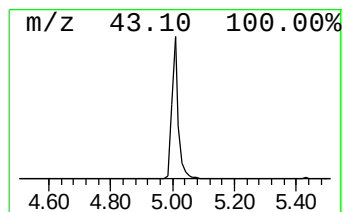
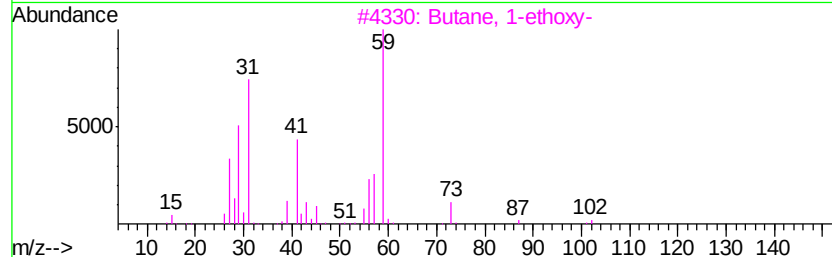
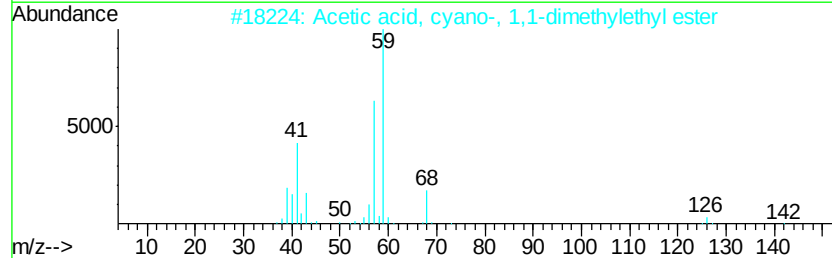
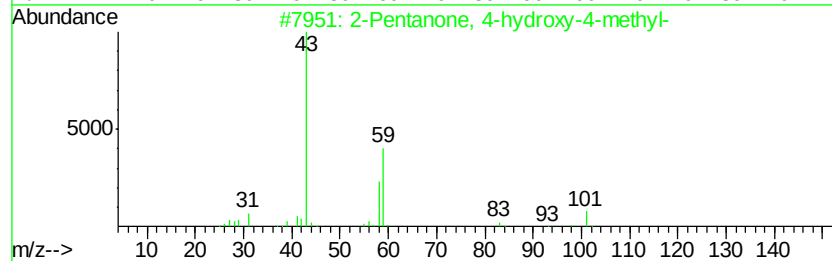
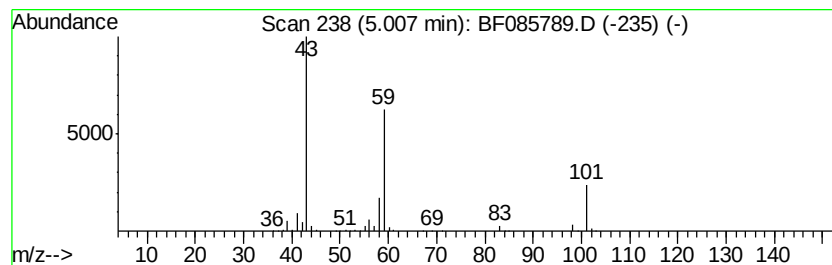
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	33.69 ng	1089800	1,4-Dichlorobenzene-d4	6.82

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	25
3		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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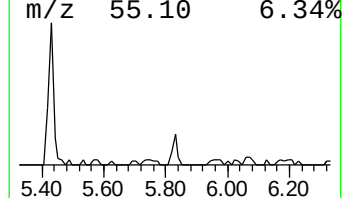
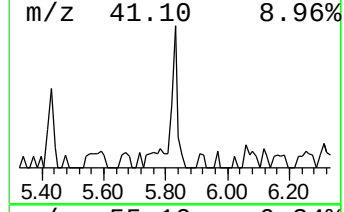
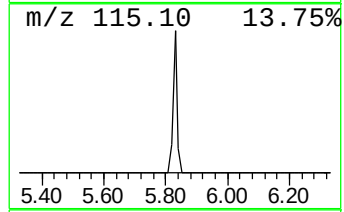
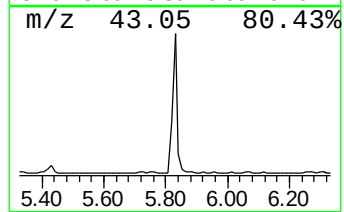
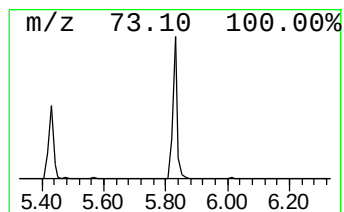
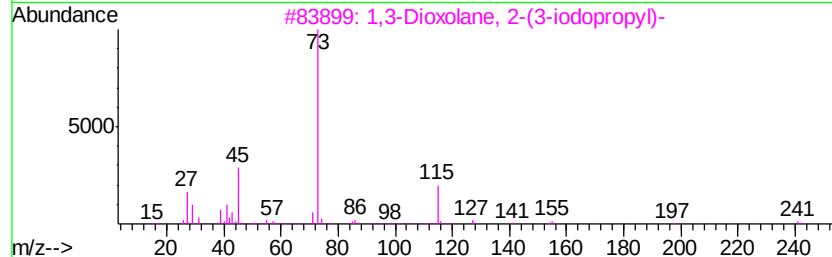
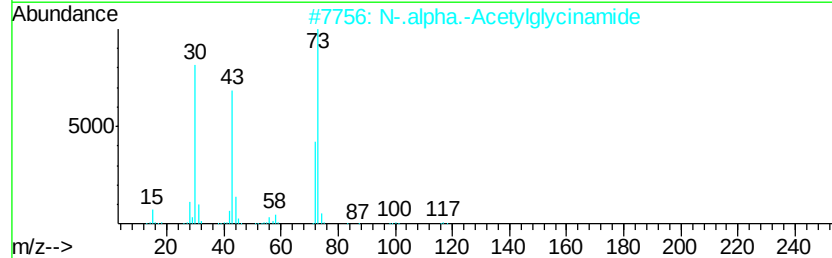
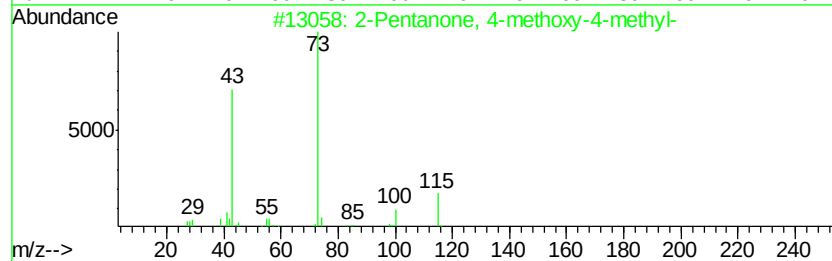
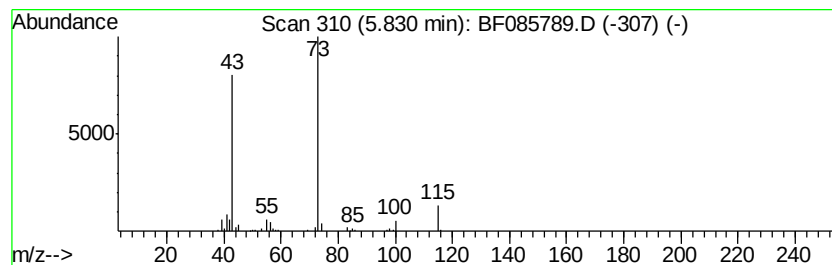
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 2-Pentanone, 4-methoxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.83	3.10 ng	100407	1,4-Dichlorobenzene-d4	6.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	83
2			N-.alpha.-Acetylglucinamide	116	C4H8N2O2	002620-63-5	28
3			1,3-Dioxolane, 2-(3-iodopropyl)-	242	C6H11IO2	058135-25-4	28
4			Butanamide, N-formyl-2-hydroxy-3...	187	C9H17NO3	056440-42-7	25
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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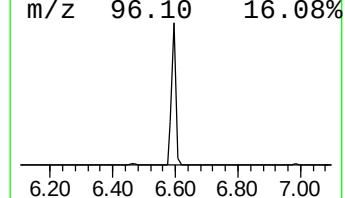
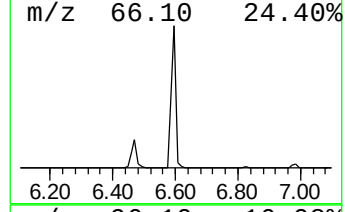
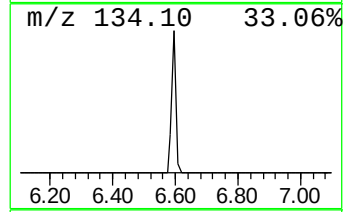
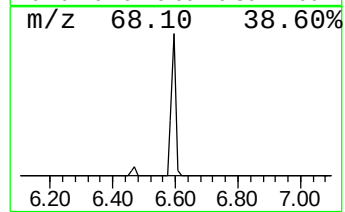
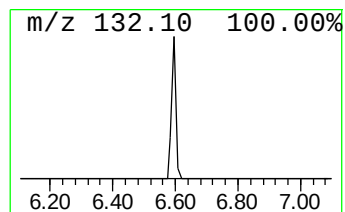
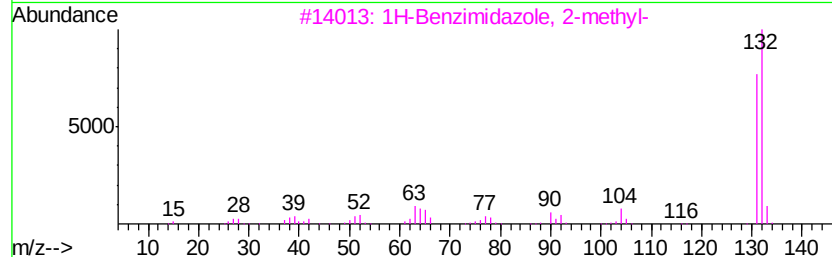
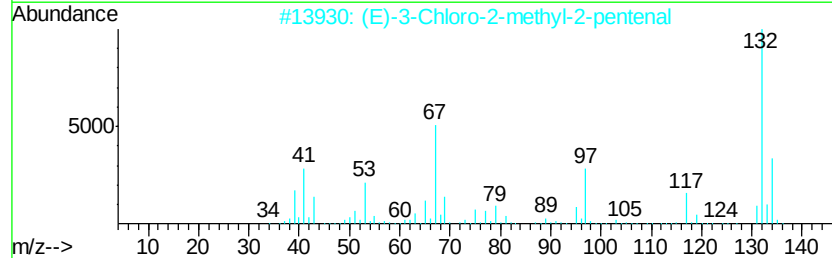
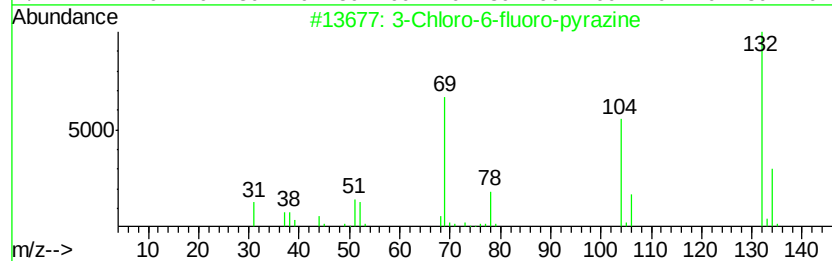
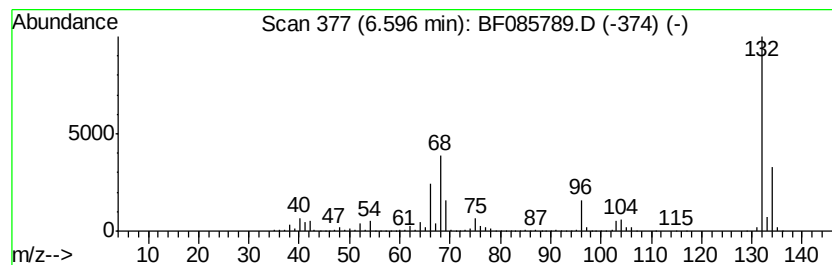
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 unknown6.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	89.28 ng	2887820	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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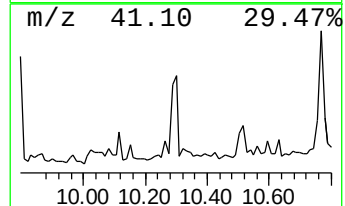
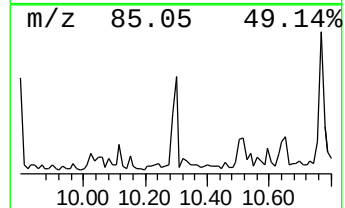
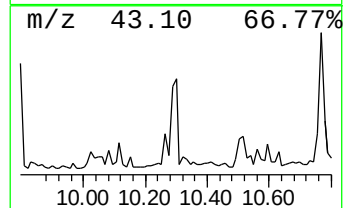
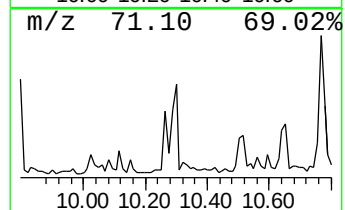
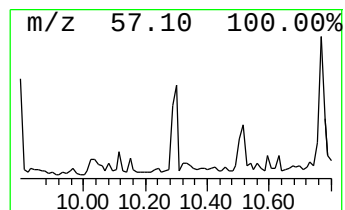
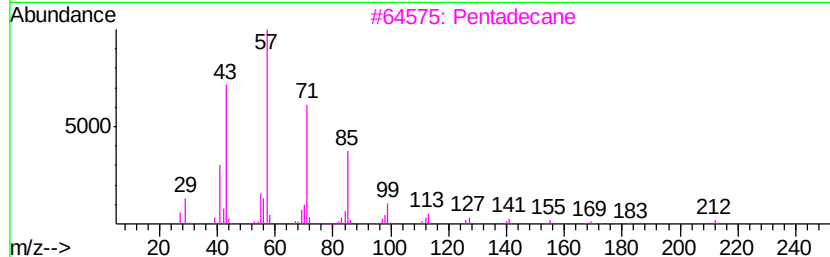
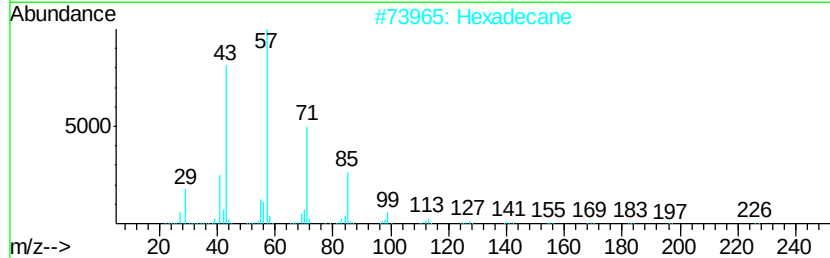
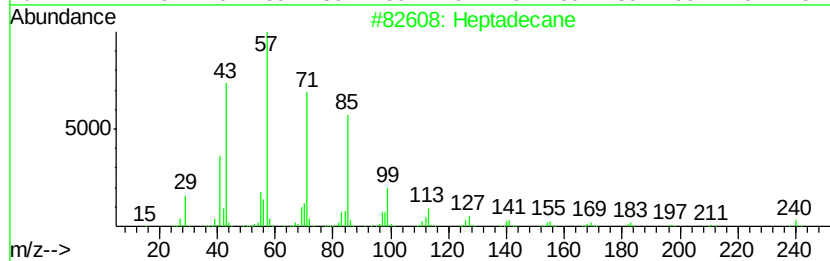
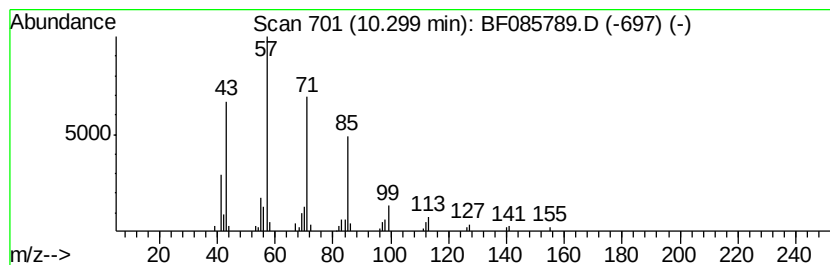
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Peak Number 5 Heptadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.30	3.11 ng	110539	Acenaphthene-d10		9.86
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	91
2	Hexadecane	226	C16H34	000544-76-3	90
3	Pentadecane	212	C15H32	000629-62-9	87
4	Tetradecane	198	C14H30	000629-59-4	87
5	Heneicosane	296	C21H44	000629-94-7	86



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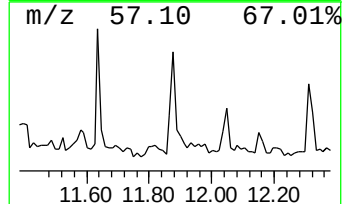
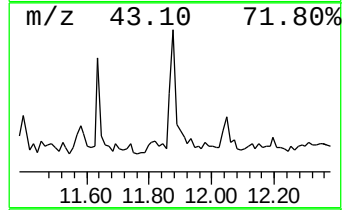
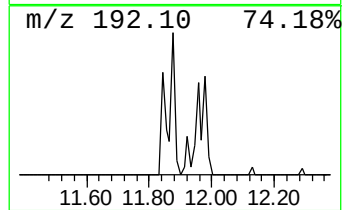
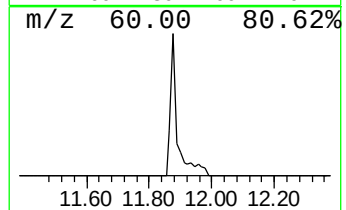
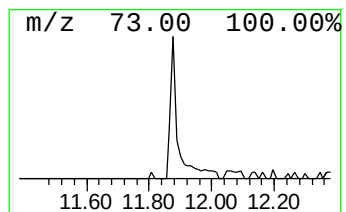
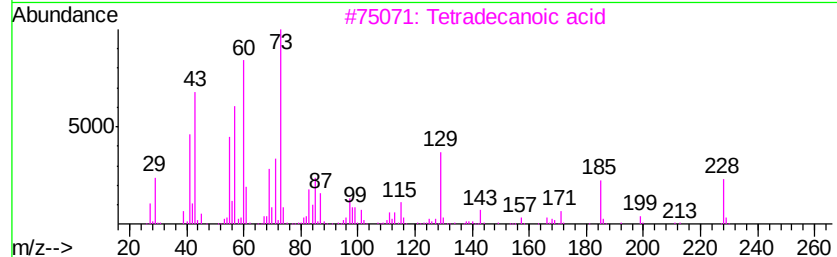
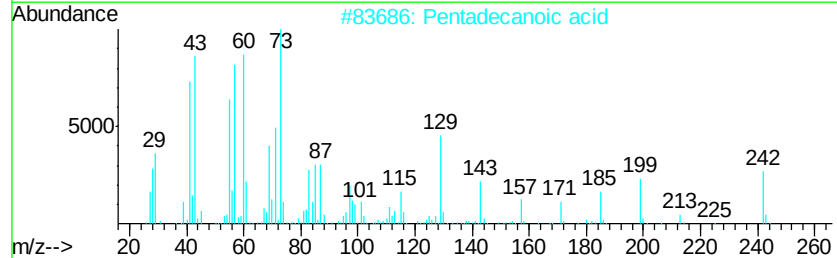
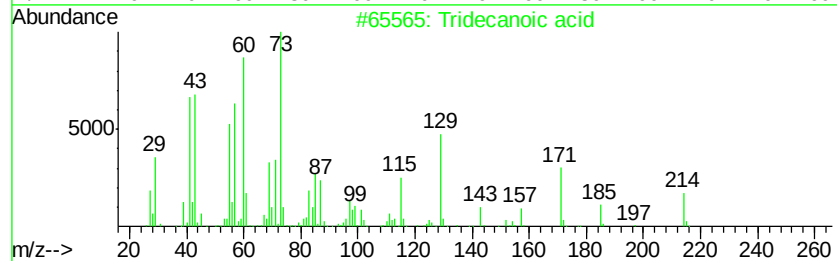
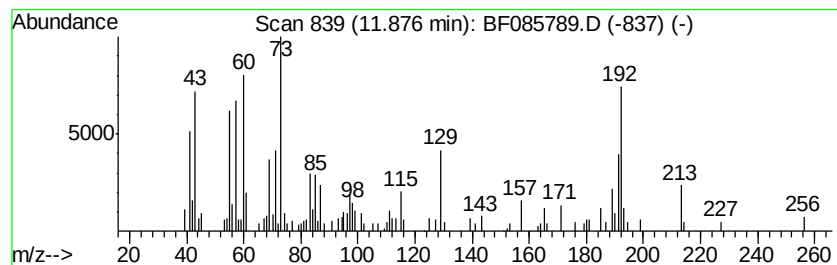
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S-PG-B(0-6)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tridecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.88	2.75 ng	99665	Phenanthrene-d10		11.34
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanoic acid	214	C13H26O2	000638-53-9	70
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	49
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	49
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	45
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	42



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032516\
Data File : BF085789.D
Acq On : 26 Mar 2016 4:21
Operator : UM/SJ
Sample : H1923-02
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
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ClientSampleId :
S-PG-B(0-6)

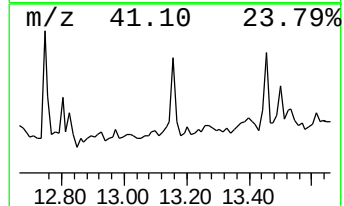
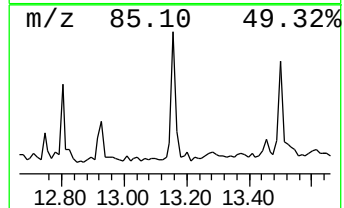
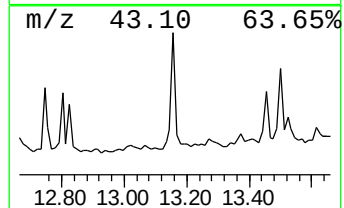
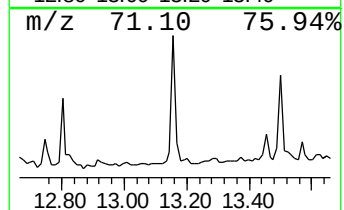
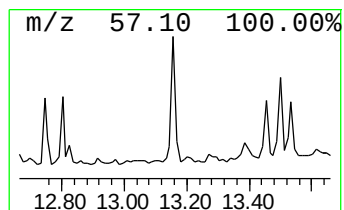
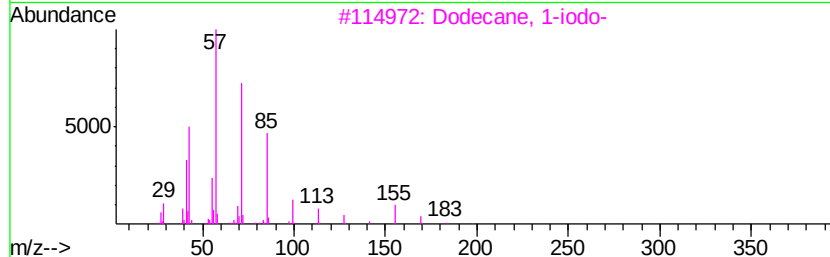
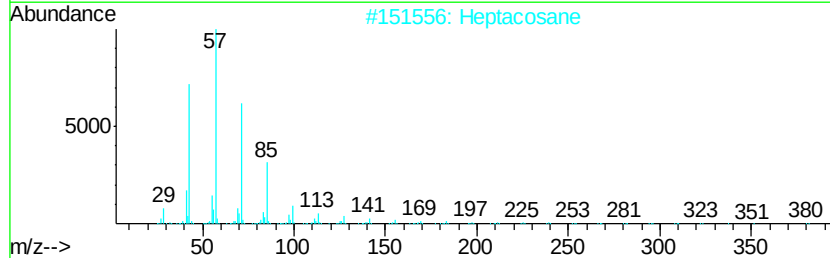
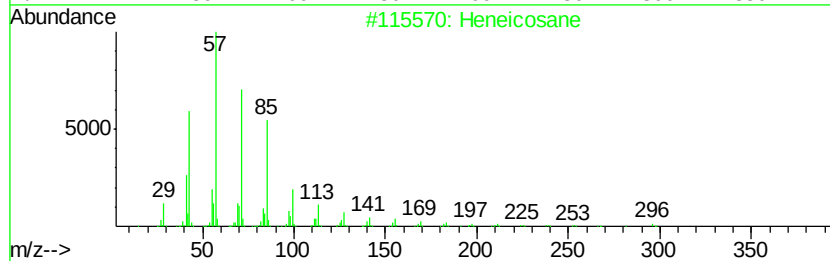
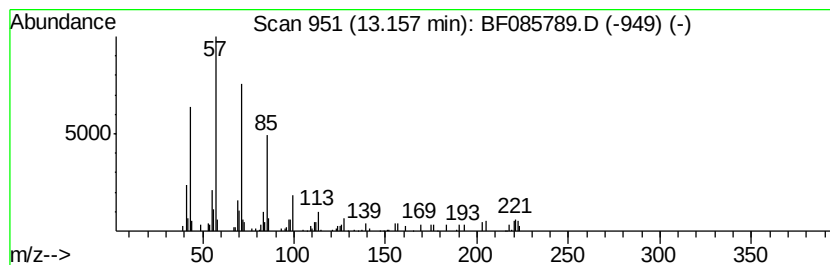
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heneicosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.16	2.03 ng	84619	Chrysene-d12	13.98

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	83
2	Heptacosane		380	C27H56	000593-49-7	80
3	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	80
4	Nonacosane		408	C29H60	000630-03-5	80
5	Tetracosane		338	C24H50	000646-31-1	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032516\
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Sample : H1923-02
Misc :
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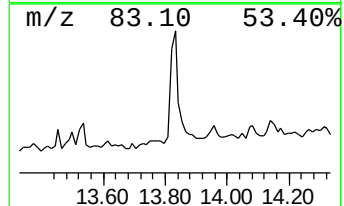
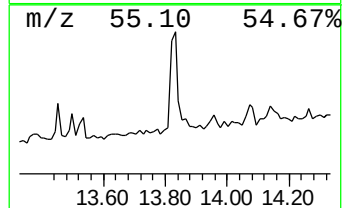
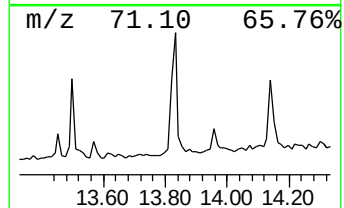
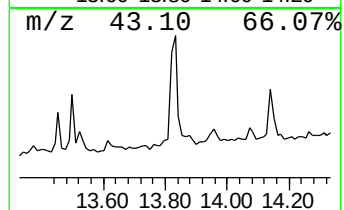
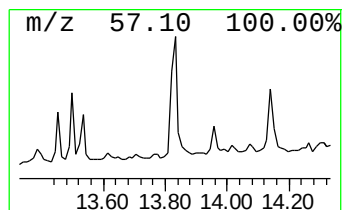
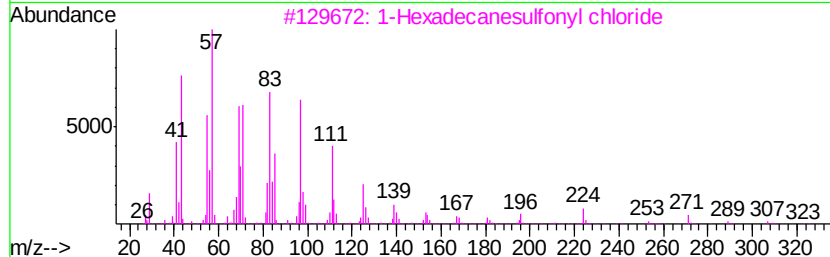
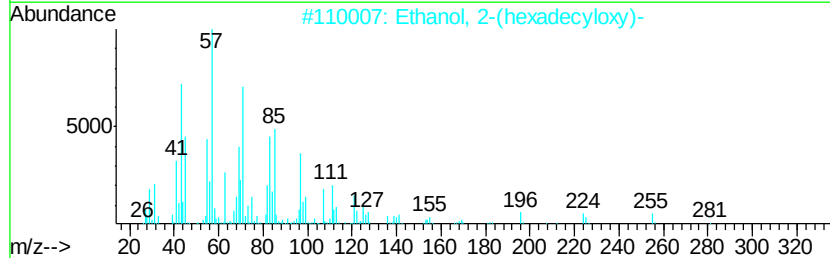
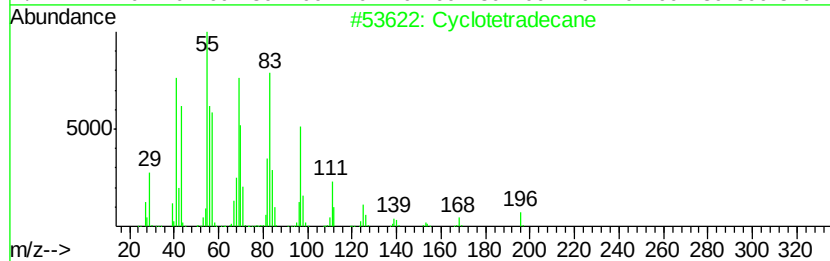
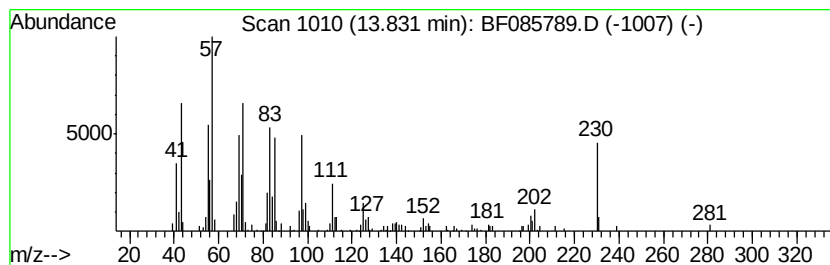
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Cyclotetradecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.83	4.46 ng	185873	Chrysene-d12	13.98	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetradecane	196	C14H28	000295-17-0	81
2	Ethanol, 2-(hexadecyloxy)-	286	C18H38O2	002136-71-2	74
3	1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	62
4	Heptafluorobutanoic acid, hepta...	452	C21H35F7O2	1000282-97-3	60
5	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	60



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032516\
Data File : BF085789.D
Acq On : 26 Mar 2016 4:21
Operator : UM/SJ
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Instrument :
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ClientSampleId :
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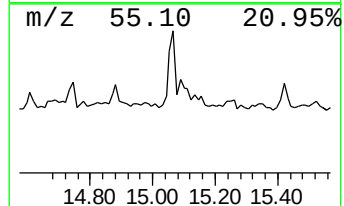
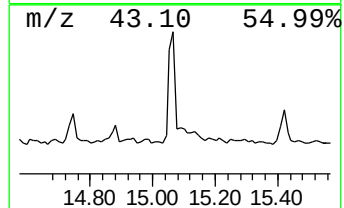
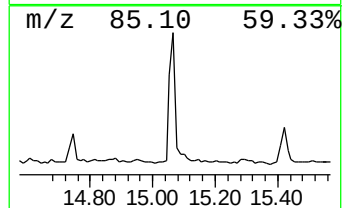
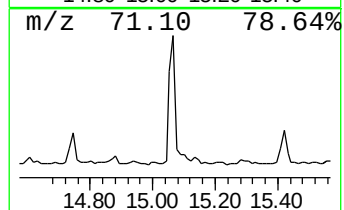
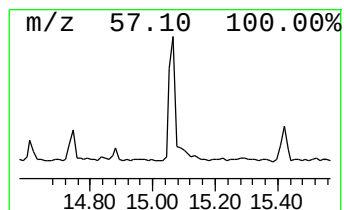
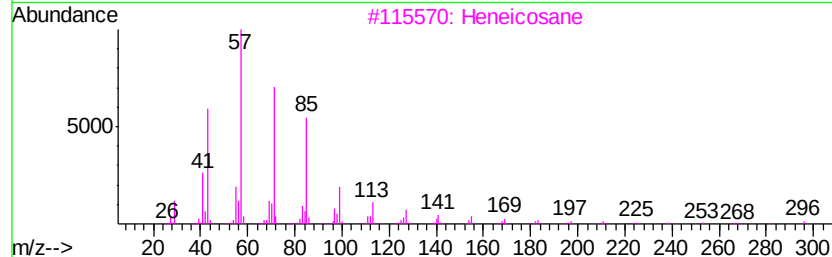
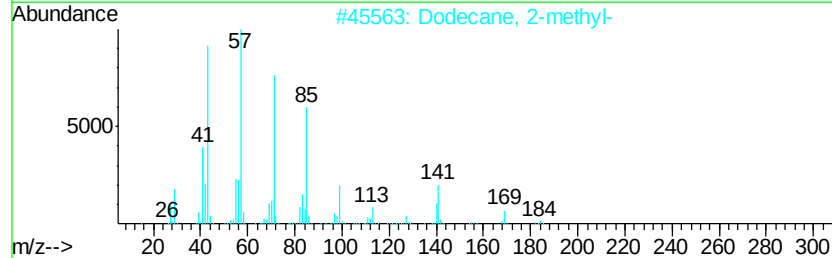
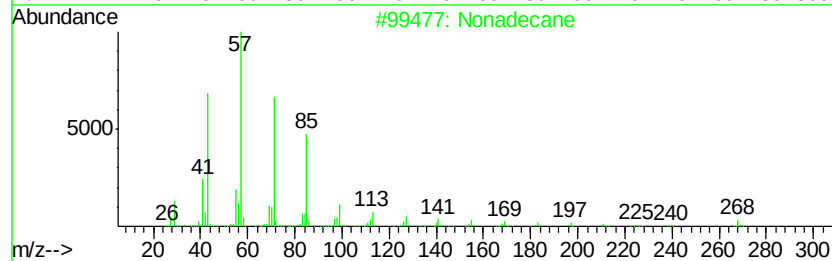
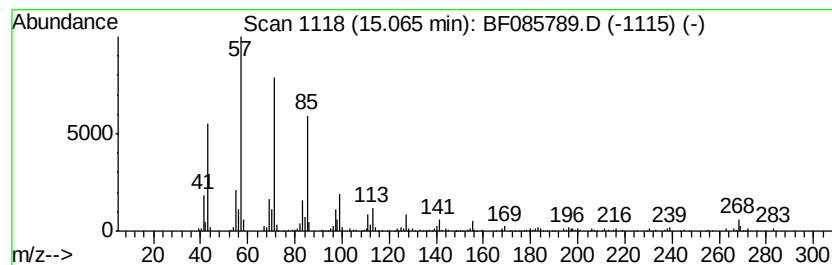
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Nonadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	6.05 ng	173631	Perylene-d12	15.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	91
2		Dodecane, 2-methyl-	184	C13H28	001560-97-0	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	87
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032516\
Data File : BF085789.D
Acq On : 26 Mar 2016 4:21
Operator : UM/SJ
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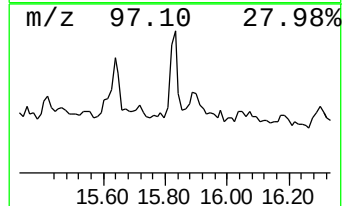
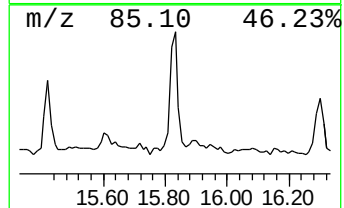
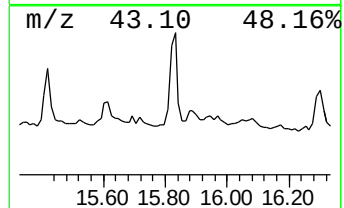
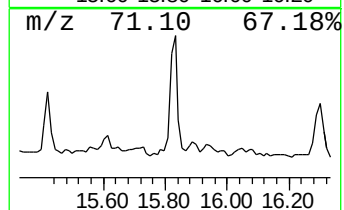
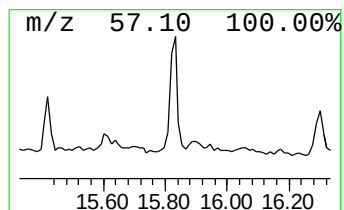
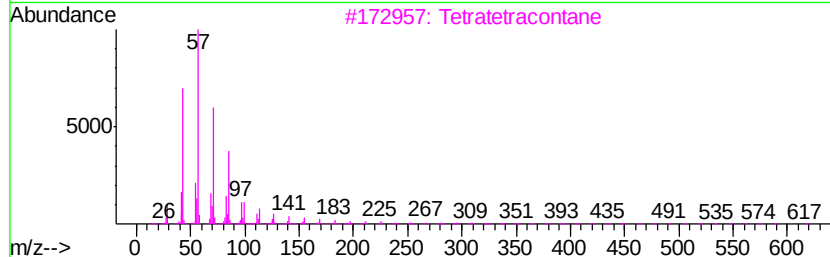
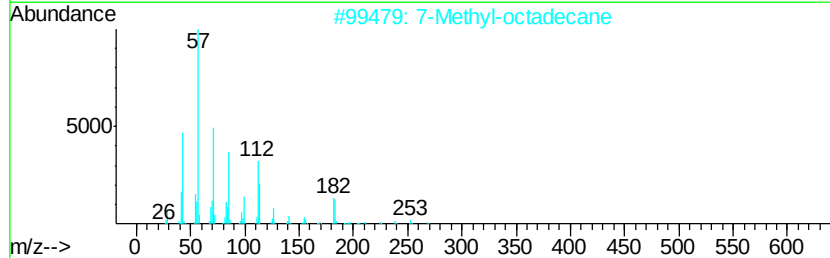
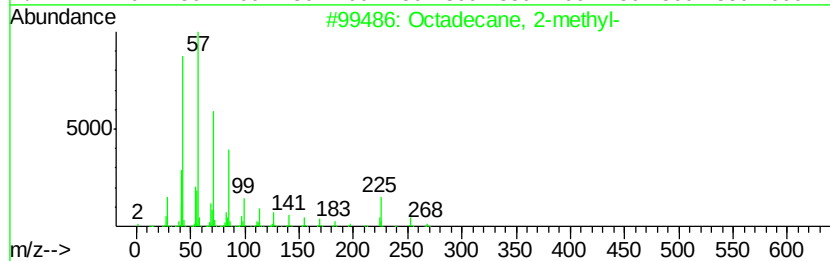
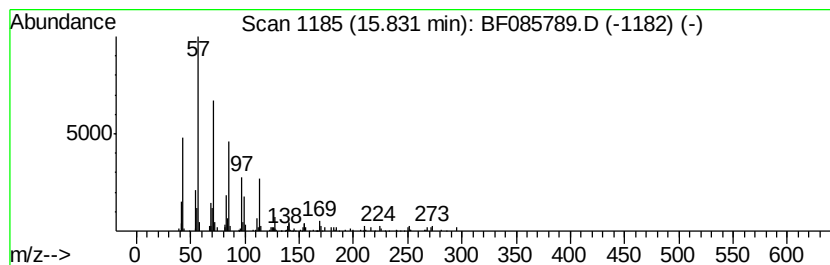
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Octadecane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.83	3.78 ng	108411	Perylene-d12	15.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane, 2-methyl-	268	C19H40	001560-88-9	81
2		7-Methyl-octadecane	268	C19H40	1000192-63-3	72
3		Tetratetracontane	619	C44H90	007098-22-8	64
4		Nonadecane	268	C19H40	000629-92-5	64
5		Eicosane	282	C20H42	000112-95-8	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032516\
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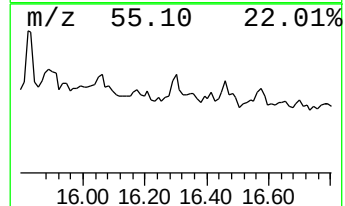
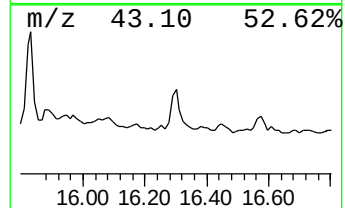
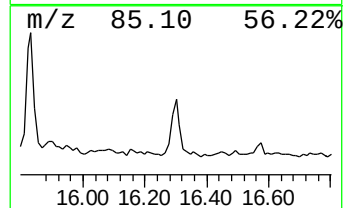
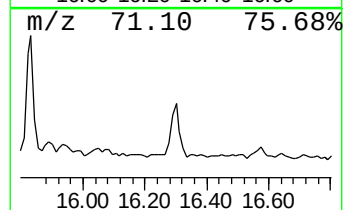
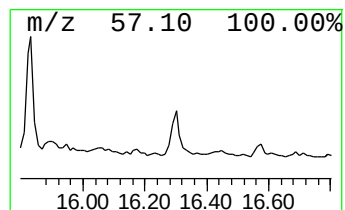
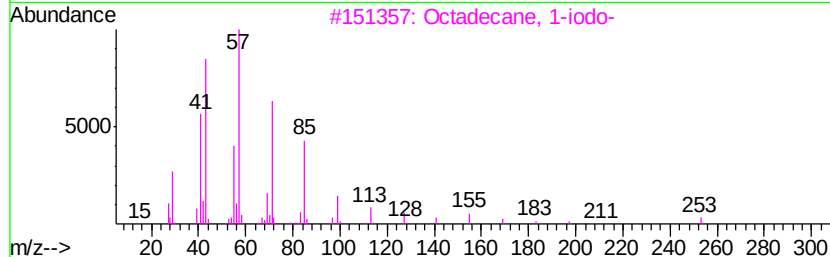
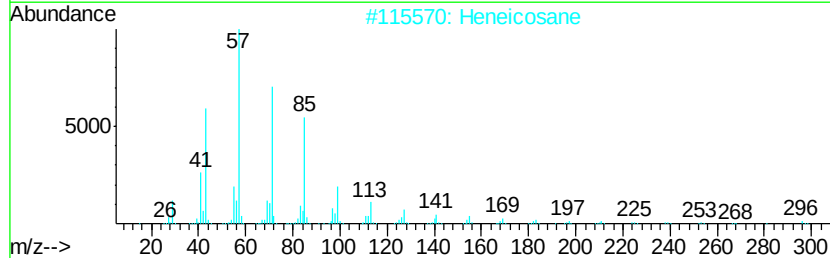
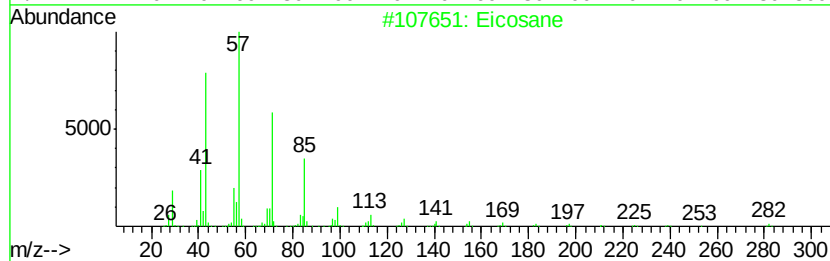
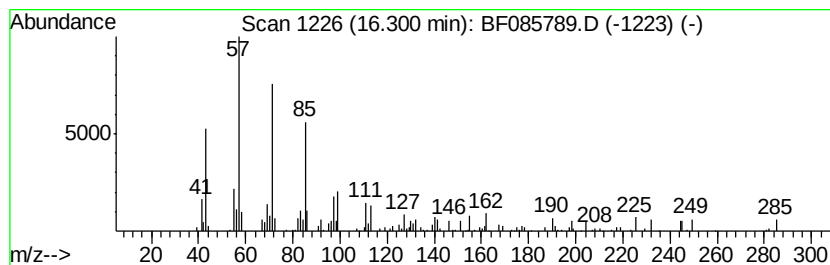
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Eicosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.30	2.27 ng	65055	Perylene-d12	15.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	76
2	Heneicosane	296 C21H44	000629-94-7	74
3	Octadecane, 1-iodo-	380 C18H37I	000629-93-6	72
4	Pentadecane, 8-heptyl-	310 C22H46	071005-15-7	72
5	Tridecanol, 2-ethyl-2-methyl-	242 C16H34O	1000115-66-1	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032516\
Data File : BF085789.D
Acq On : 26 Mar 2016 4:21
Operator : UM/SJ
Sample : H1923-02
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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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2-Pentanone, 4-me...	5.83	3.1	ng	100407	1	6.82	646924	20.0
unknown6.60	6.60	89.3	ng	2887820	1	6.82	646924	20.0
Heptadecane	10.30	3.1	ng	110539	3	9.86	711559	20.0
Tridecanoic acid	11.88	2.8	ng	99665	4	11.34	723888	20.0
Heneicosane	13.16	2.0	ng	84619	5	13.98	832771	20.0
Cyclotetradecane	13.83	4.5	ng	185873	5	13.98	832771	20.0
Nonadecane	15.07	6.0	ng	173631	6	15.40	573593	20.0
Octadecane, 2-met...	15.83	3.8	ng	108411	6	15.40	573593	20.0
Eicosane	16.30	2.3	ng	65055	6	15.40	573593	20.0