

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121516\
 Data File : BF091644.D
 Acq On : 15 Dec 2016 22:32
 Operator : UM/SJ
 Sample : H5940-05
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS07-1FT

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-BF120916.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	4.956	248	251	261	rBV	1782360	2455146	37.18%	3.932%
2	5.390	286	289	291	rBV	6163727	6431251	97.39%	10.301%
3	6.430	377	380	382	rVB	5720881	6343062	96.06%	10.160%
4	6.487	382	385	387	rVB	188581	231549	3.51%	0.371%
5	6.556	388	391	393	rBV	6803465	6603322	100.00%	10.577%
6	6.785	409	411	413	rBV	2021373	1684659	25.51%	2.698%
7	6.910	420	422	423	rBV	517111	619064	9.38%	0.992%
8	6.945	423	425	427	rVB	5131131	5034822	76.25%	8.064%
9	7.356	457	461	463	rBV	2942772	4250941	64.38%	6.809%
10	7.459	466	470	474	rVB	82272	128302	1.94%	0.206%
11	8.065	521	523	526	rBV	1586586	1929070	29.21%	3.090%
12	9.150	615	618	620	rBV	6775940	6017025	91.12%	9.638%
13	9.265	627	628	630	rVB	113381	98117	1.49%	0.157%
14	9.390	637	639	643	rBV	64222	79115	1.20%	0.127%
15	9.550	650	653	655	rBV	1332974	1787193	27.07%	2.863%
16	9.768	670	672	673	rVB	84727	77762	1.18%	0.125%
17	9.825	673	677	679	rBV	2528608	2253862	34.13%	3.610%
18	10.259	714	715	717	rVB	152662	160333	2.43%	0.257%
19	10.488	731	735	738	rBV2	76985	192518	2.92%	0.308%
20	10.545	738	740	741	rVV	461031	402214	6.09%	0.644%
21	10.613	743	746	748	rBV	3311551	3581510	54.24%	5.737%
22	10.739	755	757	762	rBV	229641	330298	5.00%	0.529%
23	11.185	794	796	797	rBV	174955	158354	2.40%	0.254%
24	11.299	804	806	809	rBV	1597120	1635669	24.77%	2.620%
25	11.539	824	827	832	rBV	65591	104484	1.58%	0.167%
26	11.836	851	853	858	rBV2	132535	231925	3.51%	0.371%
27	12.419	902	904	907	rVB	364771	396296	6.00%	0.635%
28	12.511	910	912	914	rBV	153370	141505	2.14%	0.227%
29	12.739	928	932	933	rBV	101212	121740	1.84%	0.195%
30	12.888	942	945	948	rBV	3938871	3758273	56.91%	6.020%
31	13.471	993	996	999	rBV2	78650	135392	2.05%	0.217%
32	13.791	1021	1024	1027	rBV	374113	438811	6.65%	0.703%
33	13.939	1034	1037	1039	rBV	1108133	1587123	24.04%	2.542%
34	14.420	1078	1079	1082	rBV	77004	112082	1.70%	0.180%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.785	1109	1111	1113	rBV	99427	104433	1.58%	0.167%
36	14.854	1115	1117	1121	rVB2	71143	91013	1.38%	0.146%
37	14.945	1123	1125	1129	rVB	69170	153613	2.33%	0.246%
38	15.037	1130	1133	1136	rBV2	128369	204502	3.10%	0.328%
39	15.220	1147	1149	1152	rVB	44414	66645	1.01%	0.107%
40	15.277	1152	1154	1157	rVV	43773	72661	1.10%	0.116%
41	15.345	1157	1160	1162	rVV	909997	1226153	18.57%	1.964%
42	15.391	1162	1164	1168	rVB	79061	133770	2.03%	0.214%
43	15.791	1196	1199	1202	rBV	122800	206798	3.13%	0.331%
44	16.260	1237	1240	1244	rBV2	68648	139227	2.11%	0.223%
45	16.797	1285	1287	1299	rVB2	63759	199962	3.03%	0.320%
46	18.603	1440	1445	1450	rBV2	106054	320945	4.86%	0.514%

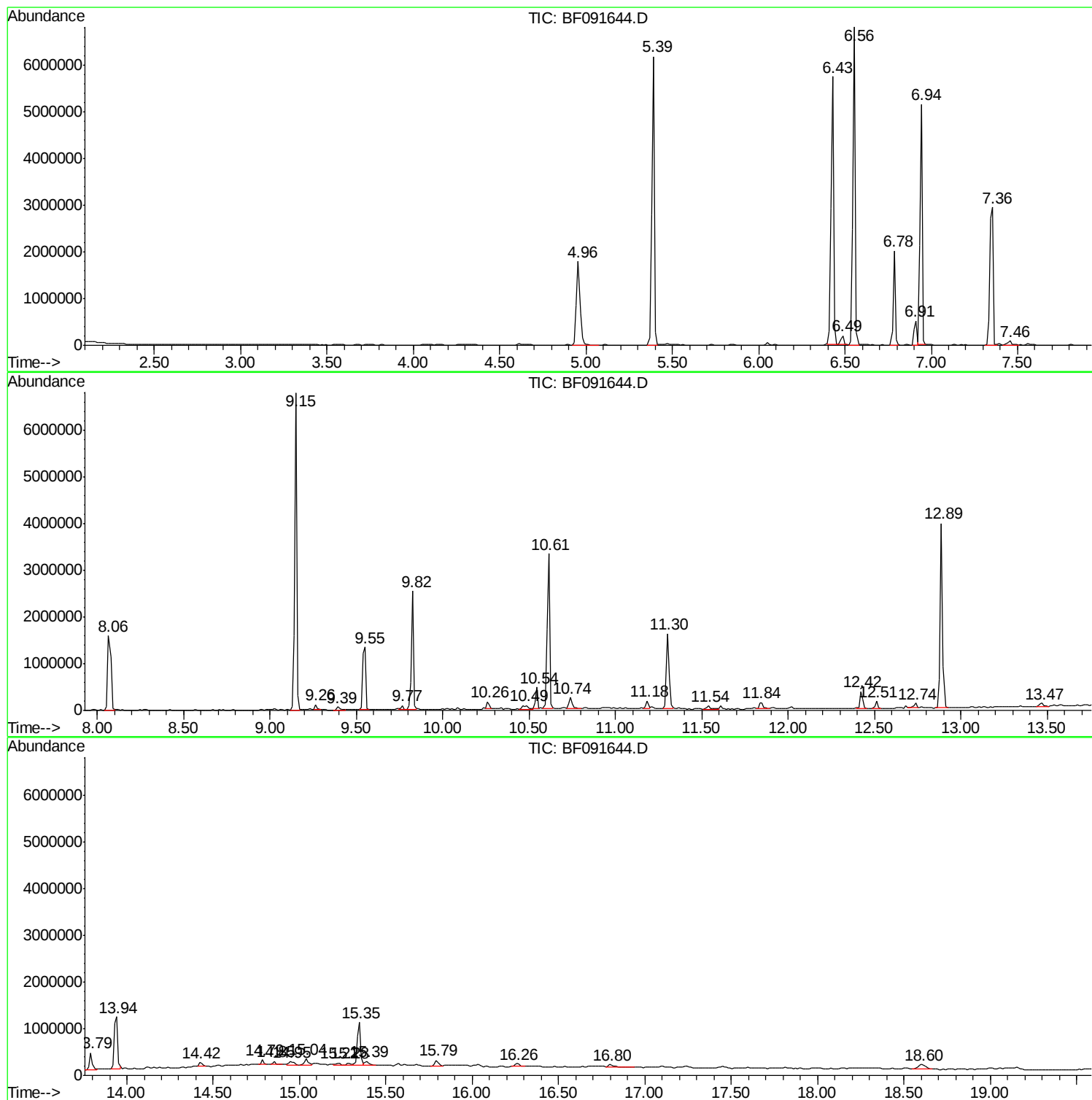
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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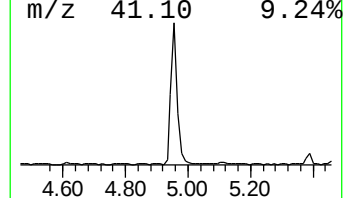
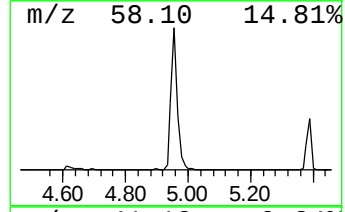
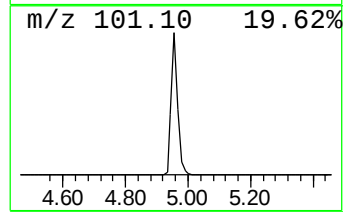
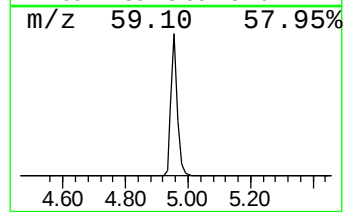
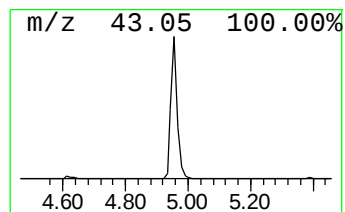
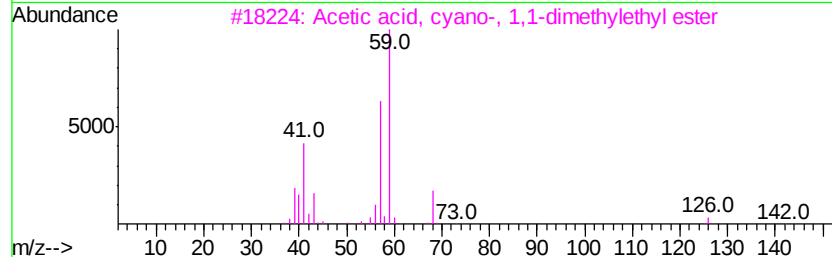
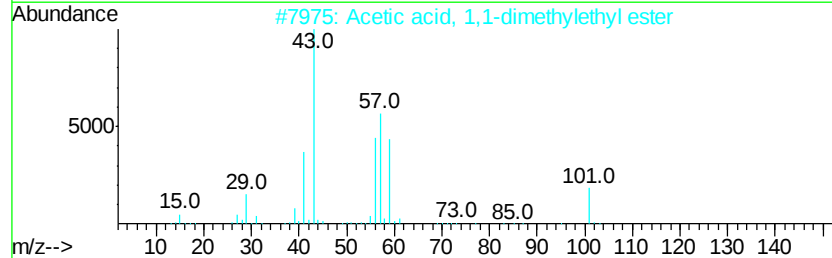
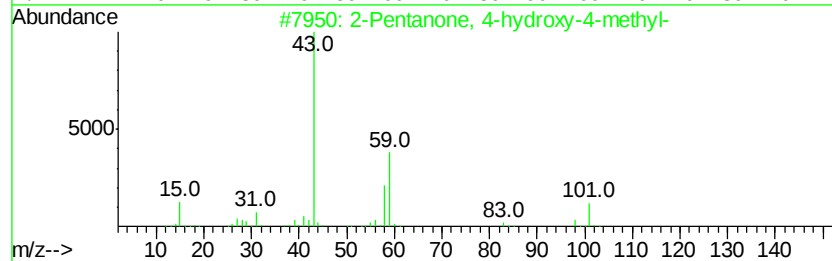
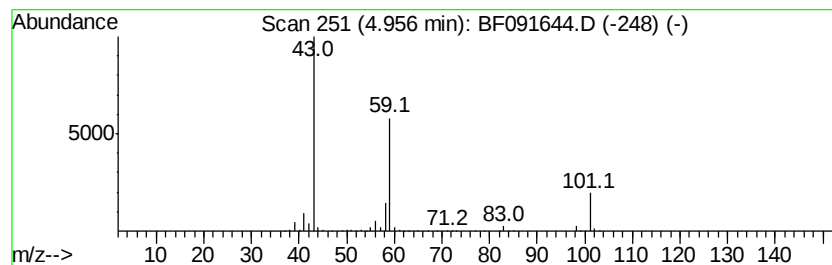
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
4.96	29.15 ng	2455150	1,4-Dichlorobenzene-d4	6.78

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56	
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39	
3	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25	
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9	
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9	



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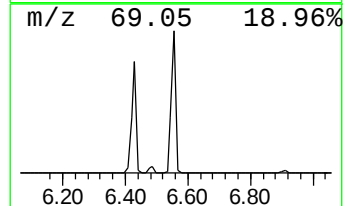
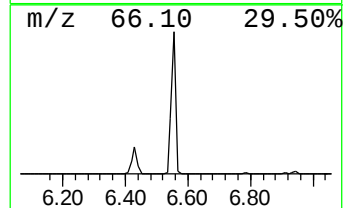
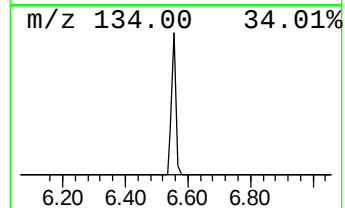
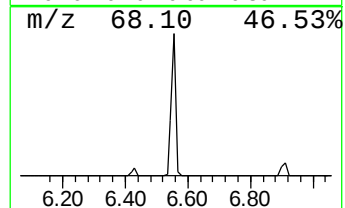
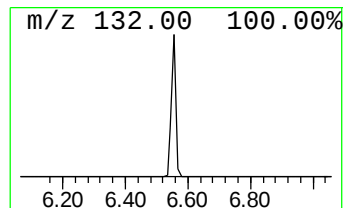
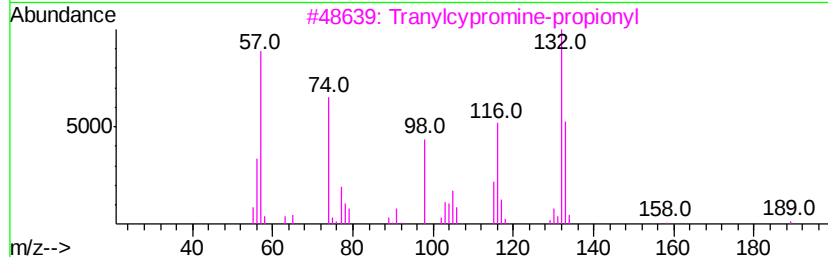
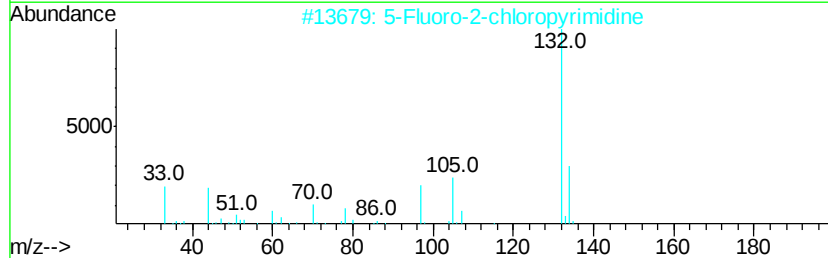
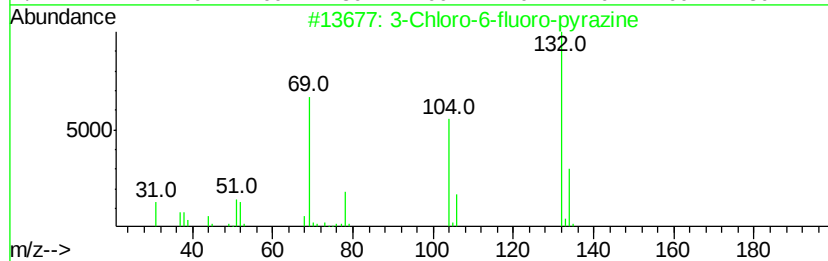
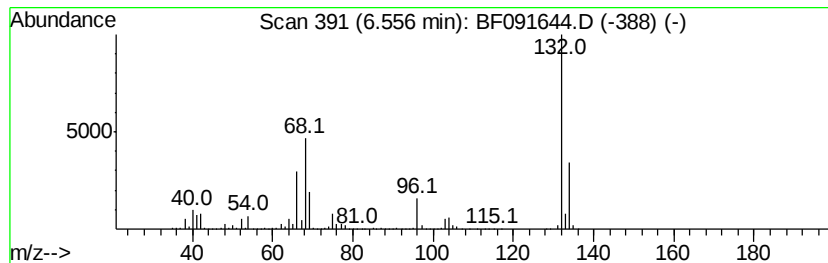
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	78.39 ng	6603320	1,4-Dichlorobenzene-d4	6.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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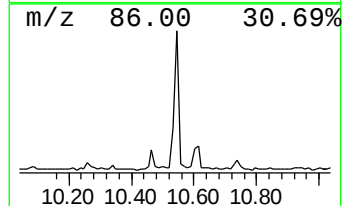
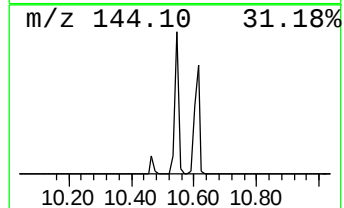
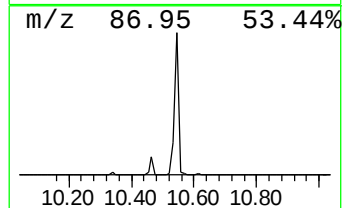
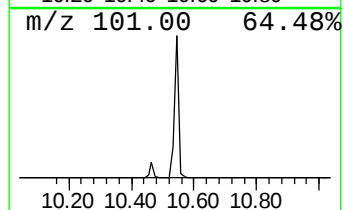
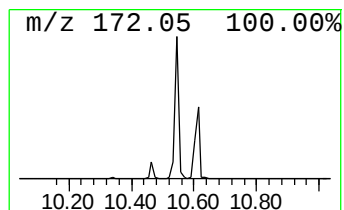
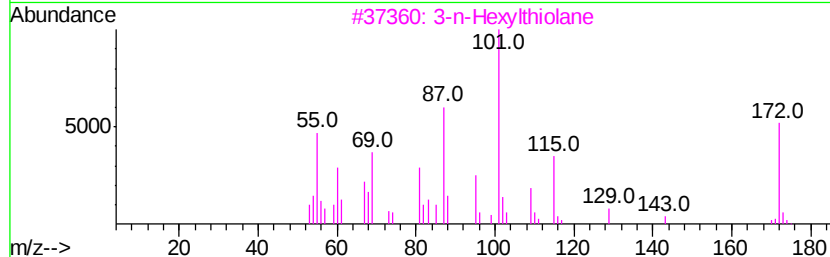
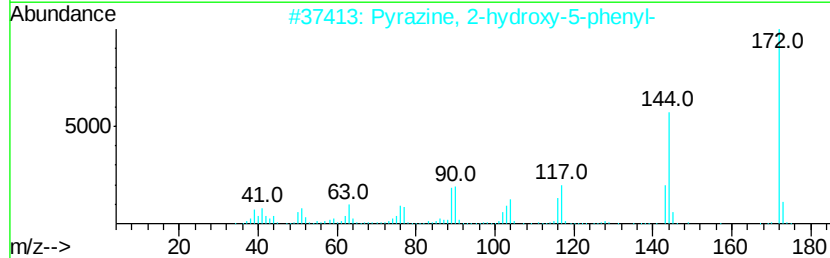
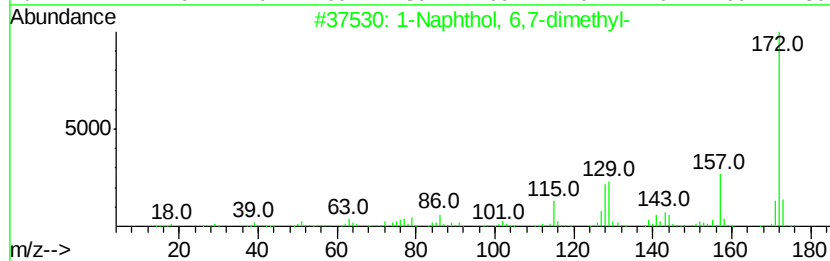
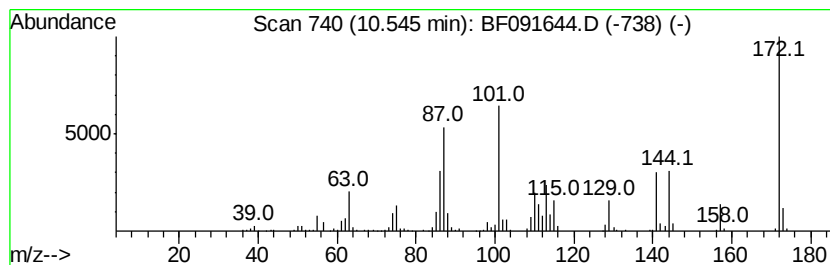
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown10.54 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.54	3.57 ng	402214	Acenaphthene-d10	9.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthol,	6,7-dimethyl-	172	C12H12O	031776-14-4	41	
2	Pyrazine,	2-hydroxy-5-phenyl-	172	C10H8N2O	025844-72-8	27	
3	3-n-Hexylthiolane		172	C10H20S	001613-50-9	23	
4	2,5-Dimethyl-3-phenylfuran		172	C12H12O	019842-57-0	22	
5	Tetroquinone		172	C6H4O6	000319-89-1	16	



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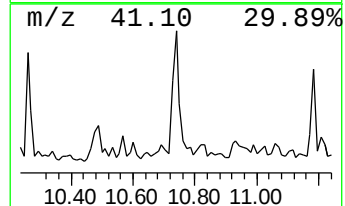
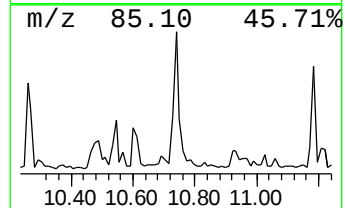
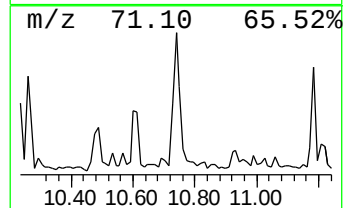
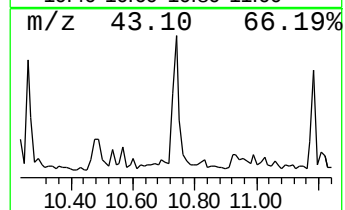
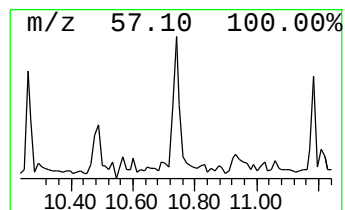
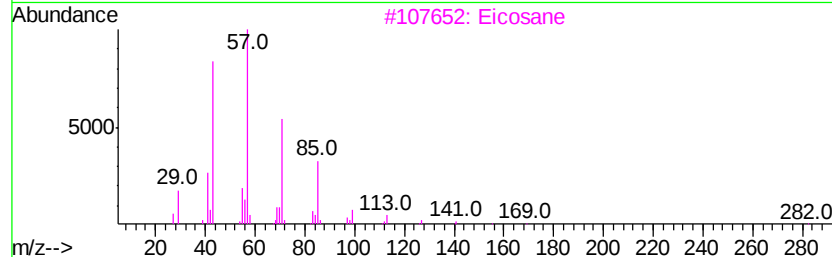
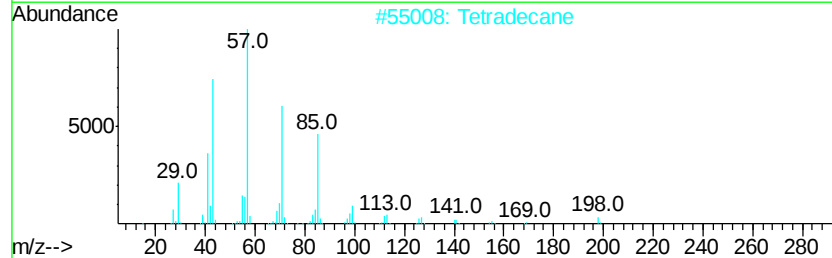
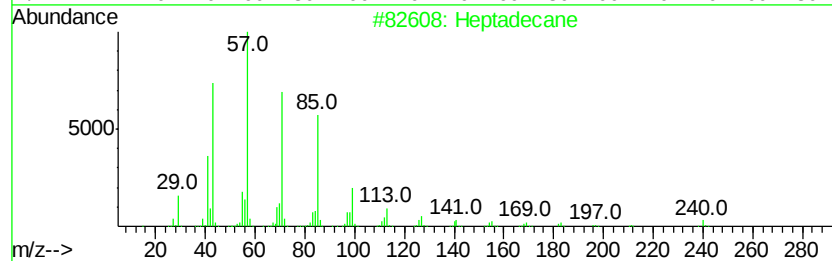
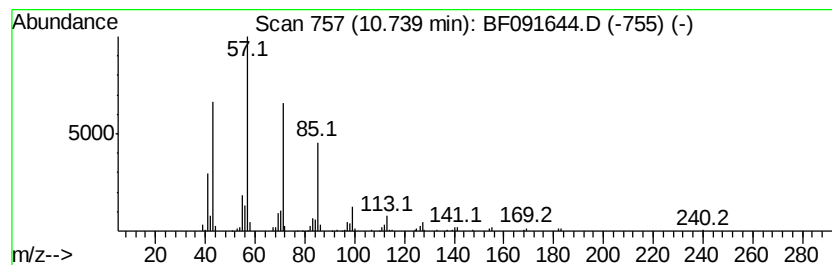
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Heptadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.74	4.04 ng	330298	Phenanthrene-d10	11.30

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	97
2	Tetradecane		198	C14H30	000629-59-4	91
3	Eicosane		282	C20H42	000112-95-8	91
4	Hexadecane		226	C16H34	000544-76-3	90
5	Heneicosane		296	C21H44	000629-94-7	90



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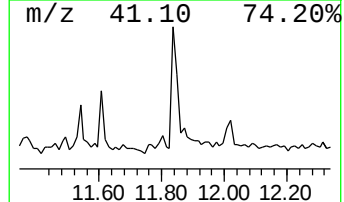
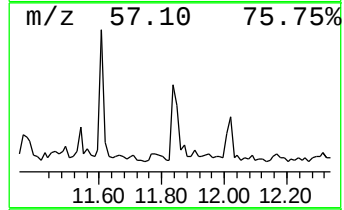
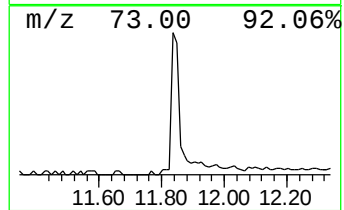
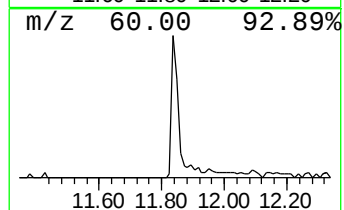
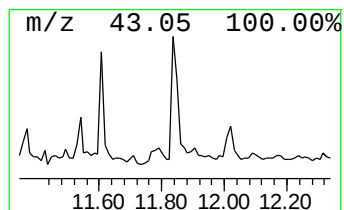
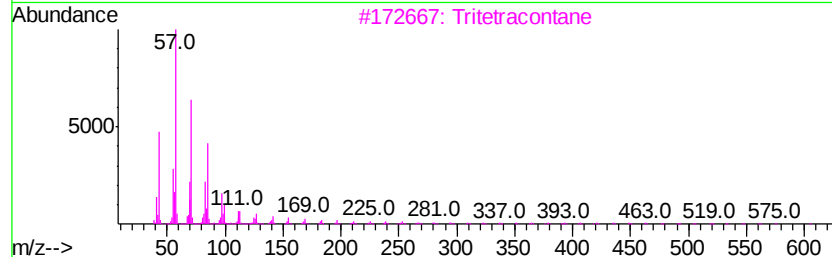
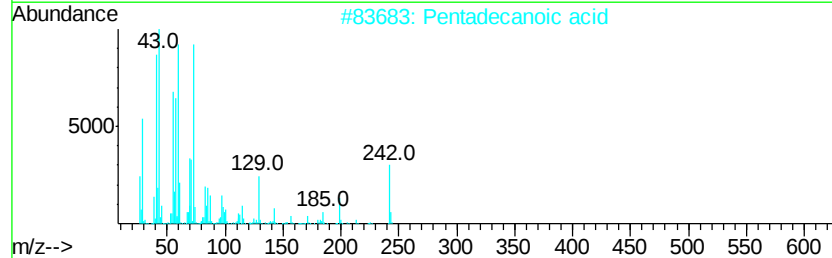
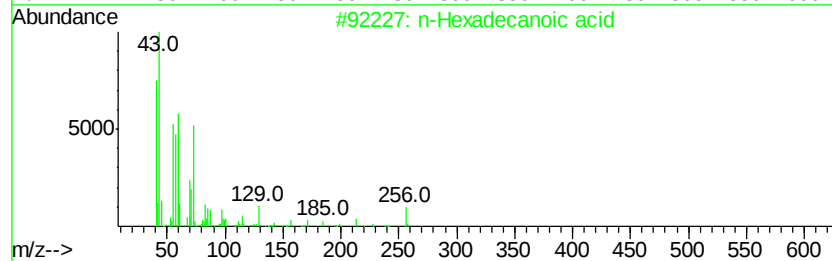
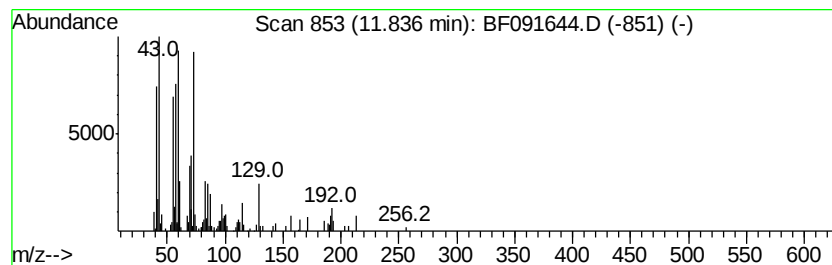
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.84	2.84 ng	231925	Phenanthrene-d10	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	86
3		Tritetracontane	605	C43H88	007098-21-7	20
4		Tetratetracontane	619	C44H90	007098-22-8	18
5		Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121516\
Data File : BF091644.D
Acq On : 15 Dec 2016 22:32
Operator : UM/SJ
Sample : H5940-05
Misc :
ALS Vial : 18 Sample Multiplier: 1

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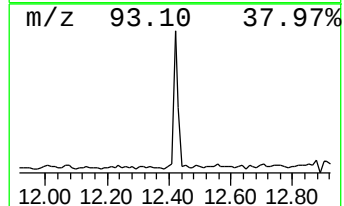
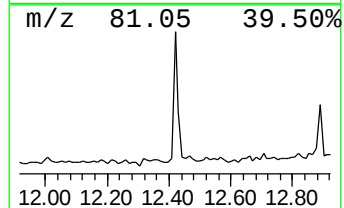
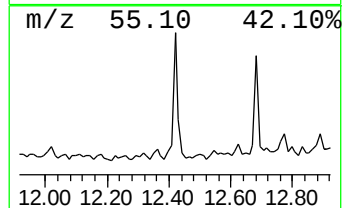
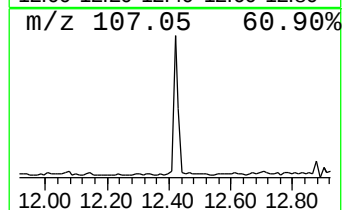
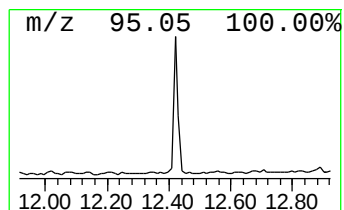
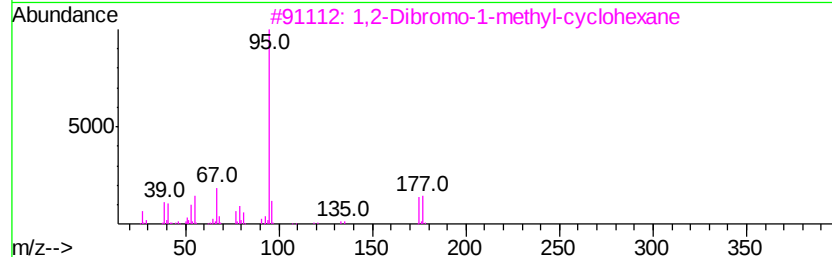
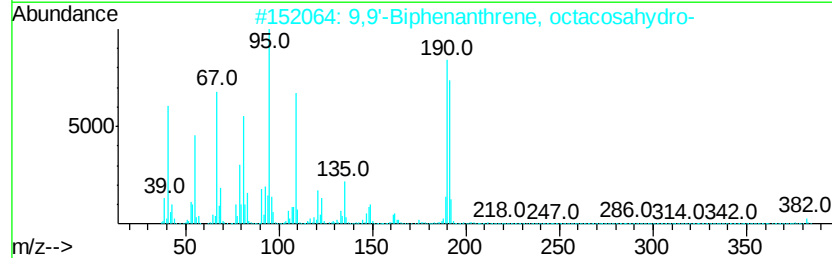
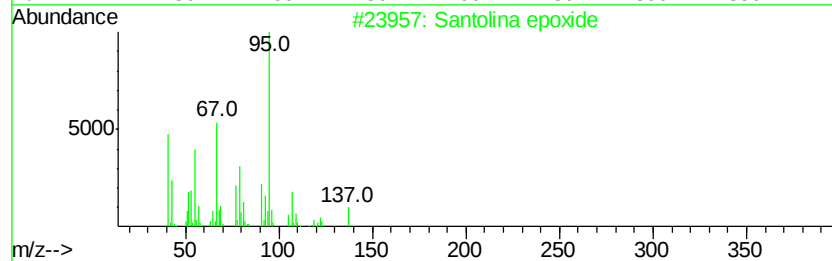
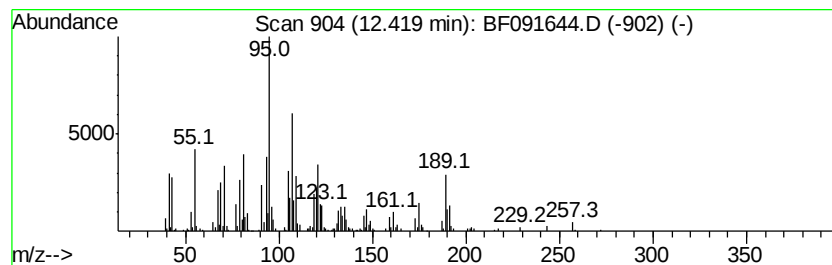
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown12.42 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.42	4.85 ng	396296	Phenanthrene-d10	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Santolina epoxide	152	C10H16O	060485-45-2	35
2		9,9'-Biphenanthrene, octacosahydro-	382	C28H46	055334-18-4	22
3		1,2-Dibromo-1-methyl-cyclohexane	254	C7H12Br2	1000192-26-1	22
4		4-(Bromomethyl)cyclohexene	174	C7H11Br	034960-41-3	22
5		trans,cis-1,7-Dimethylspiro[4.5]...	166	C12H22	1000111-72-7	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121516\
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Sample : H5940-05
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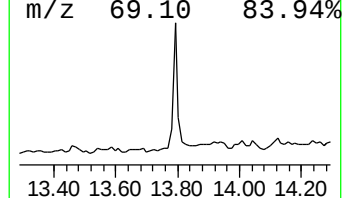
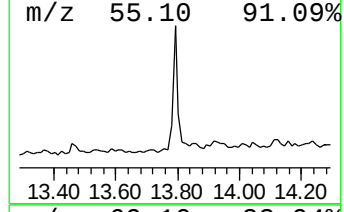
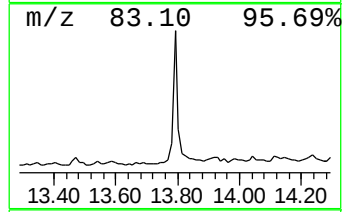
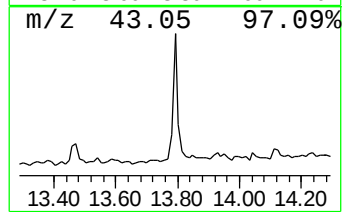
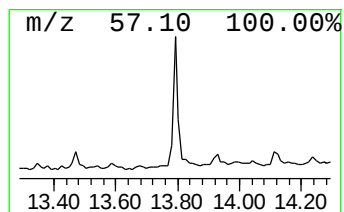
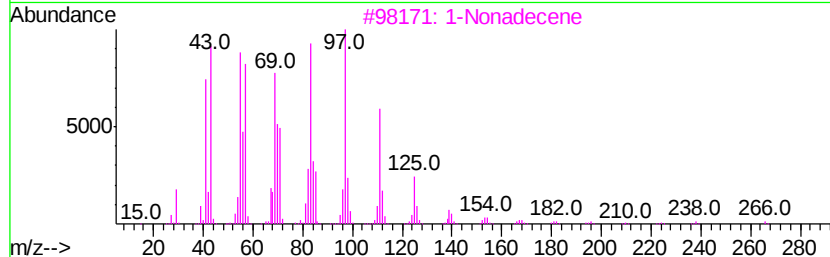
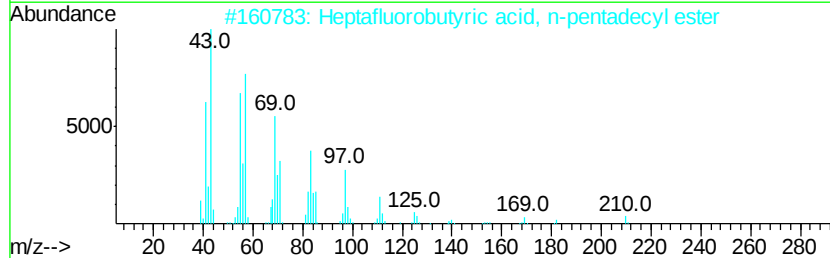
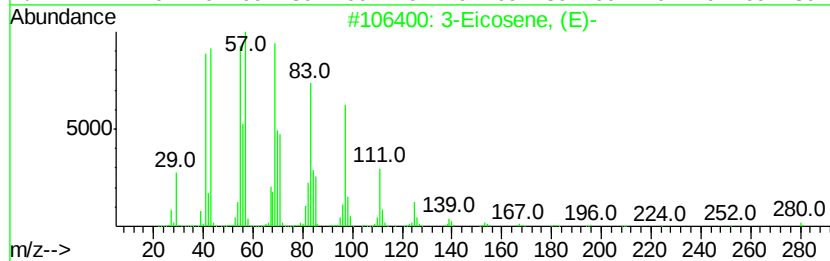
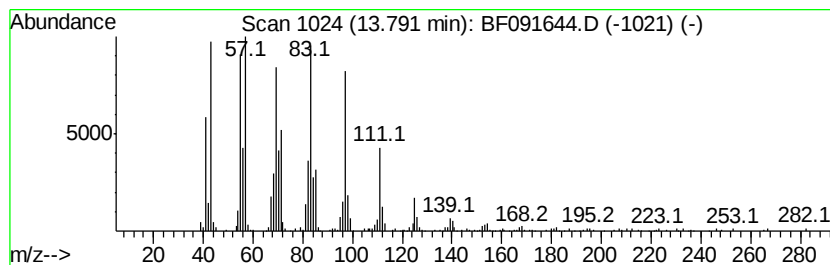
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 3-Eicosene, (E)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.79	5.53 ng	438811	Chrysene-d12	13.94	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-	280	C20H40	074685-33-9	91
2	Heptafluorobutyric acid, n-penta...	424	C19H31F7O2	1000216-79-3	91
3	1-Nonadecene	266	C19H38	018435-45-5	91
4	1-Hexadecanol	242	C16H34O	036653-82-4	90
5	1-Octadecanol	270	C18H38O	000112-92-5	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121516\
Data File : BF091644.D
Acq On : 15 Dec 2016 22:32
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Misc :
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Instrument :
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ClientSampleId :
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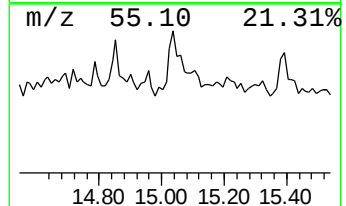
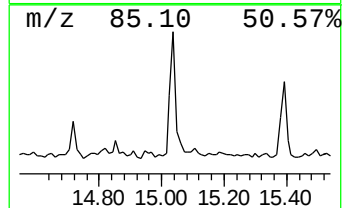
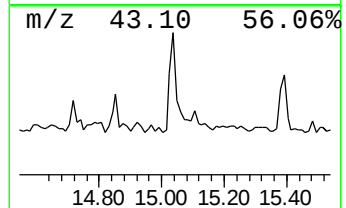
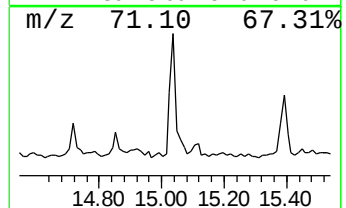
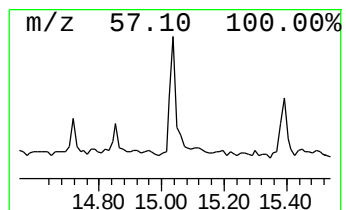
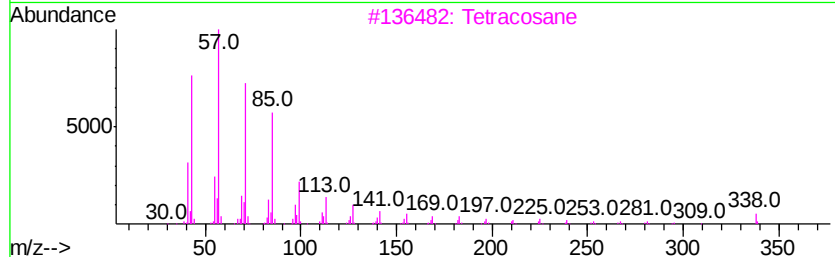
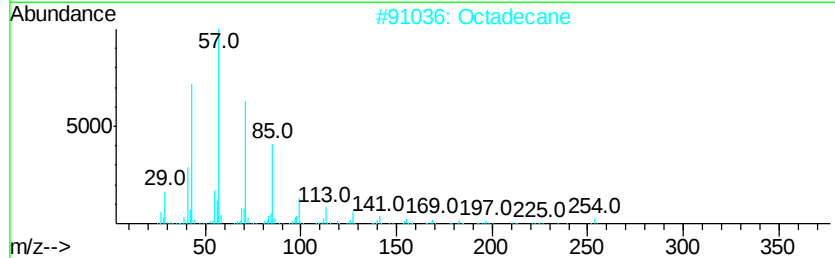
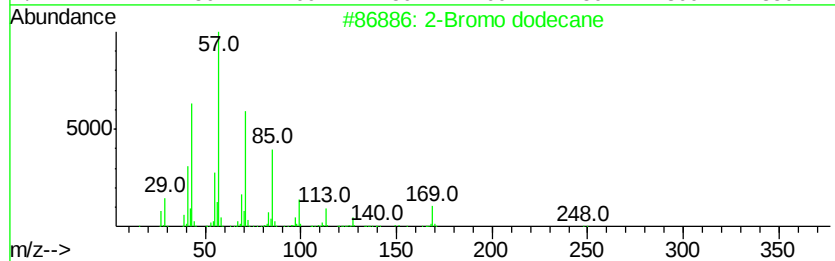
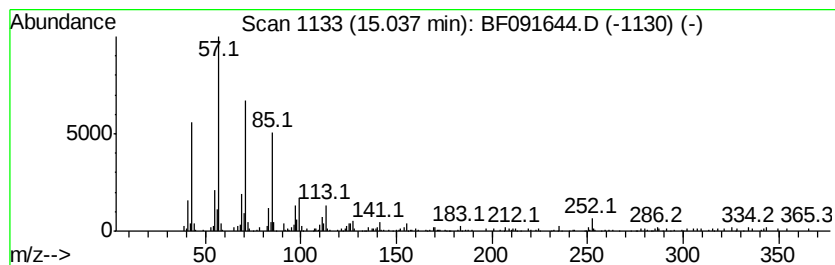
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2-Bromo dodecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.04	3.34 ng	204502	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Bromo dodecane	248	C12H25Br	013187-99-0	91
2		Octadecane	254	C18H38	000593-45-3	91
3		Tetracosane	338	C24H50	000646-31-1	87
4		Tetradecane, 2-methyl-	212	C15H32	001560-95-8	87
5		Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121516\
Data File : BF091644.D
Acq On : 15 Dec 2016 22:32
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Instrument :
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ClientSampleId :
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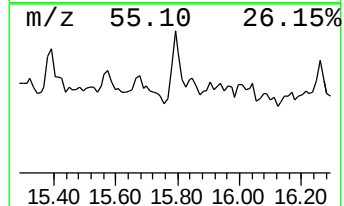
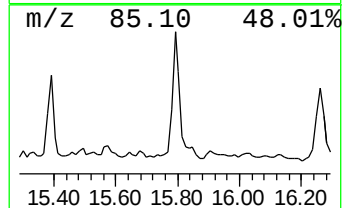
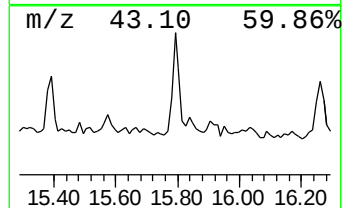
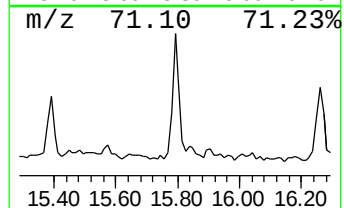
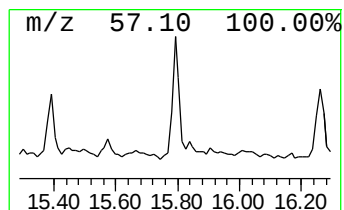
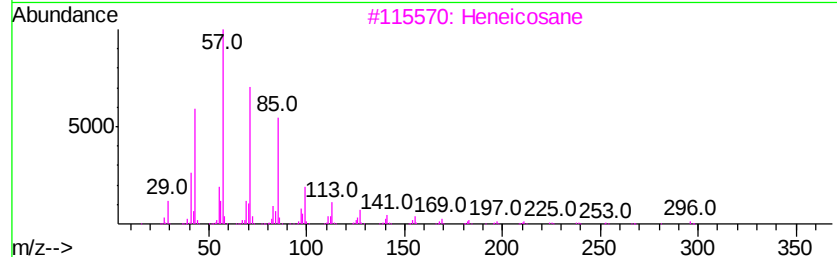
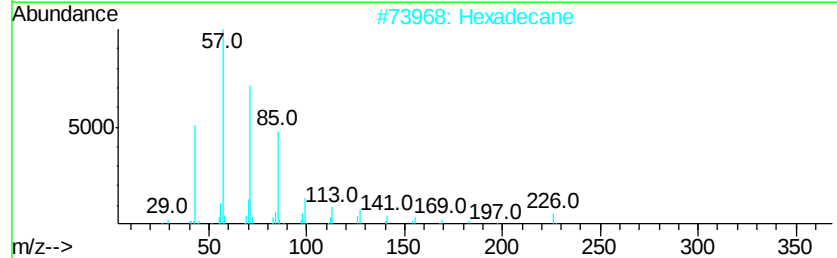
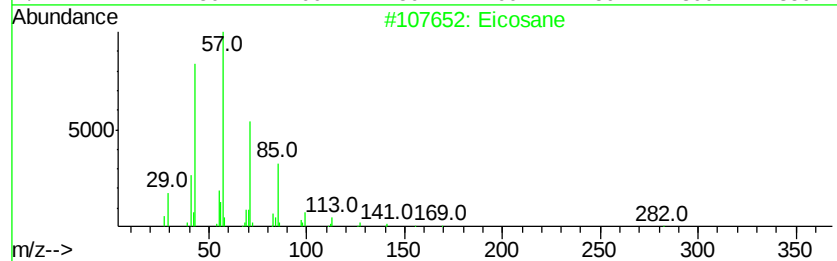
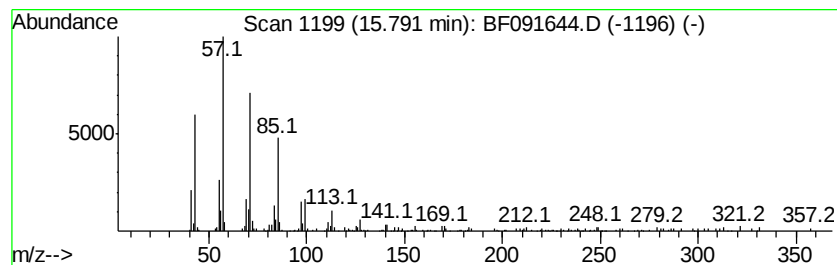
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.79	3.37 ng	206798	Perylene-d12	15.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	93
2			Hexadecane	226	C16H34	000544-76-3	91
3			Heneicosane	296	C21H44	000629-94-7	90
4			Tetradecane, 2-methyl-	212	C15H32	001560-95-8	90
5			Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121516\
Data File : BF091644.D
Acq On : 15 Dec 2016 22:32
Operator : UM/SJ
Sample : H5940-05
Misc :
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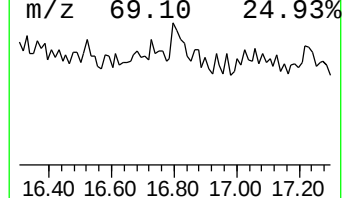
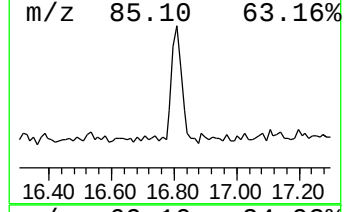
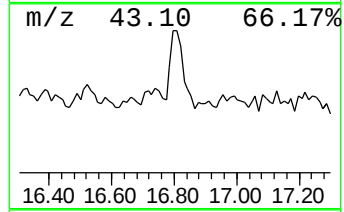
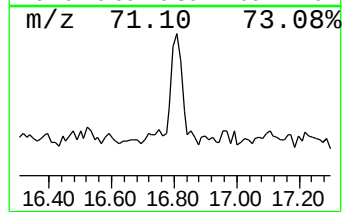
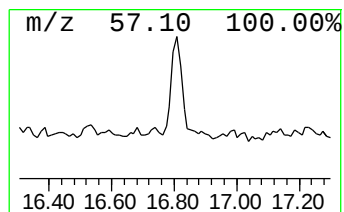
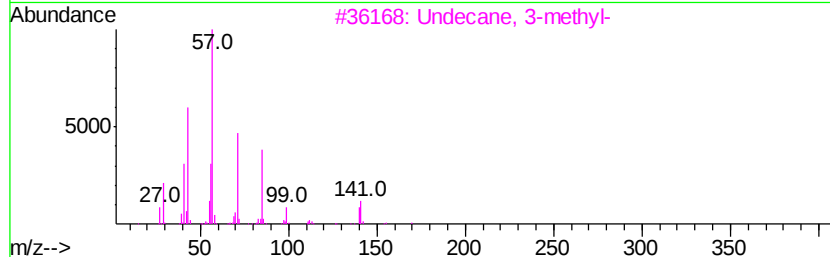
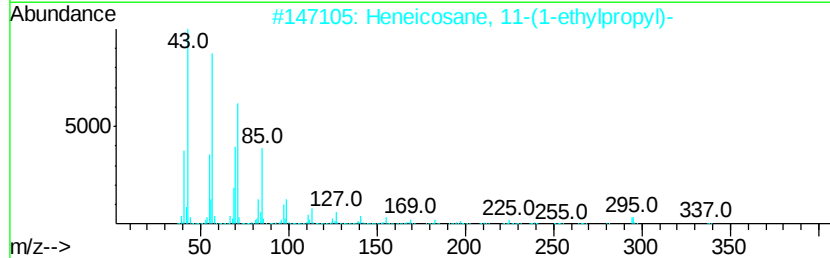
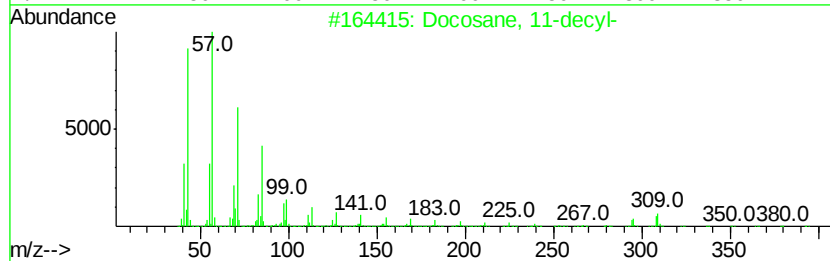
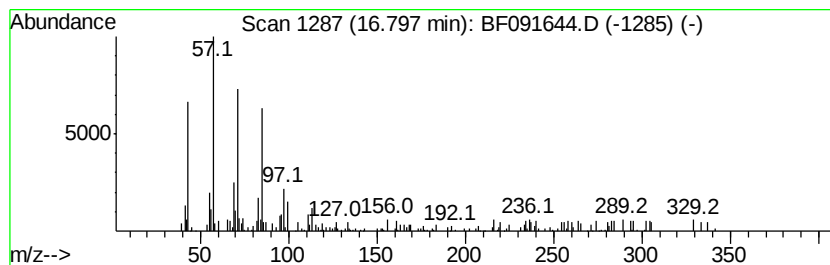
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Docosane, 11-decyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.80	3.26 ng	199962	Perylene-d12	15.35	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 11-decyl-	451	C32H66	055401-55-3	53
2	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	53
3	Undecane, 3-methyl-	170	C12H26	001002-43-3	47
4	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	47
5	Octadecane	254	C18H38	000593-45-3	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121516\
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Acq On : 15 Dec 2016 22:32
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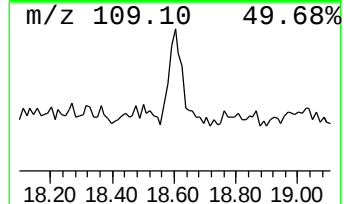
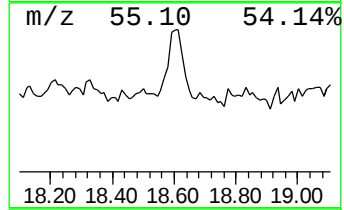
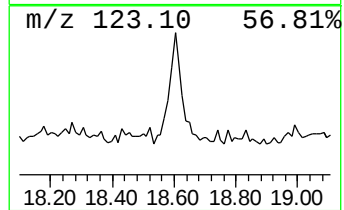
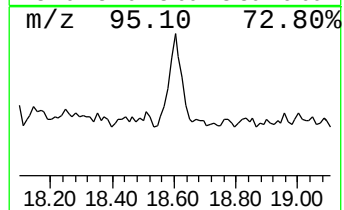
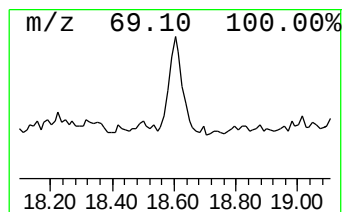
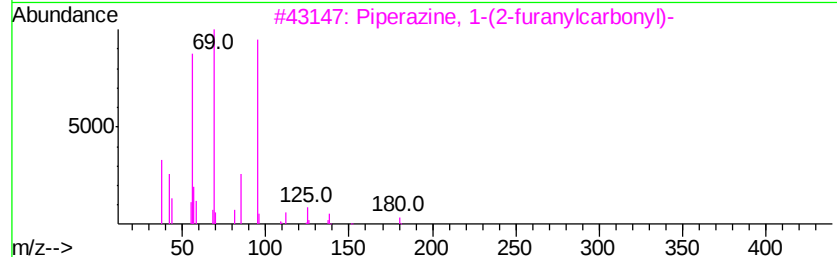
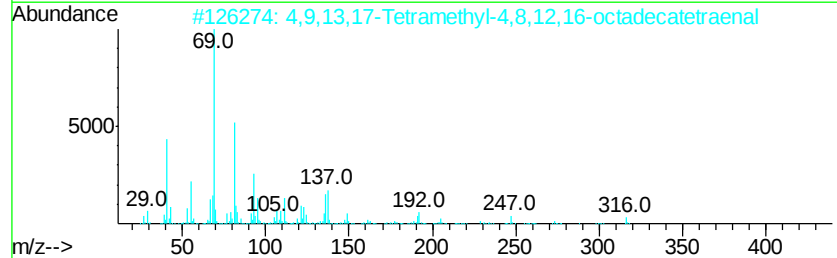
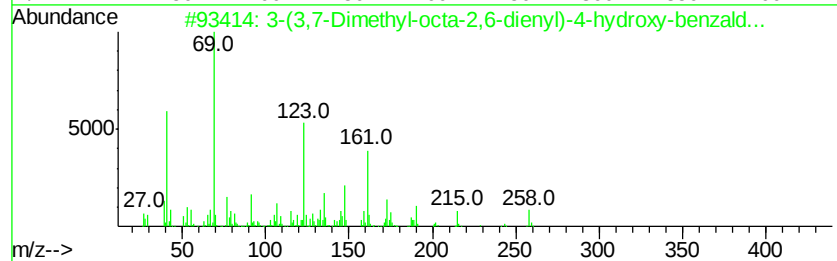
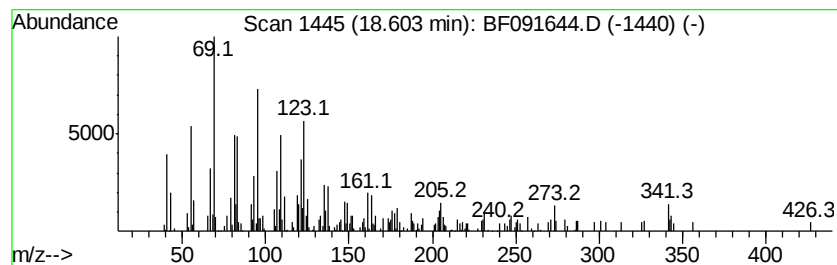
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown18.60 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.60	5.23 ng	320945	Perylene-d12	15.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3		(3,7-Dimethyl-octa-2,6-dienvl)...	258	C17H22O2	1000192-28-0	35
2	4,9,13,17		Tetramethyl-4,8,12,16-...	316	C22H36O	056882-09-8	35
3	Piperazine, 1-(2-furanvlcarbonyl)-			180	C9H12N2O2	040172-95-0	27
4	1,6,10,14,18,22-Tetracosahexaen-...			426	C30H50O	054159-46-5	25
5	Trichloro(2-pentenyl)silane			202	C5H9Cl3Si	1000109-99-7	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121516\
Data File : BF091644.D
Acq On : 15 Dec 2016 22:32
Operator : UM/SJ
Sample : H5940-05
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120916.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...	4.96	29.1 ng		2455150	1	6.78	1684660	20.0
unknown6.56	6.56	78.4 ng		6603320	1	6.78	1684660	20.0
Cyclohexene, 1-me...	6.91	7.3 ng		619064	1	6.78	1684660	20.0
unknown10.54	10.54	3.6 ng		402214	3	9.82	2253860	20.0
Heptadecane	10.74	4.0 ng		330298	4	11.30	1635670	20.0
n-Hexadecanoic acid	11.84	2.8 ng		231925	4	11.30	1635670	20.0
unknown12.42	12.42	4.8 ng		396296	4	11.30	1635670	20.0
3-Eicosene, (E)-	13.79	5.5 ng		438811	5	13.94	1587120	20.0
2-Bromo dodecane	15.04	3.3 ng		204502	6	15.35	1226150	20.0
Eicosane	15.79	3.4 ng		206798	6	15.35	1226150	20.0
Docosane, 11-decyl-	16.80	3.3 ng		199962	6	15.35	1226150	20.0
unknown18.60	18.60	5.2 ng		320945	6	15.35	1226150	20.0