

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040216\
 Data File : BF086007.D
 Acq On : 2 Apr 2016 13:05
 Operator : UM/SJ
 Sample : H2055-09
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BH-9(5-7)

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-BF040216.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	4.967	250	252	260	rBV	51702	83655	2.87%	0.285%
2	5.367	285	287	290	rBV	2052104	2350151	80.74%	8.012%
3	6.419	376	379	381	rBV	2318350	2799546	96.17%	9.545%
4	6.544	387	390	392	rBV	2430467	2828065	97.15%	9.642%
5	6.773	408	410	416	rBV	703928	644781	22.15%	2.198%
6	6.921	421	423	426	rBV	1649277	2010668	69.07%	6.855%
7	7.344	458	460	462	rBV	1555940	1850916	63.58%	6.310%
8	8.053	520	522	532	rVB	880528	931232	31.99%	3.175%
9	8.773	584	585	588	rBV3	56531	52676	1.81%	0.180%
10	8.876	591	594	597	rBV	60814	97134	3.34%	0.331%
11	9.082	604	612	614	rBV5	15108	53428	1.84%	0.182%
12	9.139	614	617	619	rBV	1958119	2764181	94.96%	9.424%
13	9.310	630	632	636	rBV3	21156	51036	1.75%	0.174%
14	9.402	638	640	644	rVB2	29659	44248	1.52%	0.151%
15	9.470	644	646	647	rBV	49534	53533	1.84%	0.183%
16	9.527	650	651	654	rVV	444492	421845	14.49%	1.438%
17	9.619	655	659	661	rVB4	12986	32559	1.12%	0.111%
18	9.676	661	664	668	rBV4	22284	44803	1.54%	0.153%
19	9.745	668	670	672	rBV	72242	62239	2.14%	0.212%
20	9.813	673	676	678	rBV	717809	975175	33.50%	3.325%
21	9.973	688	690	693	rBV3	30628	61952	2.13%	0.211%
22	10.167	706	707	709	rBV2	22791	30146	1.04%	0.103%
23	10.236	712	713	716	rVB	56139	77619	2.67%	0.265%
24	10.465	730	733	736	rVB	33353	59084	2.03%	0.201%
25	10.602	742	745	747	rBV	1910019	2082557	71.54%	7.100%
26	10.716	753	755	759	rVB	125093	169391	5.82%	0.578%
27	10.910	769	772	774	rBV3	18848	42317	1.45%	0.144%
28	11.036	782	783	785	rBV2	28797	34830	1.20%	0.119%
29	11.162	792	794	795	rBV	78929	67497	2.32%	0.230%
30	11.288	803	805	810	rBV	1011618	1132086	38.89%	3.860%
31	11.368	810	812	819	rVB5	24461	69850	2.40%	0.238%
32	11.825	850	852	858	rVV2	192276	218262	7.50%	0.744%
33	11.905	858	859	863	rVB2	32500	53238	1.83%	0.182%
34	12.373	899	900	904	rVB4	27093	44806	1.54%	0.153%

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35	12.511	909	912	917	rVB	88477	143423	4.93%	0.489%
36	12.602	918	920	923	rBV2	60964	91791	3.15%	0.313%
37	12.693	927	928	930	rBV	109510	132600	4.56%	0.452%
38	12.728	930	931	934	rVB2	91884	125184	4.30%	0.427%
39	12.876	941	944	946	rBV	3105470	2910944	100.00%	9.924%
40	13.059	957	960	962	rVB2	14460	33291	1.14%	0.113%
41	13.105	962	964	966	rBV	85987	84076	2.89%	0.287%
42	13.356	983	986	988	rVB	124603	116157	3.99%	0.396%
43	13.402	988	990	992	rVB	112227	96993	3.33%	0.331%
44	13.448	992	994	995	rBV	89440	85602	2.94%	0.292%
45	13.779	1020	1023	1032	rVB	217905	404553	13.90%	1.379%
46	13.928	1033	1036	1038	rBV	1007028	1087994	37.38%	3.709%
47	14.088	1048	1050	1052	rBV	70993	71563	2.46%	0.244%
48	14.385	1074	1076	1078	rBV	62968	116420	4.00%	0.397%
49	14.419	1078	1079	1084	rVV2	83451	163896	5.63%	0.559%
50	14.499	1084	1086	1094	rVB2	21869	34607	1.19%	0.118%
51	14.682	1100	1102	1105	rBV	42317	48463	1.66%	0.165%
52	14.945	1123	1125	1128	rBV	37463	74969	2.58%	0.256%
53	14.991	1128	1129	1132	rVB	39230	46982	1.61%	0.160%
54	15.048	1132	1134	1140	rBV3	16192	38719	1.33%	0.132%
55	15.277	1152	1154	1156	rBV	26278	33100	1.14%	0.113%
56	15.334	1156	1159	1166	rVB	600692	842052	28.93%	2.871%
57	15.734	1192	1194	1200	rBV	25526	41276	1.42%	0.141%
58	16.191	1231	1234	1241	rVB2	18516	39345	1.35%	0.134%
59	16.728	1276	1281	1287	rBV2	22309	57481	1.97%	0.196%
60	17.357	1331	1336	1342	rVB4	14147	44060	1.51%	0.150%
61	18.100	1396	1401	1408	rBV2	11761	37800	1.30%	0.129%
62	18.980	1474	1478	1489	rVB4	7460	32654	1.12%	0.111%

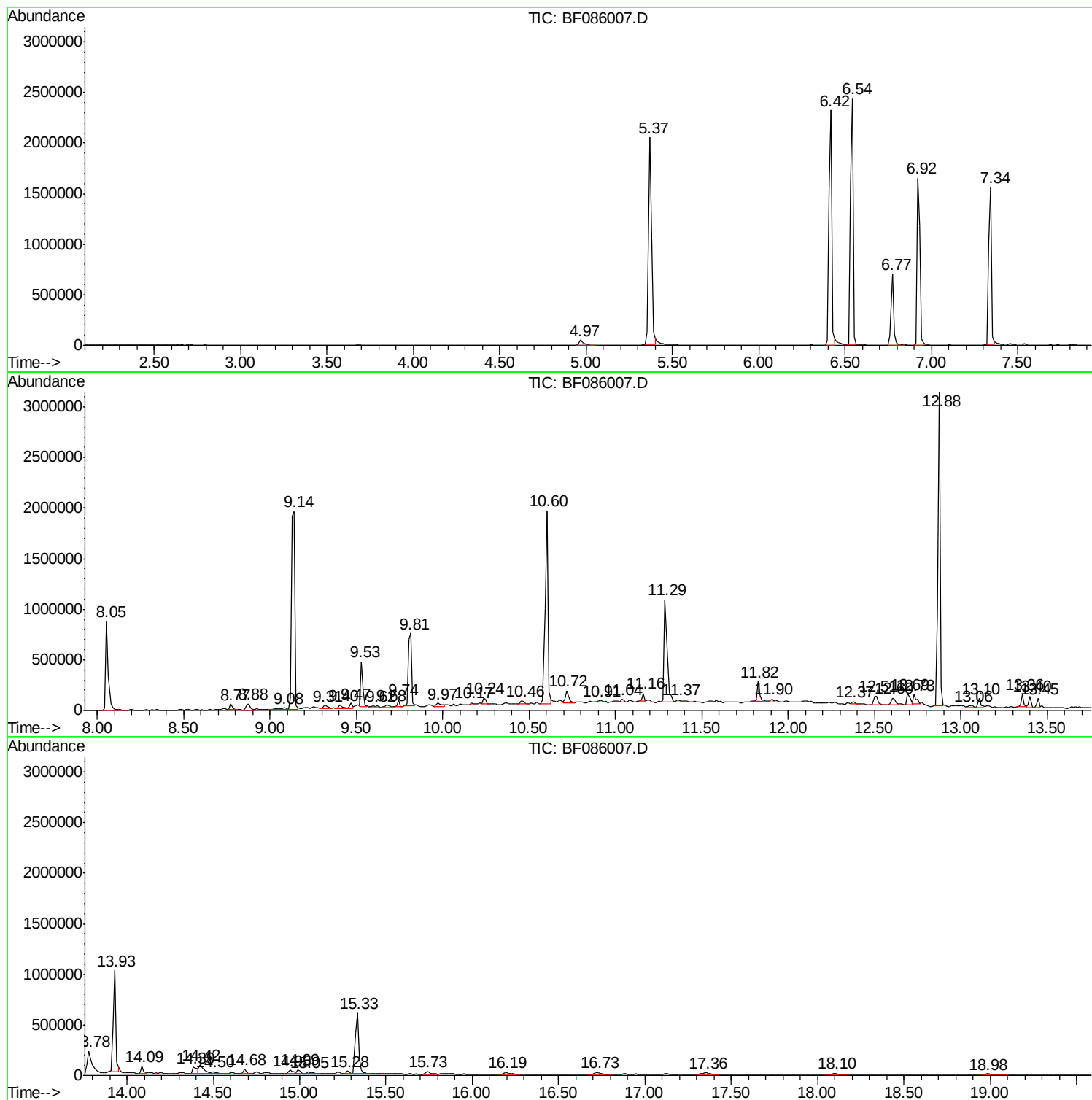
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ClientSampled :
BH-9(5-7)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.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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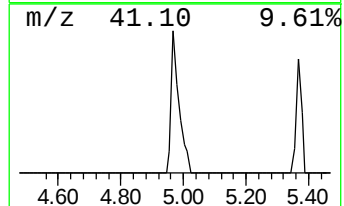
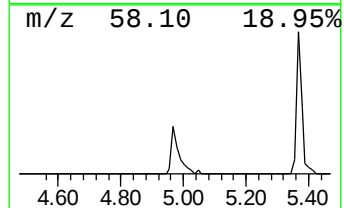
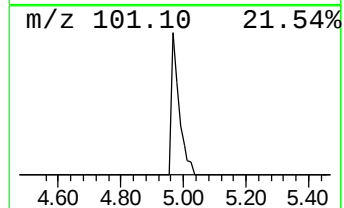
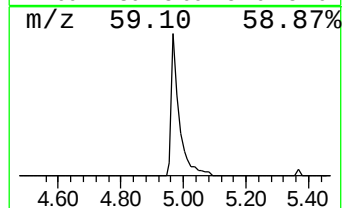
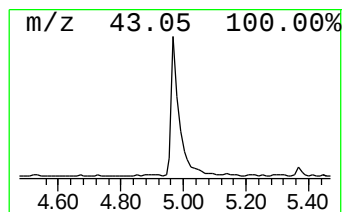
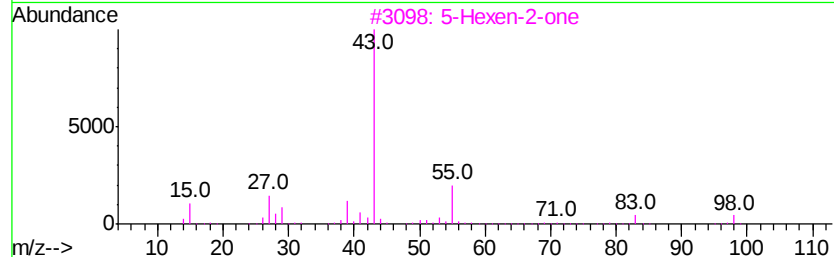
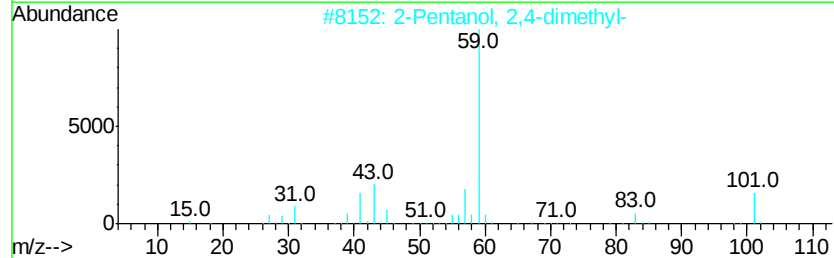
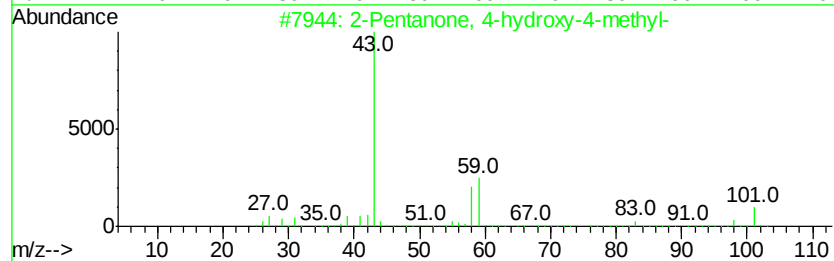
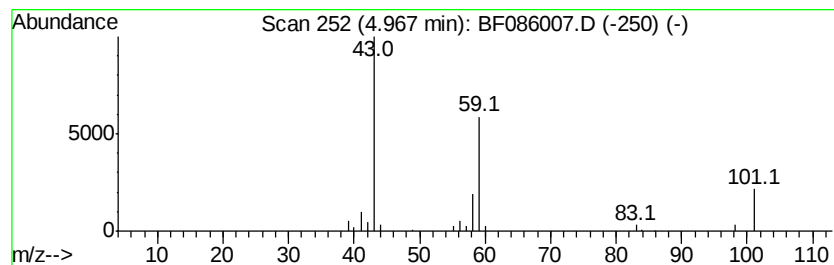
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.97	2.59 ng	83655	1,4-Dichlorobenzene-d4	6.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	23
3			5-Hexen-2-one	98	C6H10O	000109-49-9	9
4			4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	9
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9



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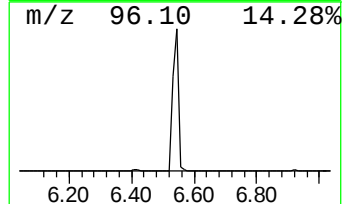
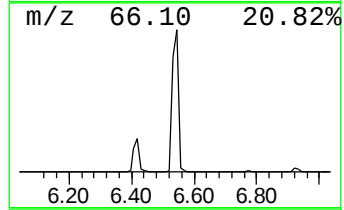
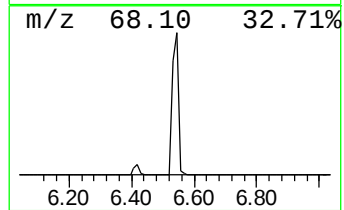
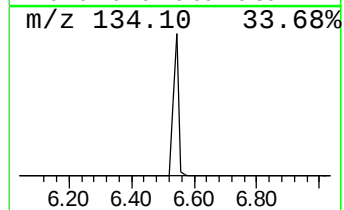
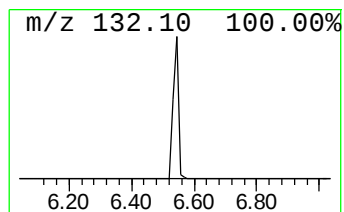
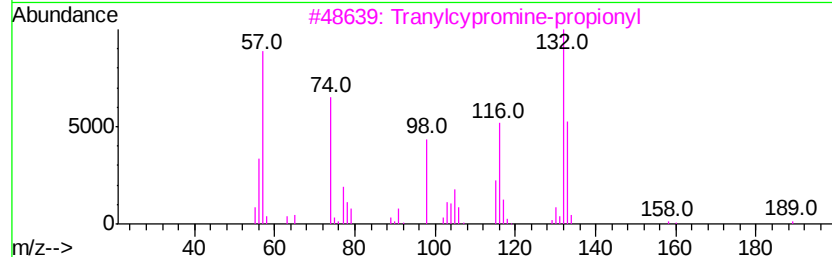
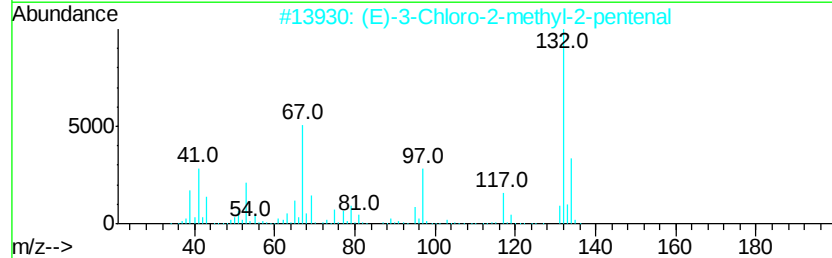
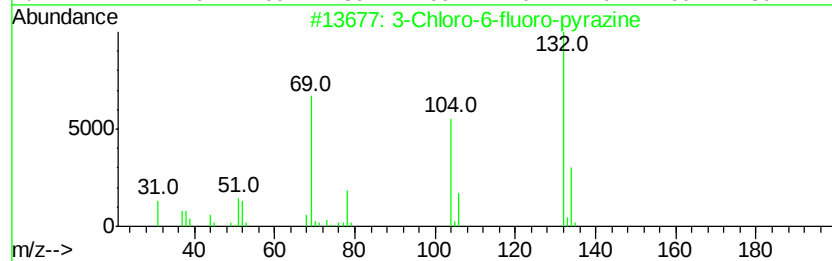
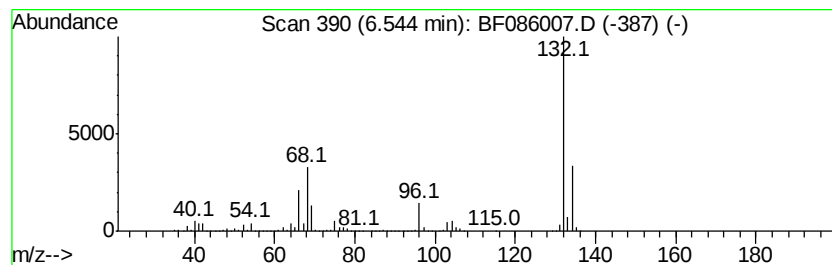
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Peak Number 2 unknown6.54 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.54	87.72 ng	2828070	1,4-Dichlorobenzene-d4	6.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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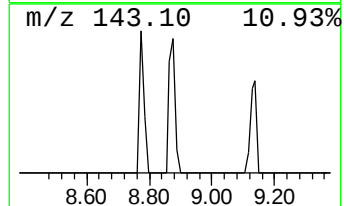
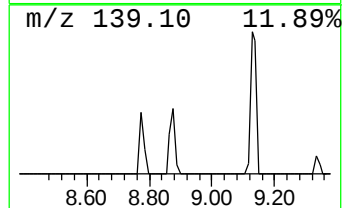
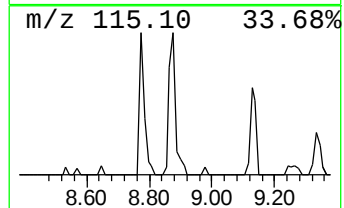
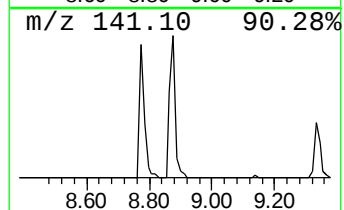
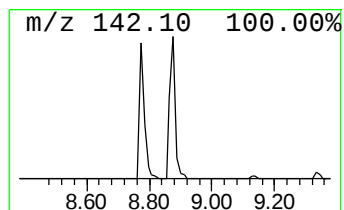
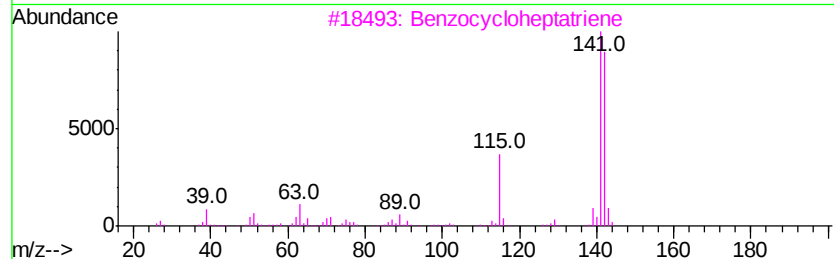
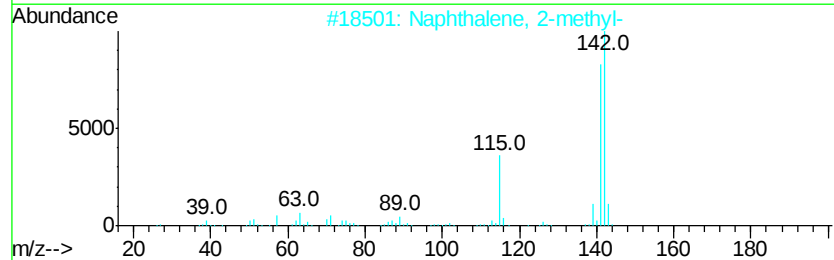
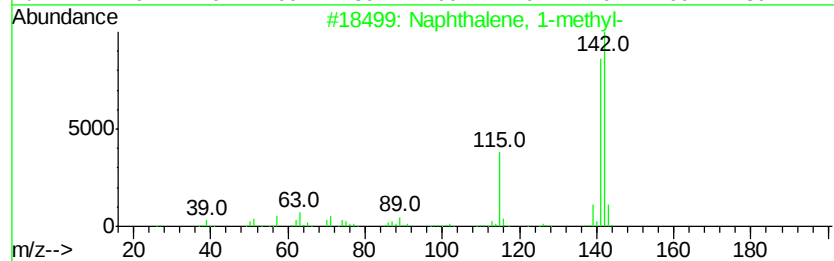
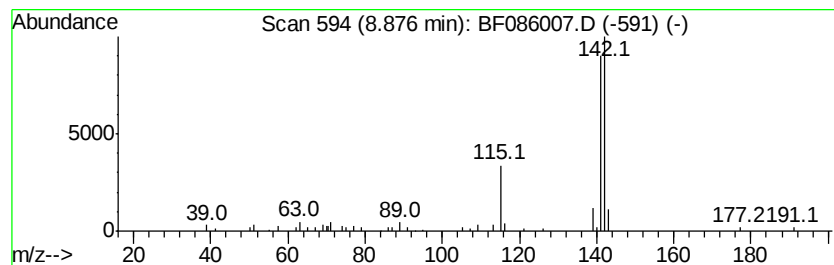
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 Naphthalene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.88	2.09 ng	97134	Naphthalene-d8	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
3		Benzocycloheptatriene	142	C11H10	000264-09-5	91
4		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	90
5		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	86



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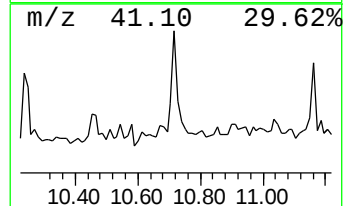
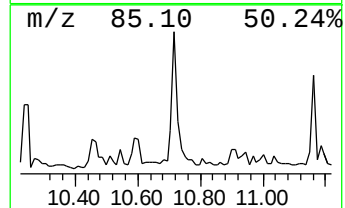
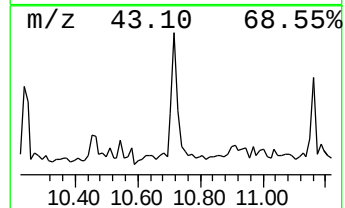
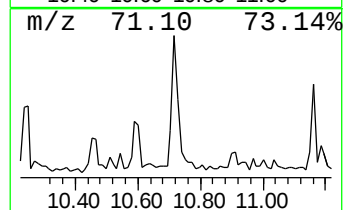
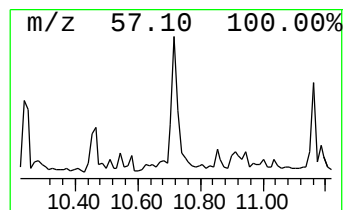
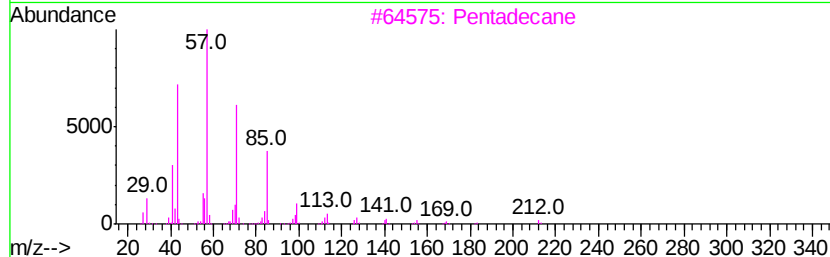
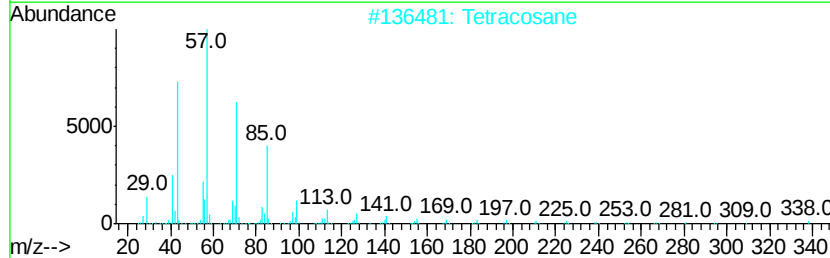
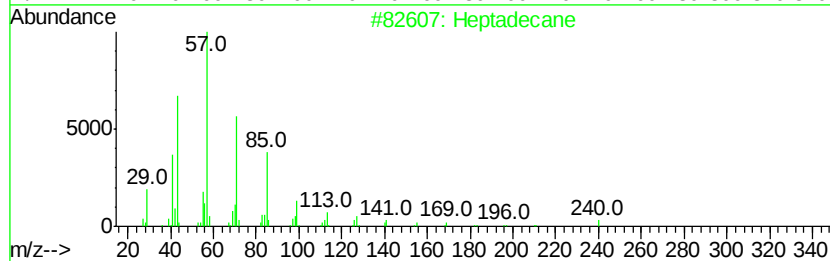
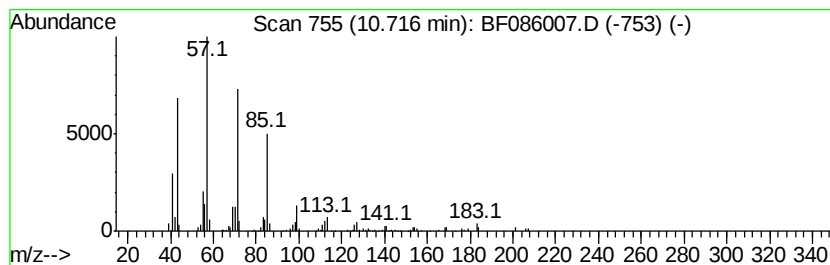
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 Heptadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.72	2.99 ng	169391	Phenanthrene-d10	11.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	91
2	Tetracosane	338	C24H50	000646-31-1	90
3	Pentadecane	212	C15H32	000629-62-9	90
4	Hexadecane	226	C16H34	000544-76-3	90
5	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	87



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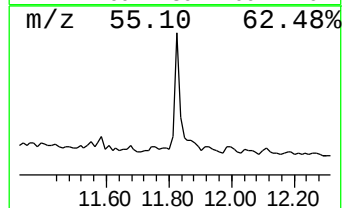
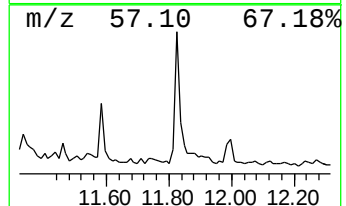
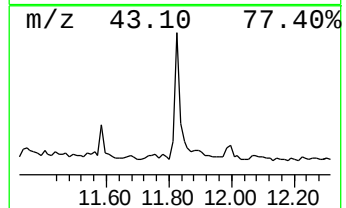
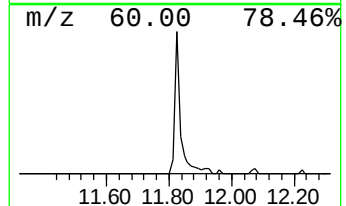
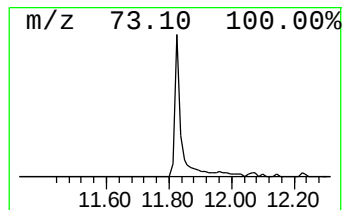
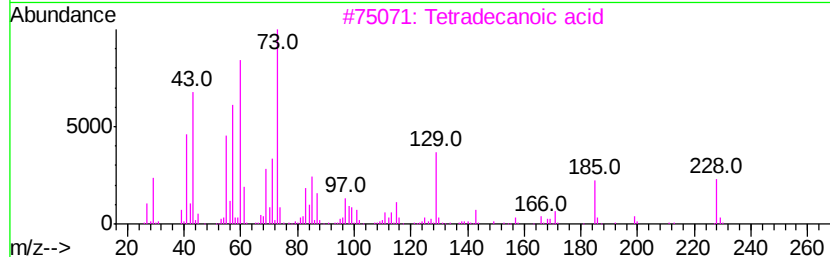
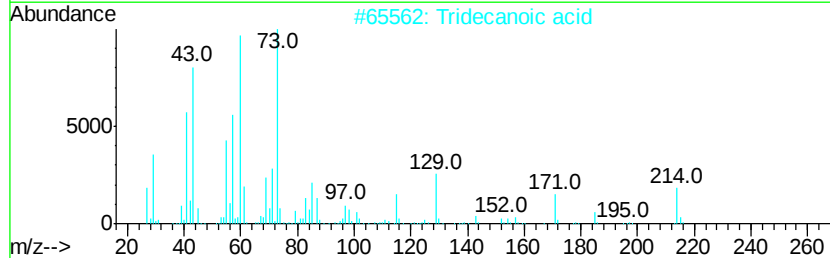
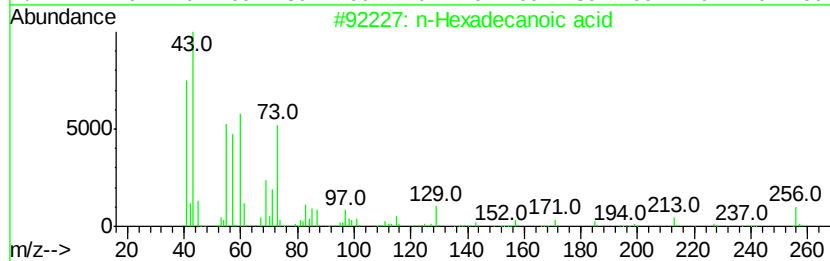
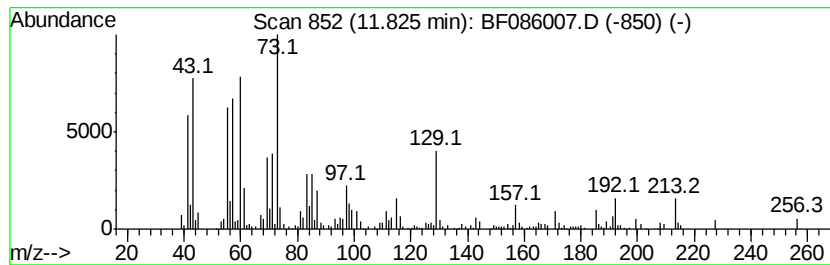
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.82	3.86 ng	218262	Phenanthrene-d10	11.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2	Tridecanoic acid	214	C13H26O2	000638-53-9	93
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	86
4	n-Decanoic acid	172	C10H20O2	000334-48-5	76
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040216\
Data File : BF086007.D
Acq On : 2 Apr 2016 13:05
Operator : UM/SJ
Sample : H2055-09
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BH-9(5-7)

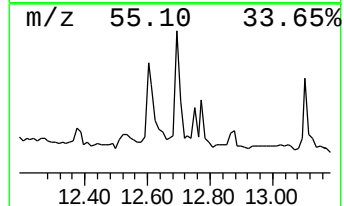
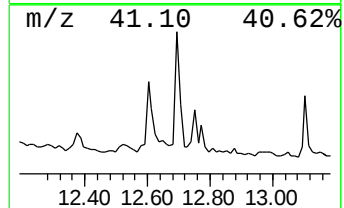
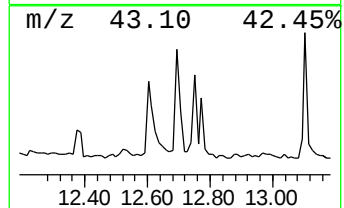
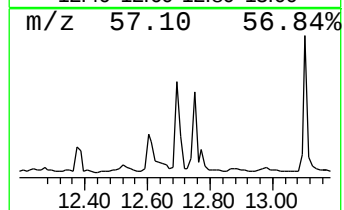
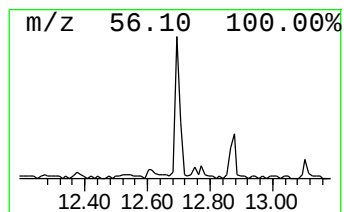
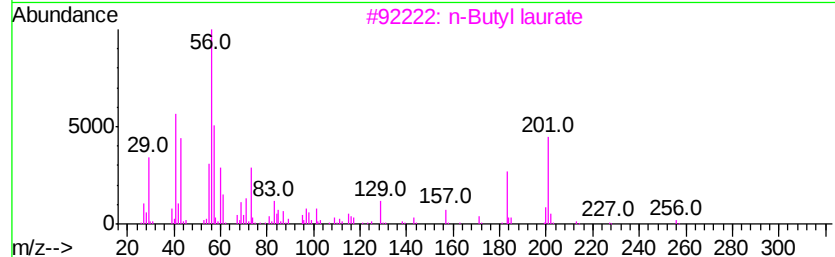
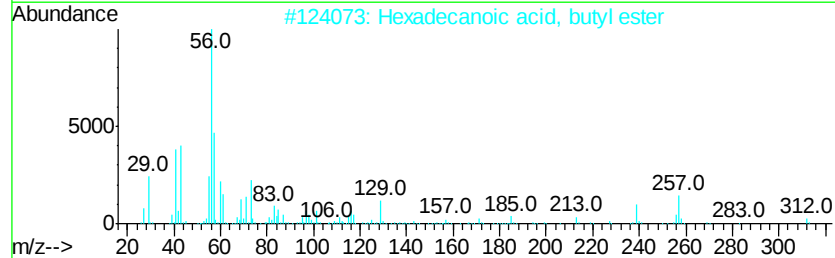
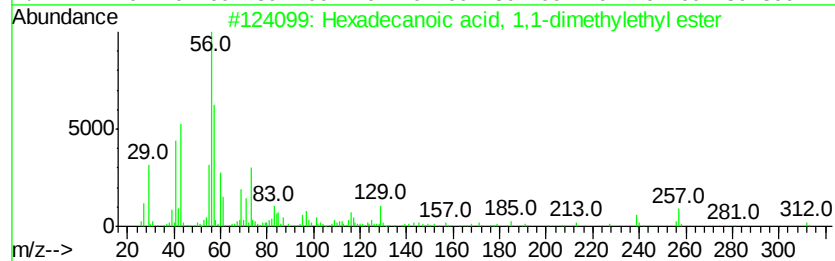
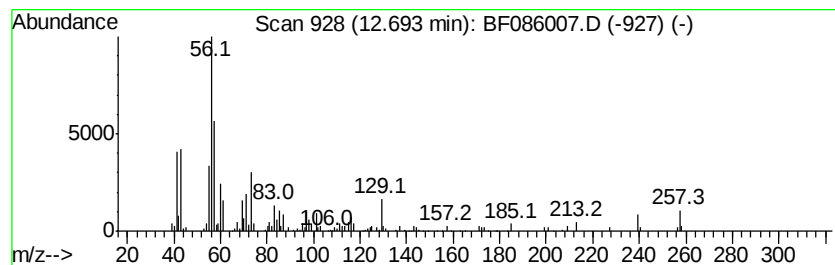
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecanoic acid, 1,1-dime... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.69	2.44 ng	132600	Chrysene-d12	13.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	87
2		Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	83
3		n-Butyl laurate	256	C16H32O2	000106-18-3	81
4		N-Nitroso-2-isopropyl-4,4-dimeth...	172	C8H16N2O2	039884-58-7	50
5		2,2-Dimethyl-1,3-butanediol	118	C6H14O2	000076-35-7	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040216\
Data File : BF086007.D
Acq On : 2 Apr 2016 13:05
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ClientSampleId :
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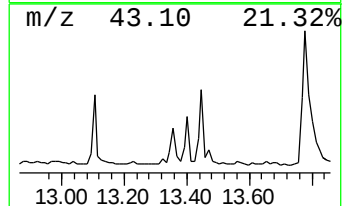
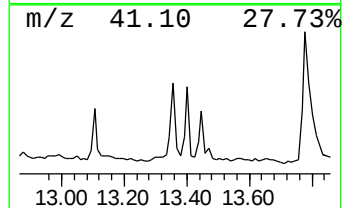
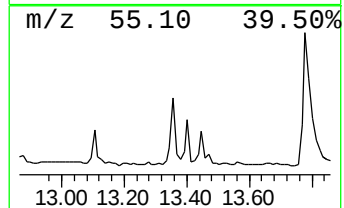
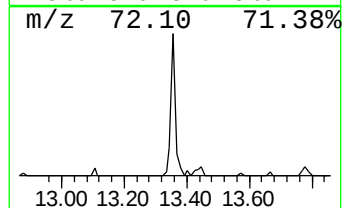
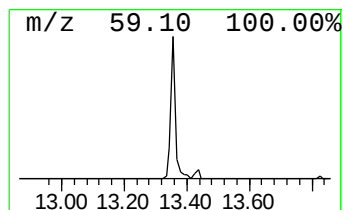
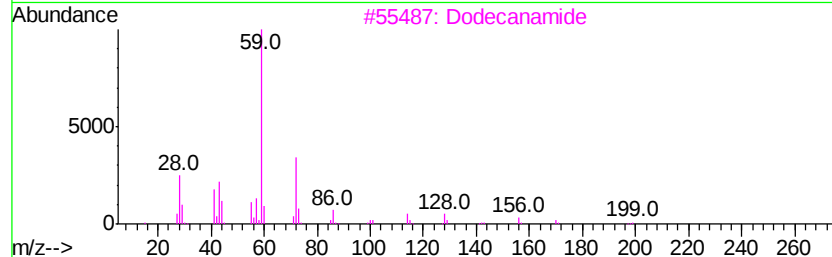
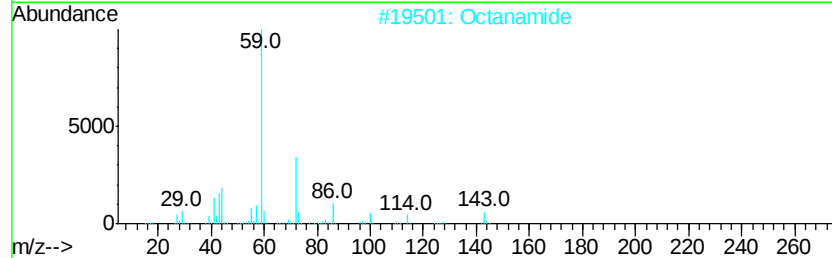
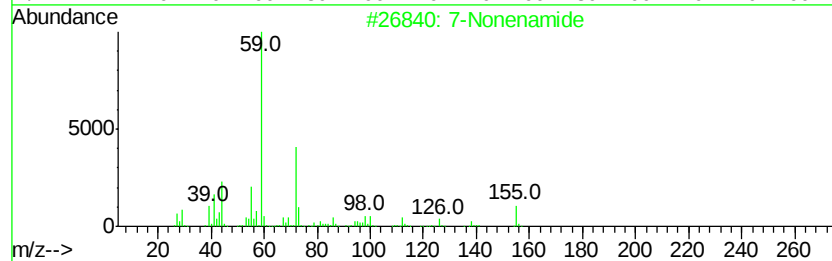
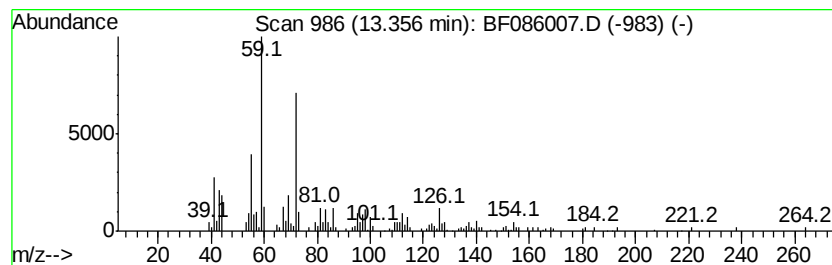
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 7-Nonenamide Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.36	2.14 ng	116157	Chrysene-d12	13.93

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Nonenamide		155	C9H17NO	090949-53-4	80
2	Octanamide		143	C8H17NO	000629-01-6	53
3	Dodecanamide		199	C12H25NO	001120-16-7	50
4	Tetradecanamide		227	C14H29NO	000638-58-4	50
5	Benzeneethanamine, 2-fluoro-.bet...		229	C11H16FN03	061338-98-5	42



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040216\
Data File : BF086007.D
Acq On : 2 Apr 2016 13:05
Operator : UM/SJ
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Misc :
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ClientSampleId :
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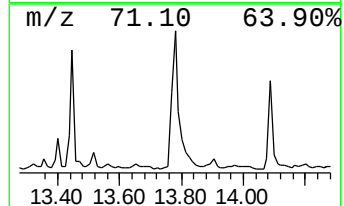
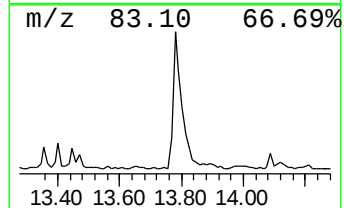
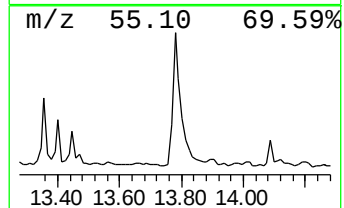
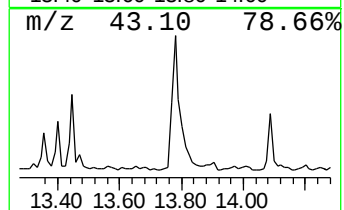
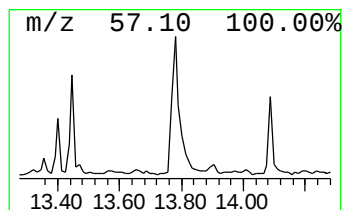
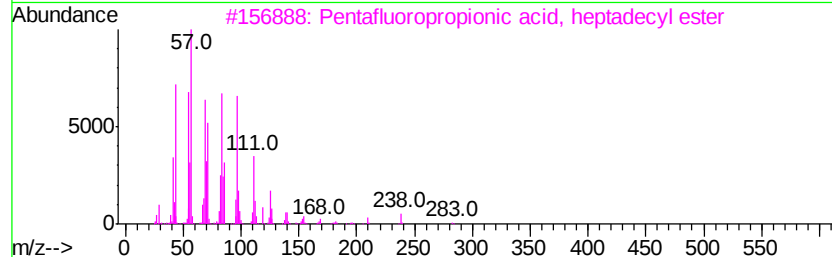
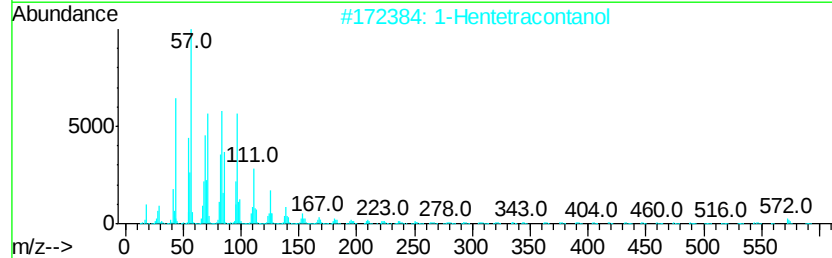
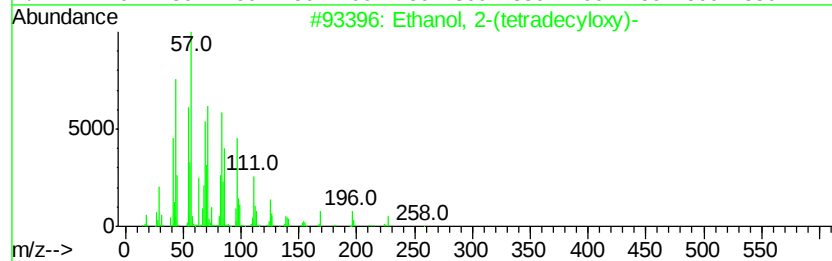
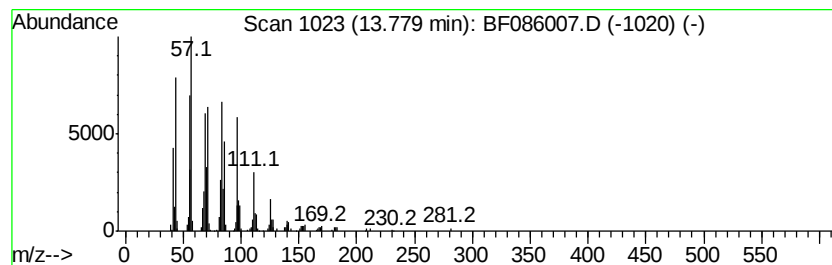
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Ethanol, 2-(tetradecyloxy)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.78	7.44 ng	404553	Chrysene-d12	13.93

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
2			1-Hentetracontanol	593	C41H84O	040710-42-7	91
3			Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	87
4			17-Pentatriacontene	491	C35H70	006971-40-0	83
5			Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040216\
Data File : BF086007.D
Acq On : 2 Apr 2016 13:05
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Misc :
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Instrument :
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ClientSampleId :
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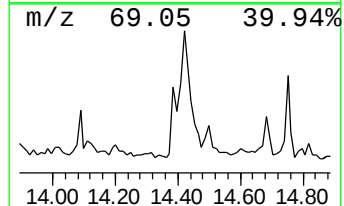
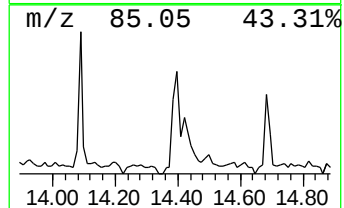
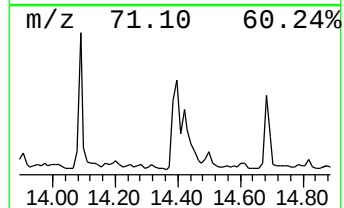
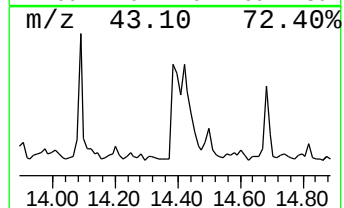
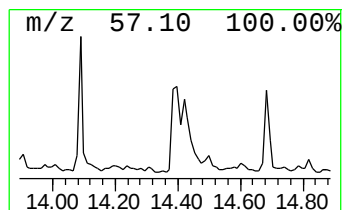
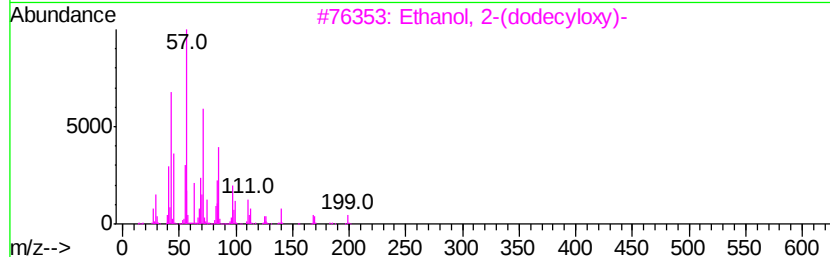
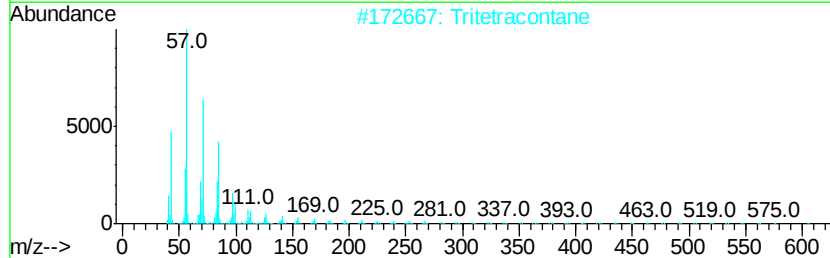
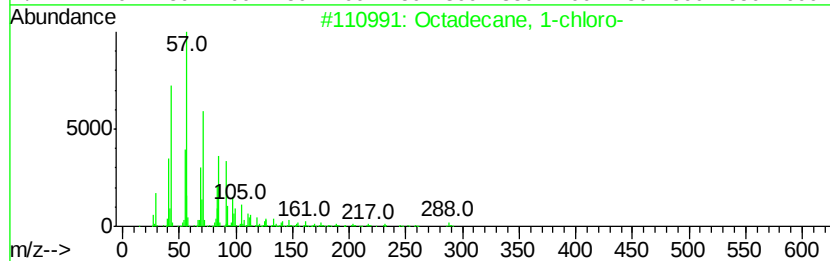
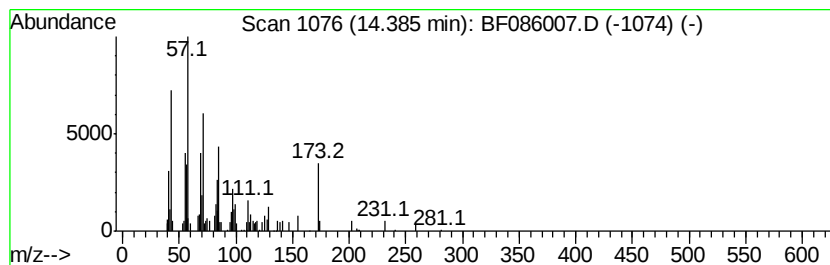
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecane, 1-chloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.39	2.14 ng	116420	Chrysene-d12	13.93

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	64
2			Tritetracontane	605	C43H88	007098-21-7	58
3			Ethanol, 2-(dodecyloxy)-	230	C14H30O2	004536-30-5	58
4			Octadecane, 1-bromo-	332	C18H37Br	000112-89-0	58
5			Nonadecane	268	C19H40	000629-92-5	49



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040216\
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Instrument :
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ClientSampleId :
BH-9(5-7)

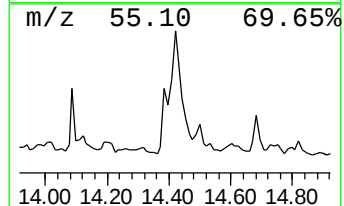
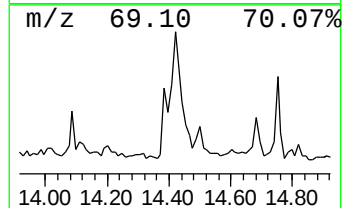
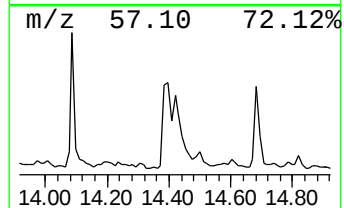
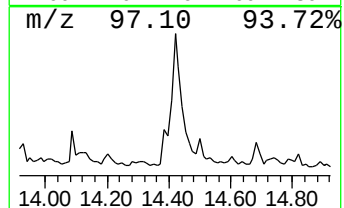
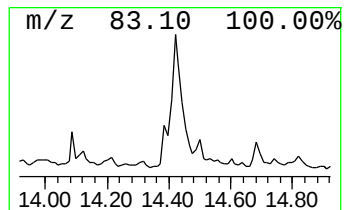
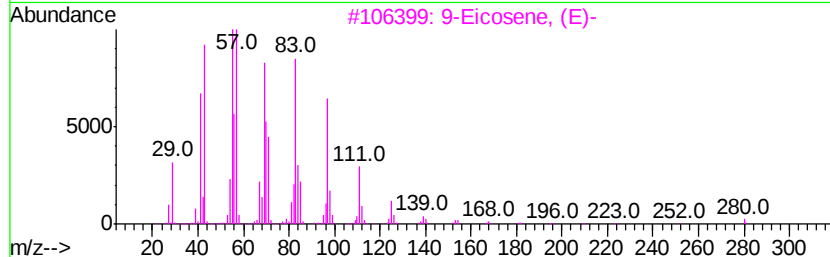
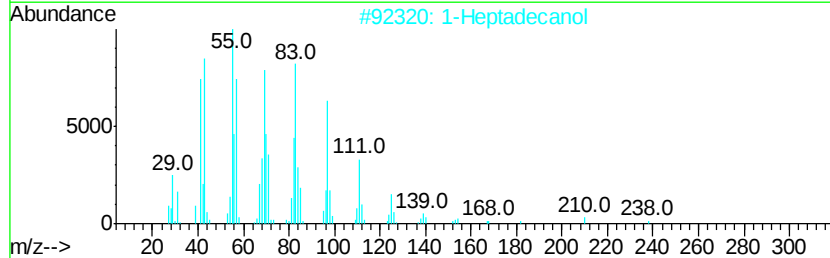
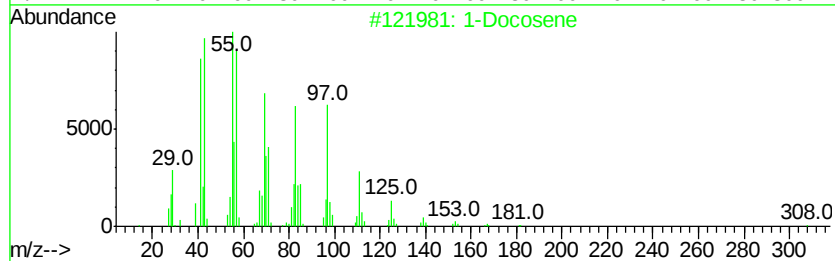
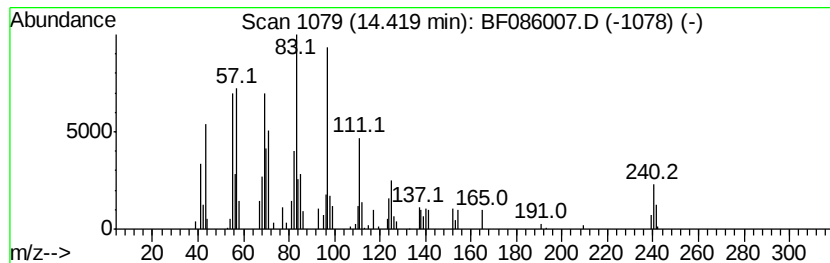
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Docosene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.42	3.01 ng	163896	Chrysene-d12	13.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Docosene	308	C22H44	001599-67-3	90
2		1-Heptadecanol	256	C17H36O	001454-85-9	83
3		9-Eicosene, (E)-	280	C20H40	074685-29-3	72
4		3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	68
5		3-Eicosene, (E)-	280	C20H40	074685-33-9	68



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040216\
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ALS Vial : 43 Sample Multiplier: 1

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ClientSampleId :
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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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unknown6.54	6.54	87.7	ng	2828070	1	6.77	644781	20.0
Naphthalene, 1-me...	8.88	2.1	ng	97134	2	8.05	931232	20.0
Heptadecane	10.72	3.0	ng	169391	4	11.29	1132090	20.0
n-Hexadecanoic acid	11.82	3.9	ng	218262	4	11.29	1132090	20.0
Hexadecanoic acid...	12.69	2.4	ng	132600	5	13.93	1087990	20.0
7-Nonenamide	13.36	2.1	ng	116157	5	13.93	1087990	20.0
Ethanol, 2-(tetra...	13.78	7.4	ng	404553	5	13.93	1087990	20.0
Octadecane, 1-chl...	14.39	2.1	ng	116420	5	13.93	1087990	20.0
1-Docosene	14.42	3.0	ng	163896	5	13.93	1087990	20.0