

Data Path : Z:\HPCHEM1\BNA F\DATA\BF030317\
 Data File : BF093381.D
 Acq On : 4 Mar 2017 1:35
 Operator : SJ/MA
 Sample : I2078-07
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C-4

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-BF013017.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	3.207	95	98	109	rVB	88747	190932	2.33%	0.403%
2	4.258	186	190	193	rBV	4046918	7454053	90.98%	15.716%
3	4.956	246	251	254	rBV	3780640	8193110	100.00%	17.274%
4	5.390	287	289	301	rBV	243647	424680	5.18%	0.895%
5	6.430	378	380	388	rBV	2439421	2485054	30.33%	5.239%
6	6.556	388	391	393	rVV	776394	861878	10.52%	1.817%
7	6.784	409	411	413	rBV	1223952	1276620	15.58%	2.692%
8	6.853	415	417	419	rBV	115174	115594	1.41%	0.244%
9	6.933	422	424	427	rBV	2655963	3328050	40.62%	7.017%
10	7.116	438	440	445	rBV	69036	100209	1.22%	0.211%
11	7.356	458	461	463	rBV	2969128	2829510	34.54%	5.966%
12	8.076	521	524	526	rBV	1440992	1650396	20.14%	3.480%
13	9.150	615	618	620	rBV	4678299	4669788	57.00%	9.846%
14	9.539	650	652	655	rBV	300122	379656	4.63%	0.800%
15	9.756	669	671	673	rBV	113170	138652	1.69%	0.292%
16	9.825	674	677	679	rVV	1627023	1653844	20.19%	3.487%
17	10.248	713	714	716	rVB	164244	126392	1.54%	0.266%
18	10.465	730	733	736	rBV	62562	98851	1.21%	0.208%
19	10.716	754	755	760	rVB	161031	282081	3.44%	0.595%
20	11.162	792	794	796	rBV	116884	147244	1.80%	0.310%
21	11.299	804	806	810	rBV	1280499	1689367	20.62%	3.562%
22	11.596	830	832	834	rBV	89960	87929	1.07%	0.185%
23	11.665	836	838	840	rVB	186450	155728	1.90%	0.328%
24	11.859	854	855	858	rVB	702883	830611	10.14%	1.751%
25	11.928	858	861	864	rVB	226408	272112	3.32%	0.574%
26	12.076	873	874	878	rVB	66149	85080	1.04%	0.179%
27	12.248	886	889	890	rBV	96466	133028	1.62%	0.280%
28	12.522	911	913	915	rBV2	128164	184897	2.26%	0.390%
29	12.751	930	933	936	rBV	91332	160880	1.96%	0.339%
30	12.888	942	945	947	rBV	3386106	3280290	40.04%	6.916%
31	13.779	1020	1023	1028	rVB2	218177	317569	3.88%	0.670%
32	13.939	1035	1037	1040	rBV	966979	1001629	12.23%	2.112%
33	14.408	1076	1078	1083	rBV	73471	151374	1.85%	0.319%
34	14.694	1101	1103	1105	rBV2	95190	124886	1.52%	0.263%

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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	15.002	1128	1130	1134	rVB	116587	194312	2.37%	0.410%
36	15.357	1158	1161	1164	rBV	952271	1166851	14.24%	2.460%
37	15.757	1193	1196	1198	rBV	204386	288265	3.52%	0.608%
38	16.214	1233	1236	1242	rVB	194173	357169	4.36%	0.753%
39	16.751	1280	1283	1286	rVB	142038	285734	3.49%	0.602%
40	17.380	1335	1338	1343	rVB	107588	255227	3.12%	0.538%

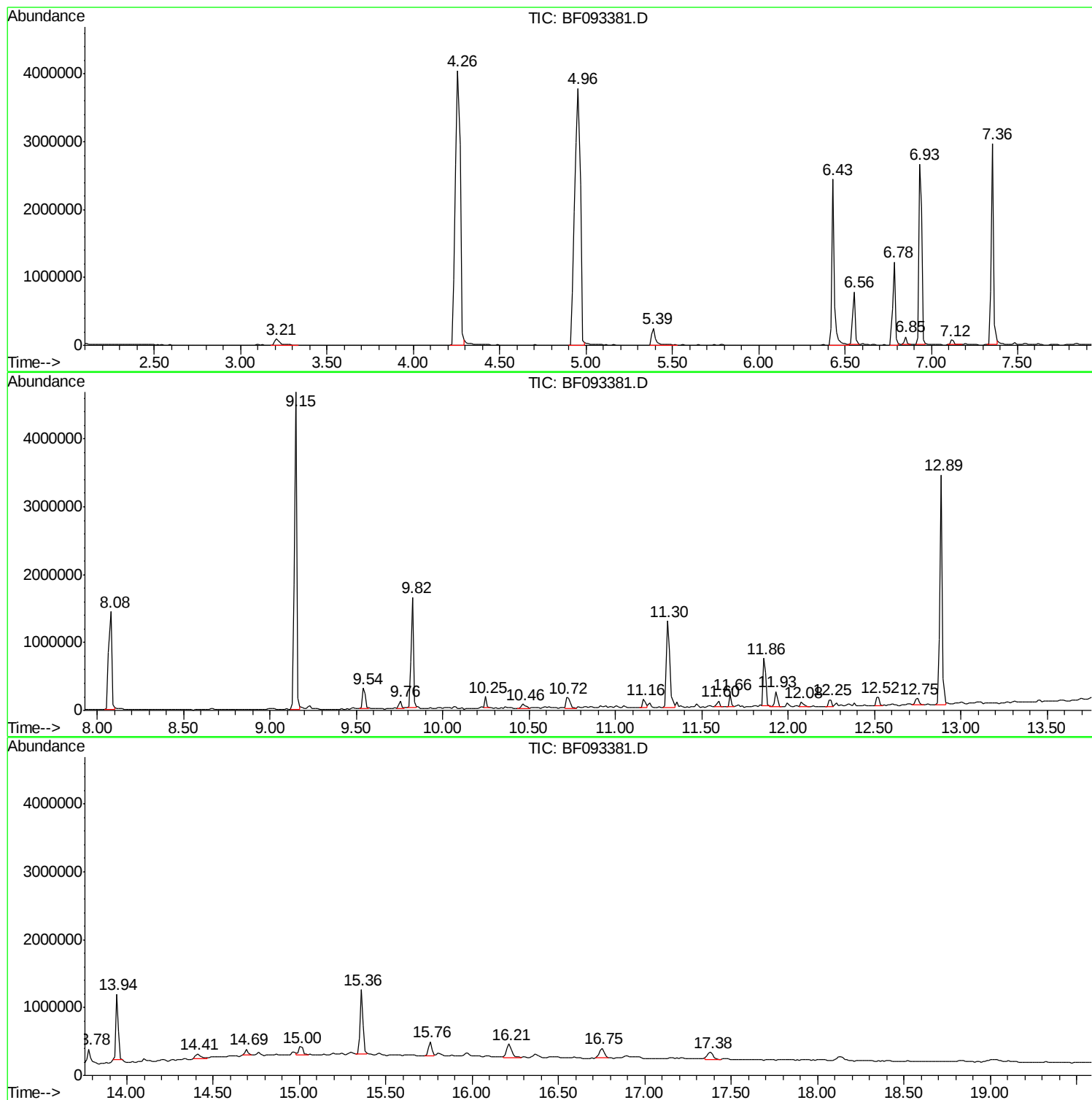
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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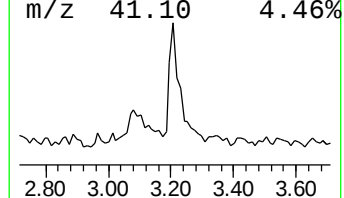
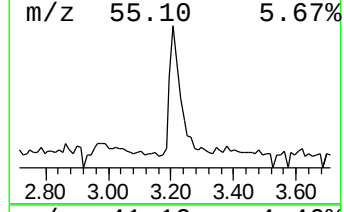
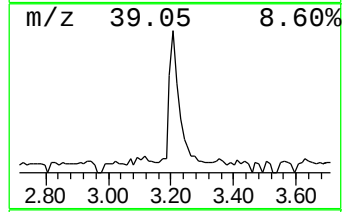
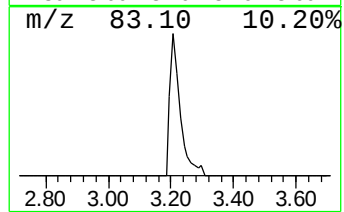
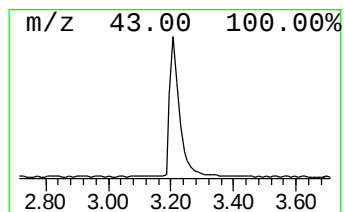
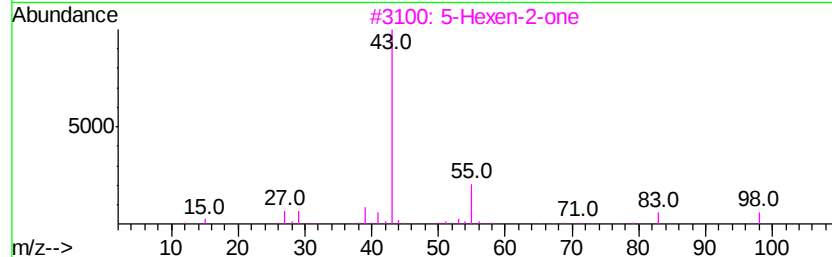
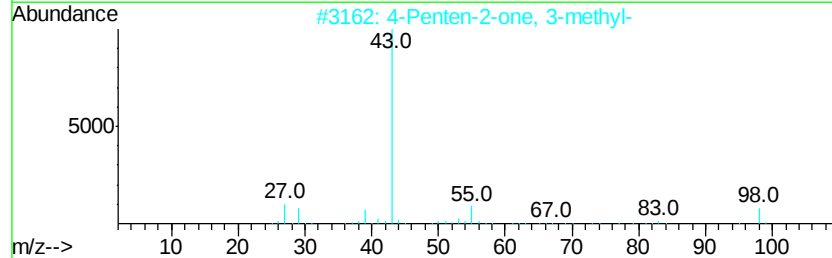
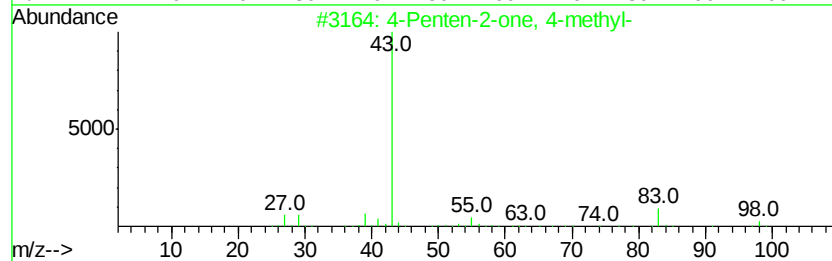
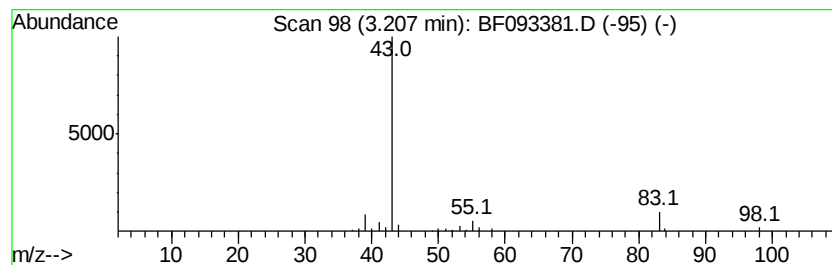
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 4-Penten-2-one, 4-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.21	2.99 ng	190932	1,4-Dichlorobenzene-d4	6.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	80
2		4-Penten-2-one, 3-methyl-	98	C6H10O	000758-87-2	50
3		5-Hexen-2-one	98	C6H10O	000109-49-9	47
4		2-Pentanone, 3-methylene-	98	C6H10O	004359-77-7	9
5		4-Hexen-2-one	98	C6H10O	025659-22-7	9



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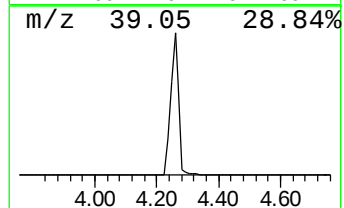
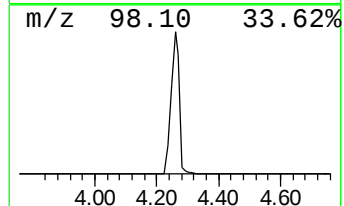
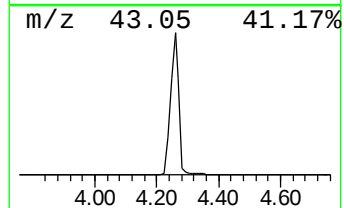
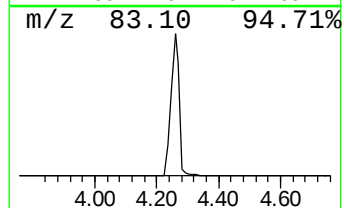
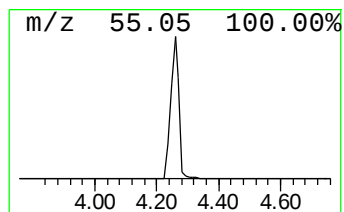
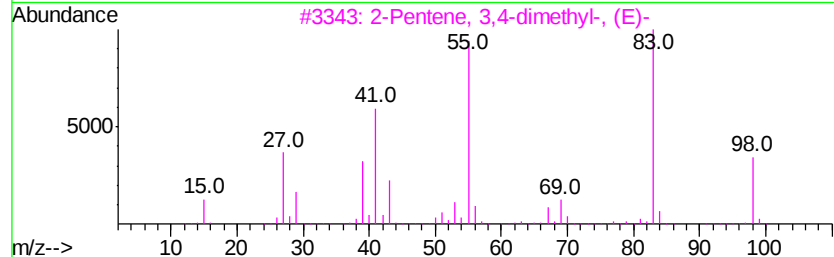
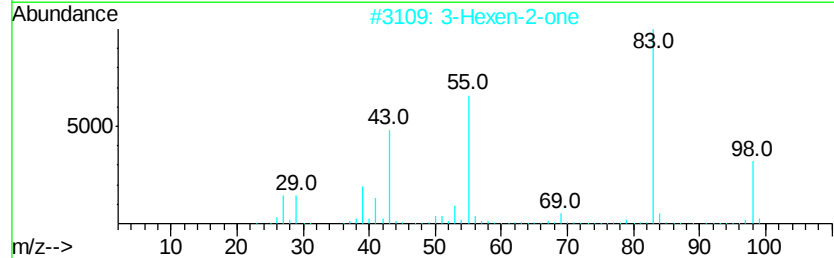
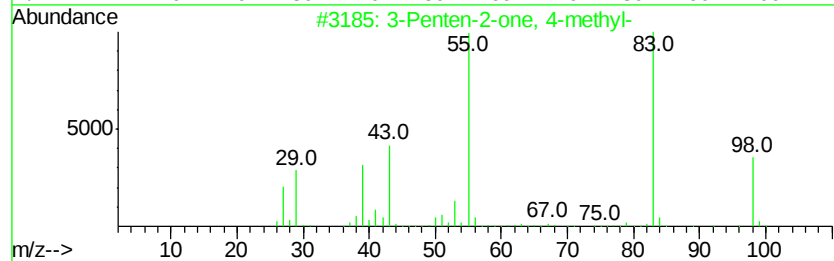
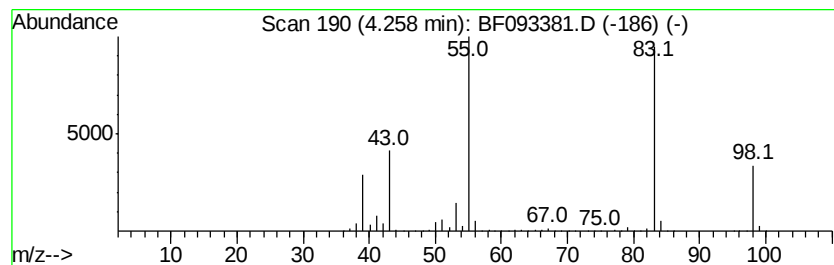
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.26	116.78 ng	7454050	1,4-Dichlorobenzene-d4	6.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
2		3-Hexen-2-one	98	C6H10O	000763-93-9	90
3		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4		2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	86
5		2-Pentene, 4,4-dimethyl-, (Z)-	98	C7H14	000762-63-0	86



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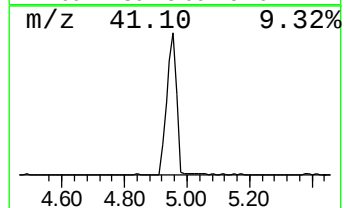
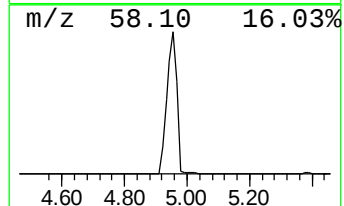
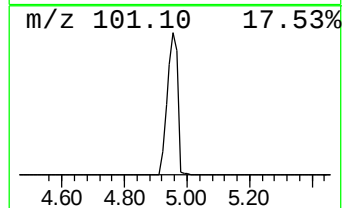
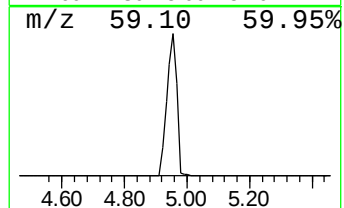
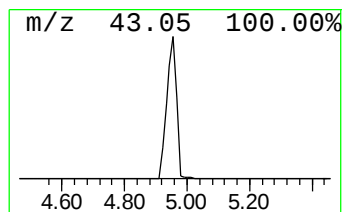
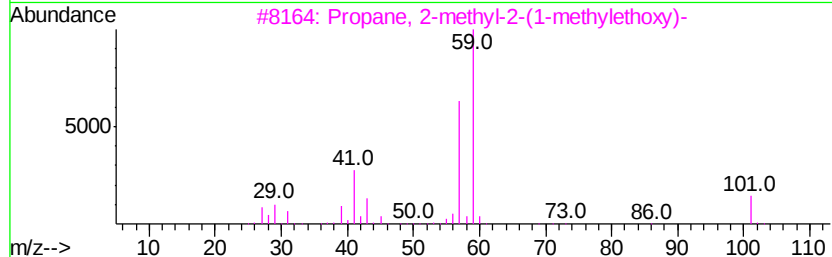
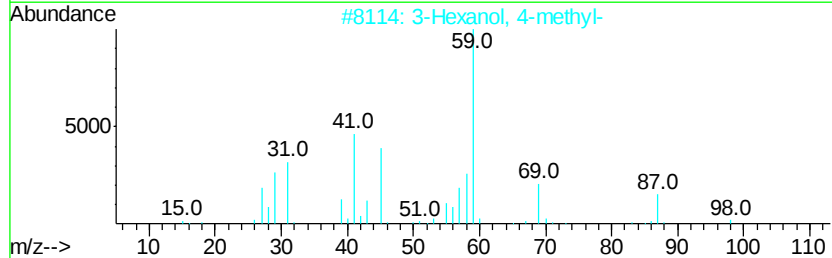
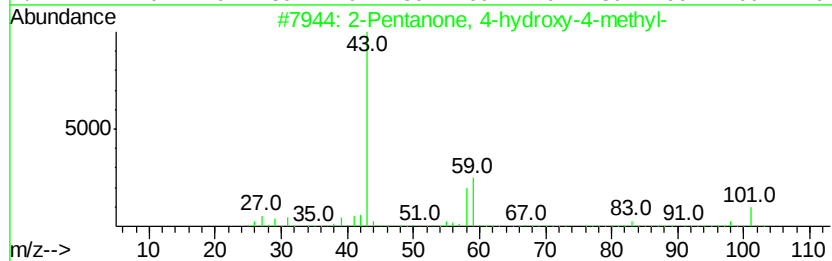
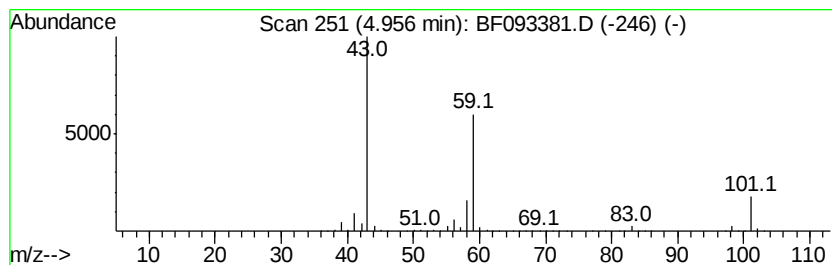
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	128.36 ng	8193110	1,4-Dichlorobenzene-d4	6.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	33
4		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	32
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28



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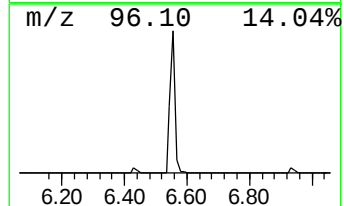
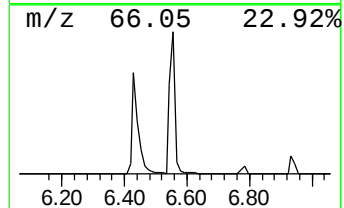
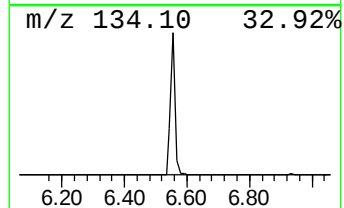
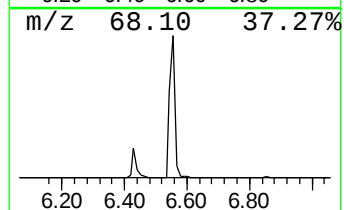
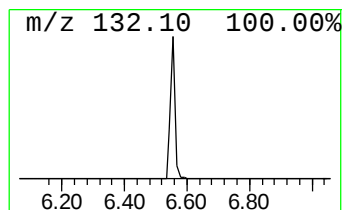
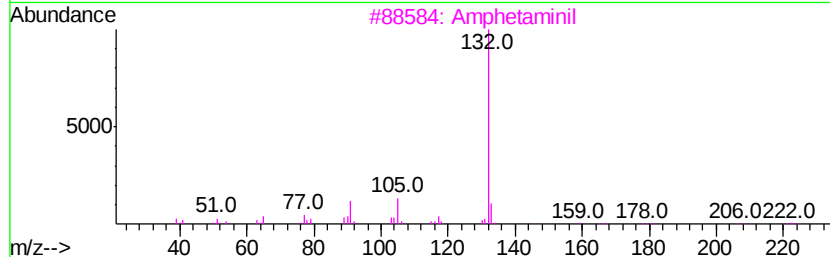
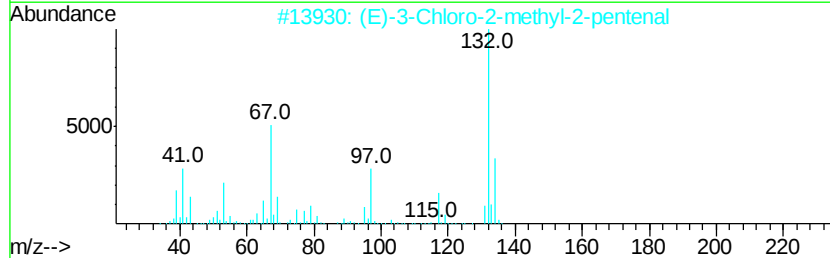
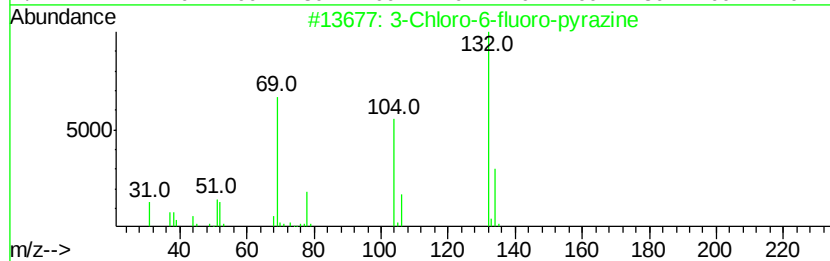
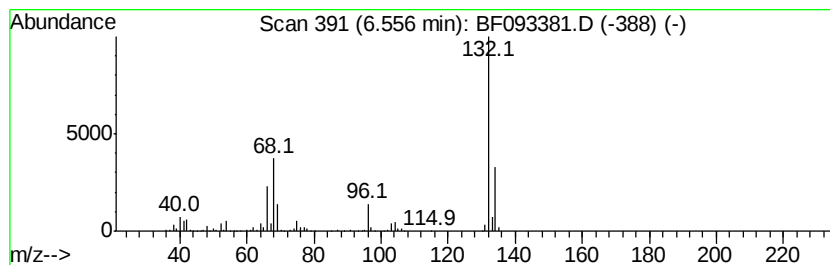
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 unknown6.56 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	13.50 ng	861878	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	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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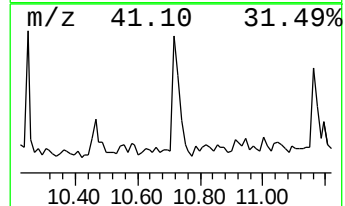
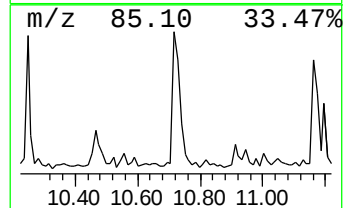
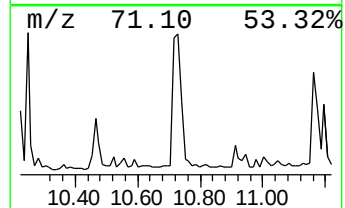
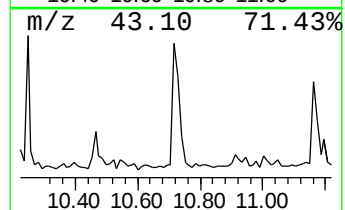
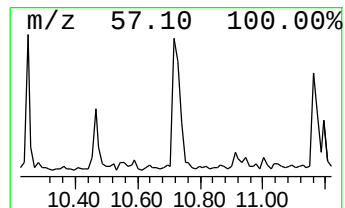
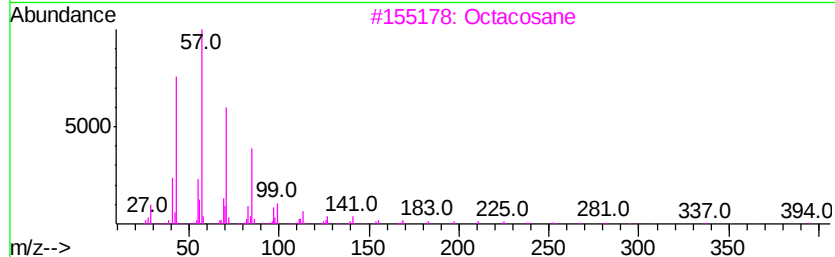
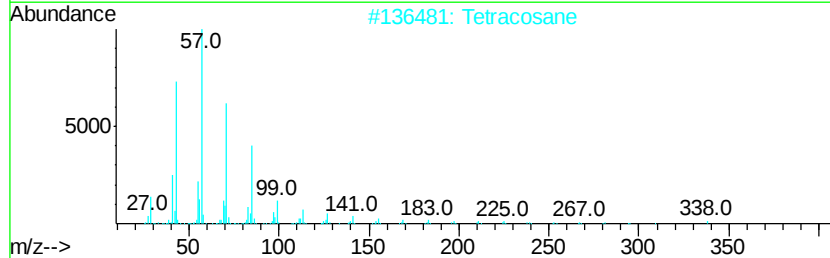
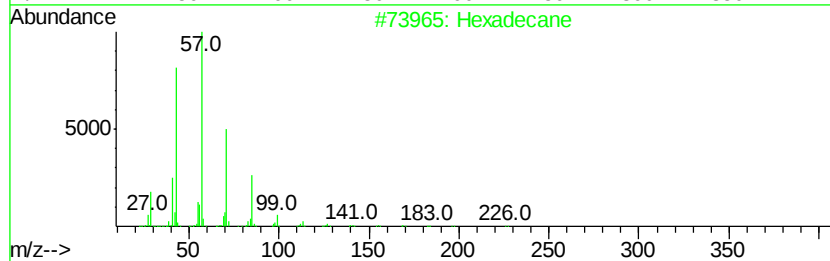
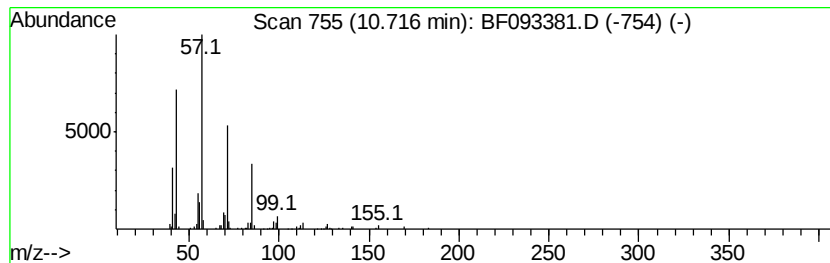
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.72	3.34 ng	282081	Phenanthrene-d10	11.30

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	91
2	Tetracosane		338	C24H50	000646-31-1	78
3	Octacosane		394	C28H58	000630-02-4	78
4	Nonane, 5-butyl-		184	C13H28	017312-63-9	72
5	Pentadecane		212	C15H32	000629-62-9	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF030317\
Data File : BF093381.D
Acq On : 4 Mar 2017 1:35
Operator : SJ/MA
Sample : I2078-07
Misc :
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Instrument :
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ClientSampled :
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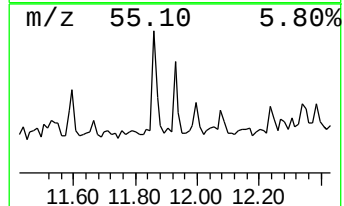
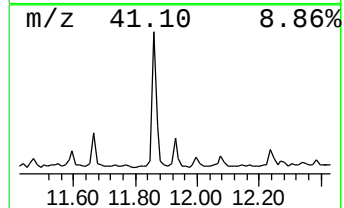
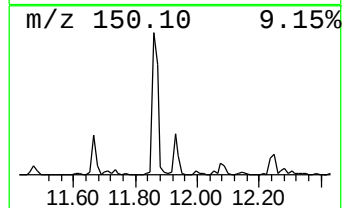
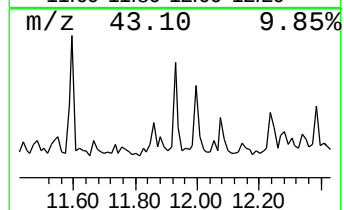
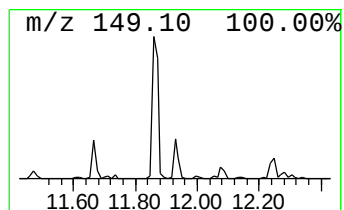
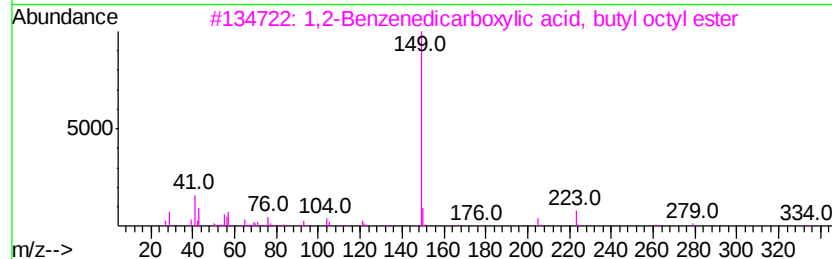
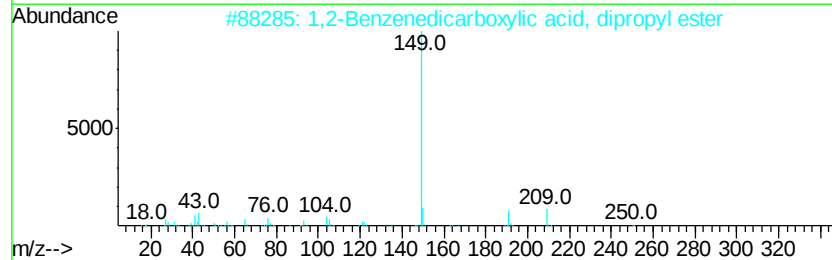
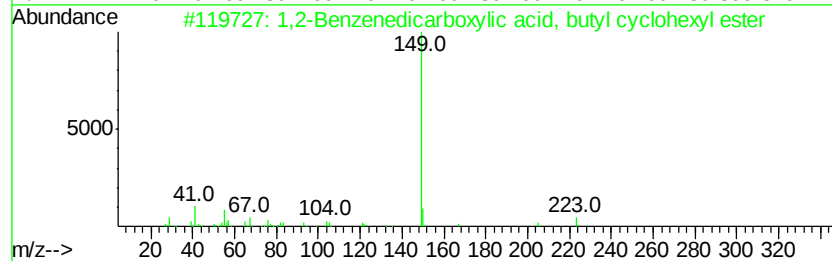
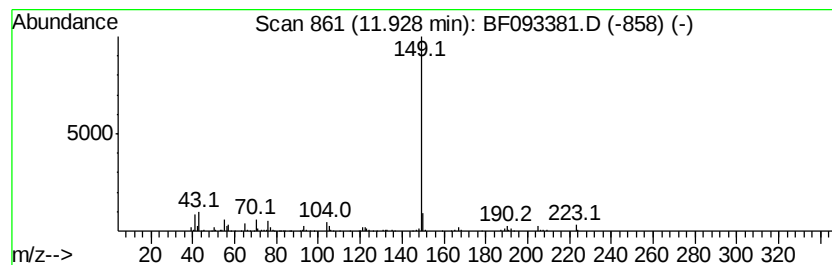
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1,2-Benzenedicarboxylic aci... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	3.22 ng	272112	Phenanthrene-d10	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2-Benzenedicarboxylic acid, bu...	304	C18H24O4	000084-64-0	86
2		1,2-Benzenedicarboxylic acid, di...	250	C14H18O4	000131-16-8	86
3		1,2-Benzenedicarboxylic acid, bu...	334	C20H30O4	000084-78-6	78
4		1,2-Benzenedicarboxylic acid, bu...	362	C22H34O4	000089-18-9	78
5		Dibutyl phthalate	278	C16H22O4	000084-74-2	78



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030317\
Data File : BF093381.D
Acq On : 4 Mar 2017 1:35
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Misc :
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Instrument :
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ClientSampleId :
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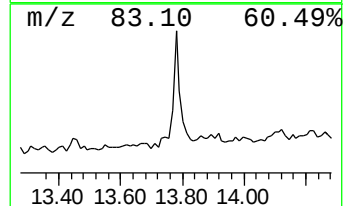
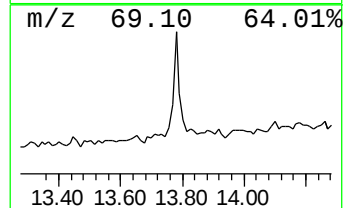
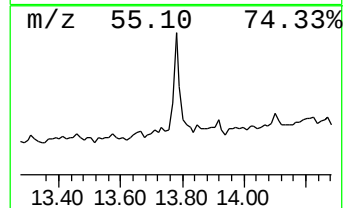
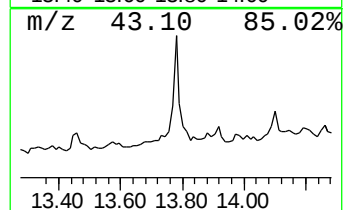
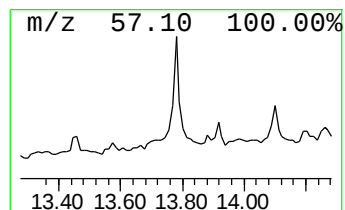
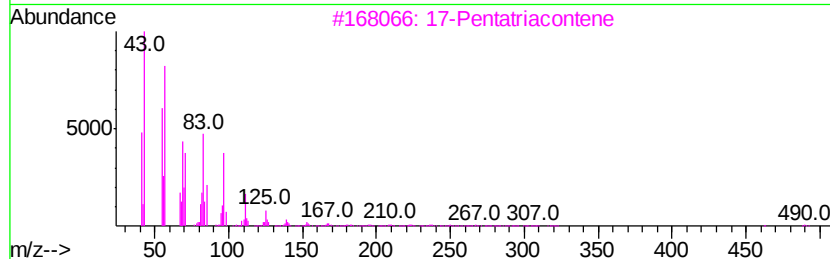
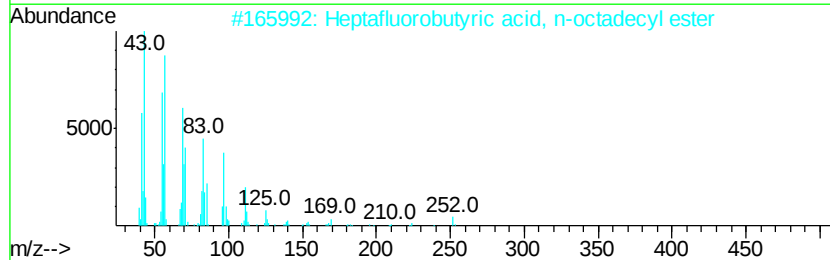
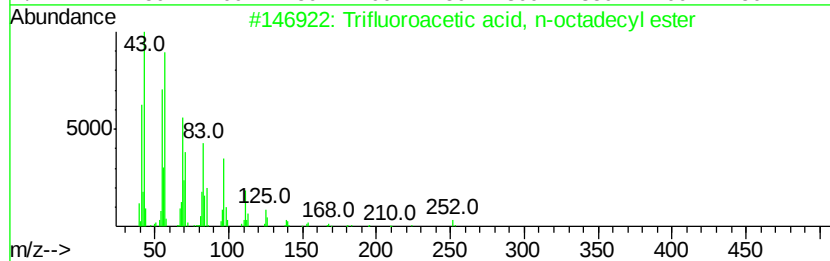
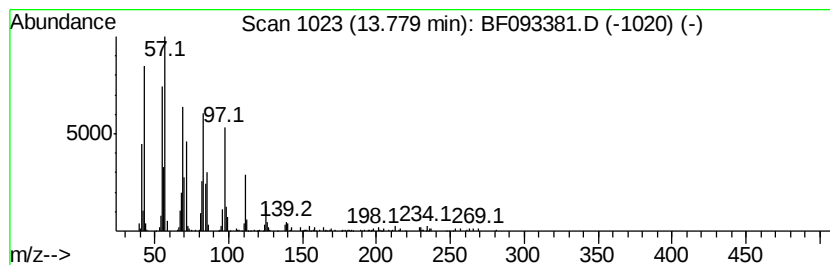
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Trifluoroacetic acid, n-oct... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.78	6.34 ng	317569	Chrysene-d12	13.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trifluoroacetic acid, n-octadecy...	366	C20H37F3O2	079392-43-1	91
2		Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	91
3		17-Pentatriacontene	491	C35H70	006971-40-0	90
4		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	90
5		Dichloroacetic acid, tetradecyl ...	324	C16H30Cl2O2	083005-02-1	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030317\
Data File : BF093381.D
Acq On : 4 Mar 2017 1:35
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Misc :
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Instrument :
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ClientSampleId :
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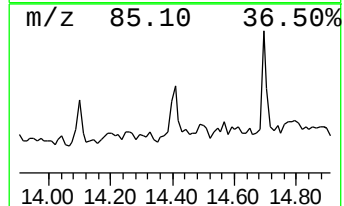
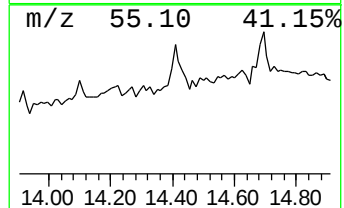
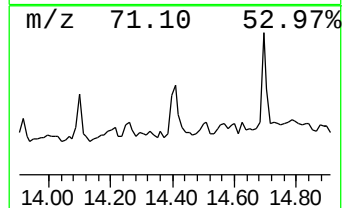
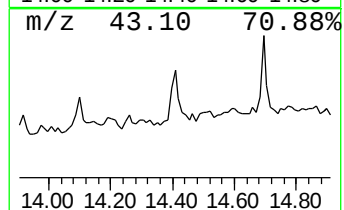
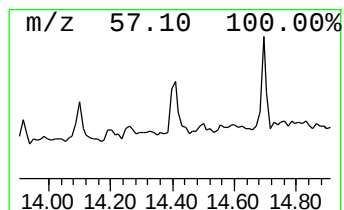
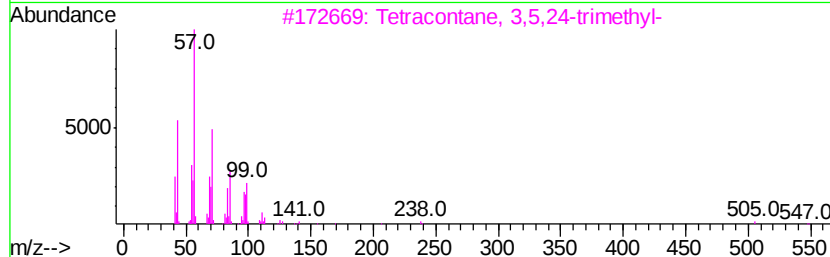
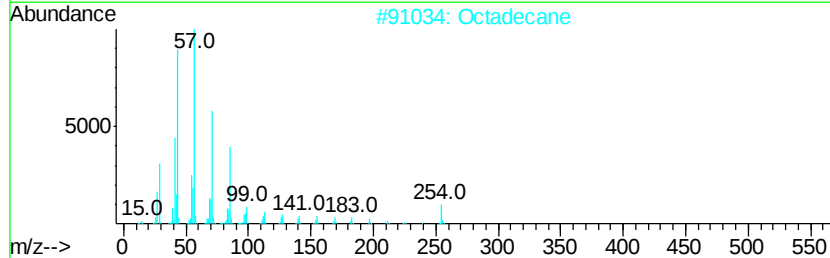
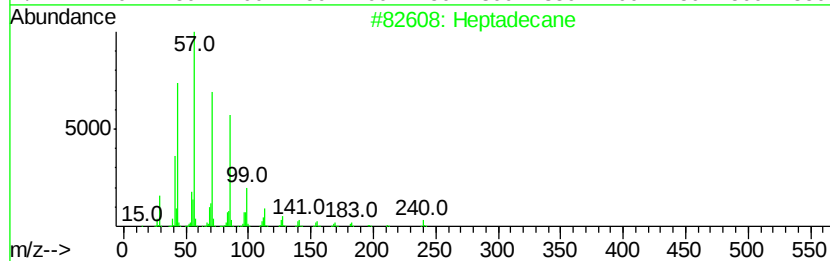
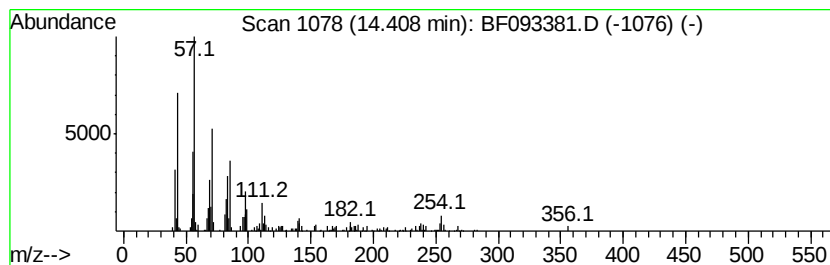
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heptadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.41	3.02 ng	151374	Chrysene-d12	13.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	90
2		Octadecane	254	C18H38	000593-45-3	87
3		Tetracontane, 3,5,24-trimethyl-	605	C43H88	055162-61-3	74
4		Nonadecane	268	C19H40	000629-92-5	70
5		Octadecane, 1-bromo-	332	C18H37Br	000112-89-0	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF030317\
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Acq On : 4 Mar 2017 1:35
Operator : SJ/MA
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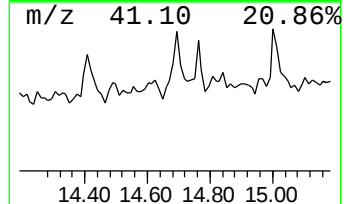
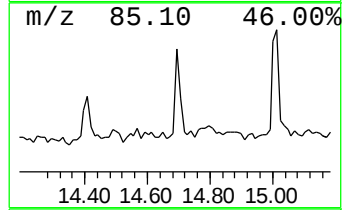
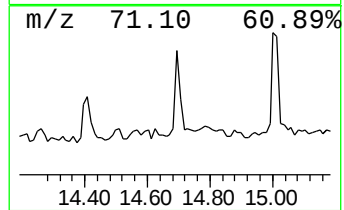
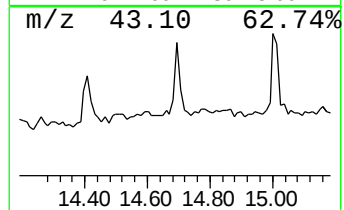
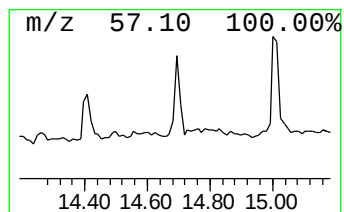
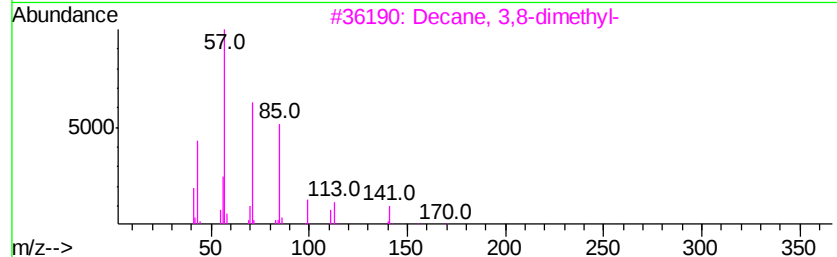
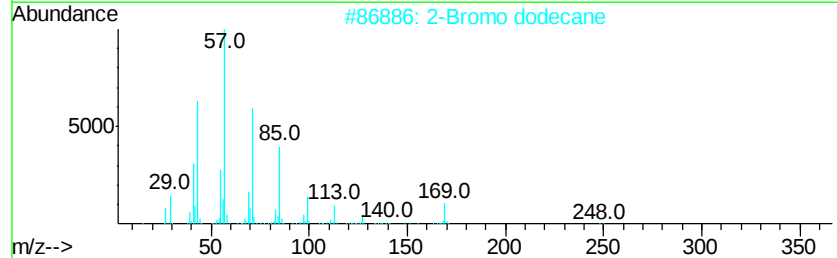
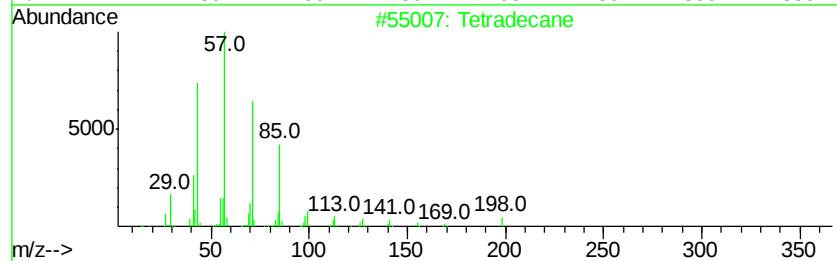
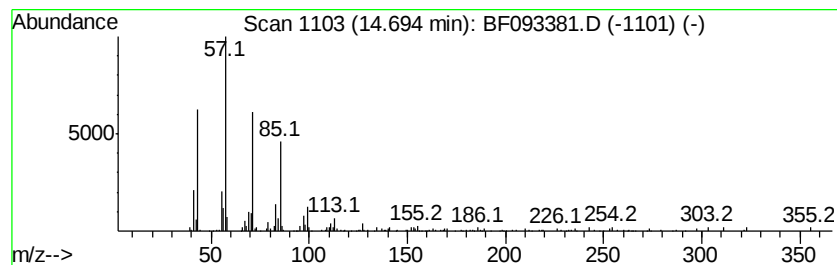
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tetradecane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	2.14 ng	124886	Perylene-d12	15.36
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tetradecane	198 C14H30	000629-59-4 87
2		2-Bromo dodecane	248 C12H25Br	013187-99-0 83
3		Decane, 3,8-dimethyl-	170 C12H26	017312-55-9 80
4		Eicosane, 10-methyl-	296 C21H44	054833-23-7 80
5		Hexadecane	226 C16H34	000544-76-3 78



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030317\
Data File : BF093381.D
Acq On : 4 Mar 2017 1:35
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Misc :
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Instrument :
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ClientSampleId :
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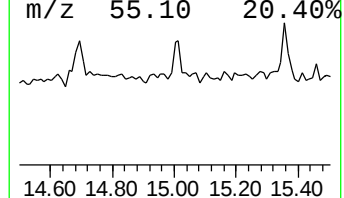
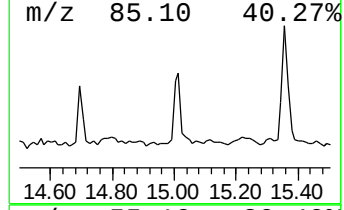
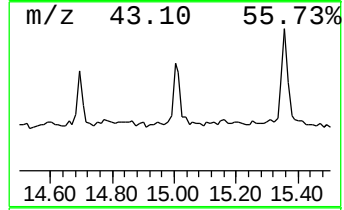
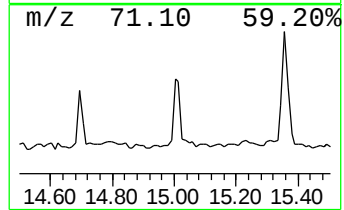
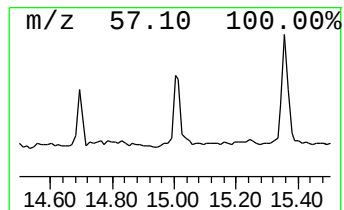
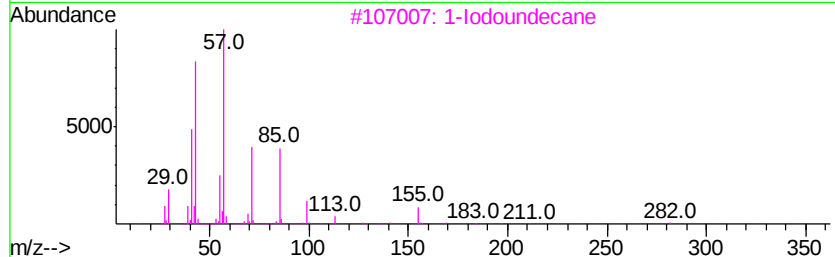
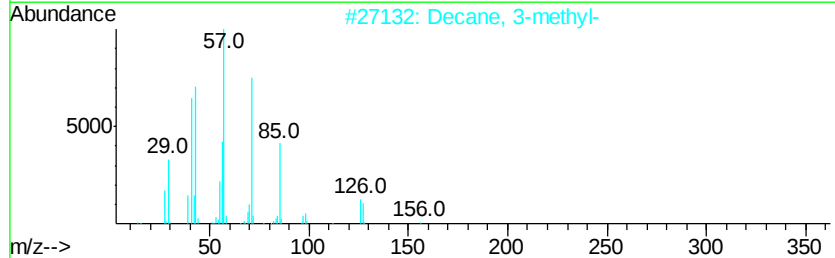
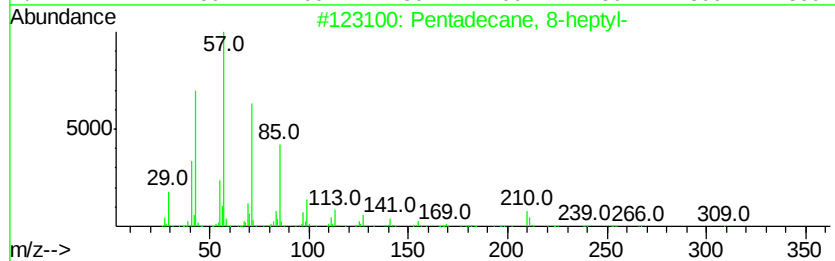
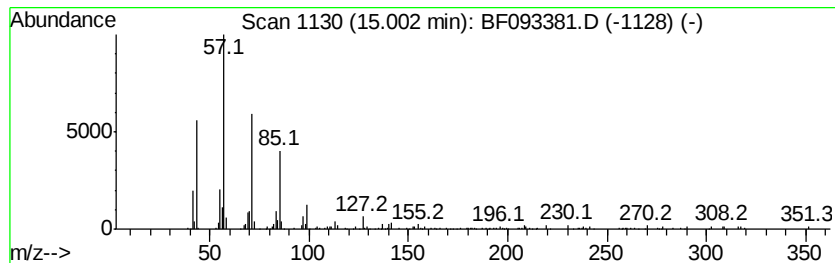
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Pentadecane, 8-heptyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.00	3.33 ng	194312	Perylene-d12	15.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	80
2		Decane, 3-methyl-	156	C11H24	013151-34-3	72
3		1-Iodoundecane	282	C11H23I	004282-44-4	72
4		Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	72
5		Hexadecane	226	C16H34	000544-76-3	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030317\
Data File : BF093381.D
Acq On : 4 Mar 2017 1:35
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Misc :
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ClientSampleId :
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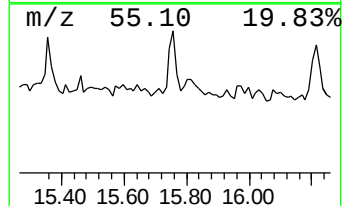
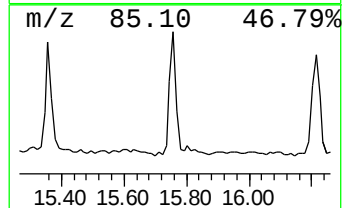
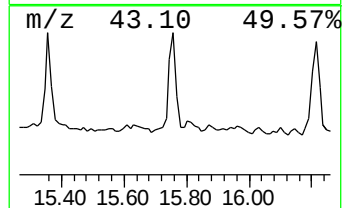
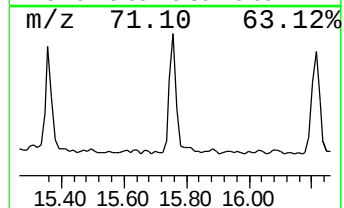
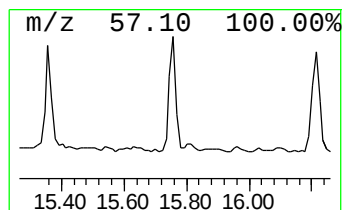
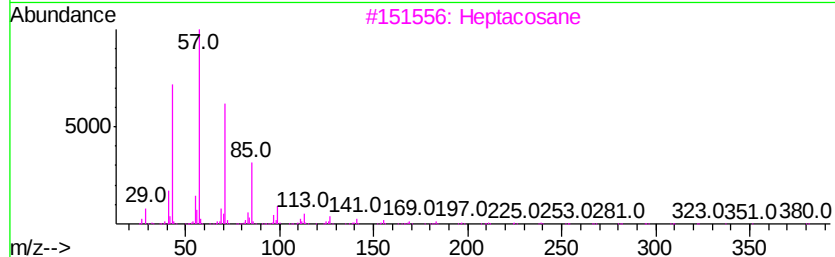
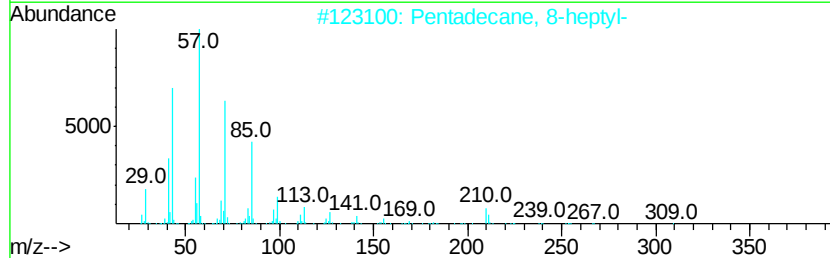
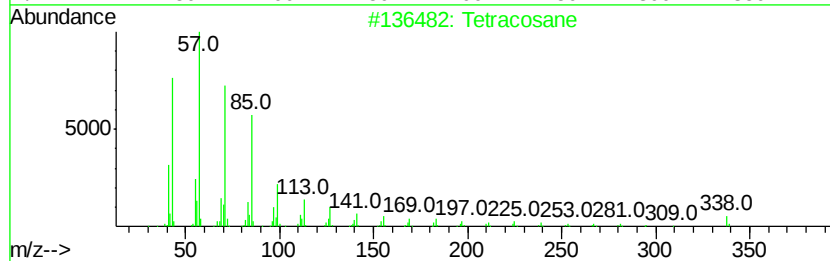
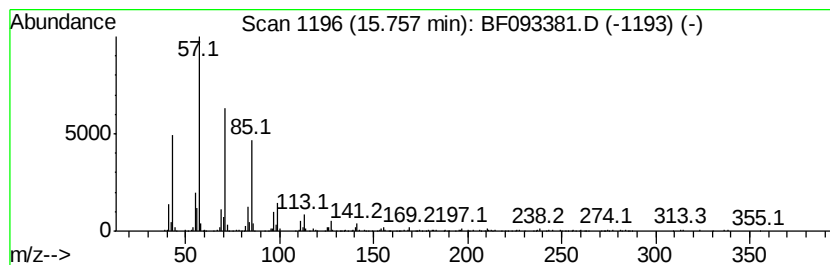
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Tetracosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.76	4.94 ng	288265	Perylene-d12	15.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	95
2		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		Tetratetracontane	619	C44H90	007098-22-8	90
5		Hexadecane	226	C16H34	000544-76-3	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030317\
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Acq On : 4 Mar 2017 1:35
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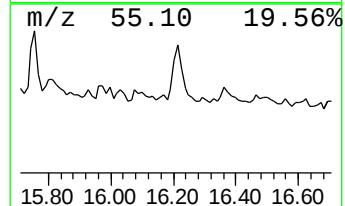
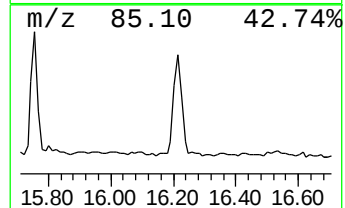
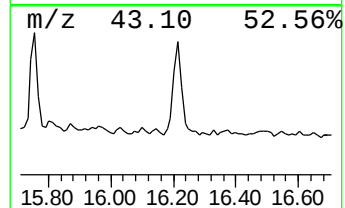
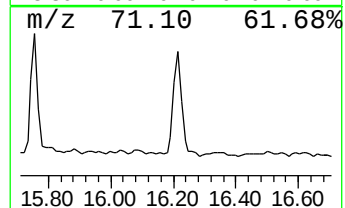
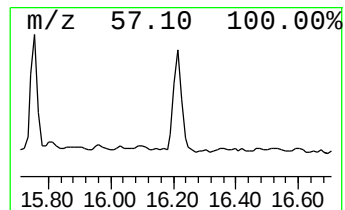
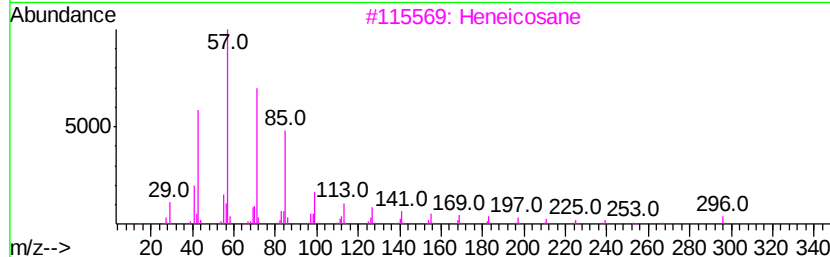
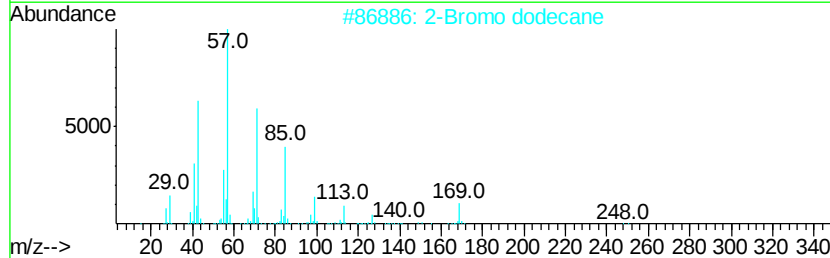
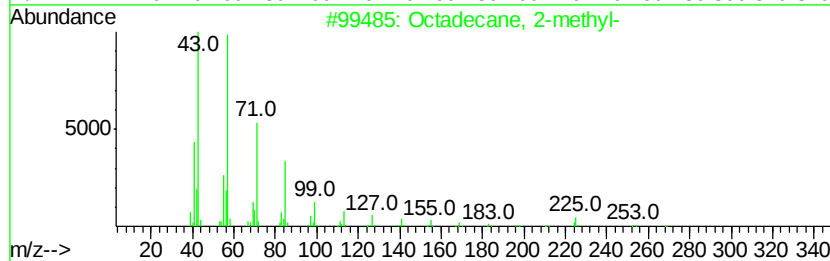
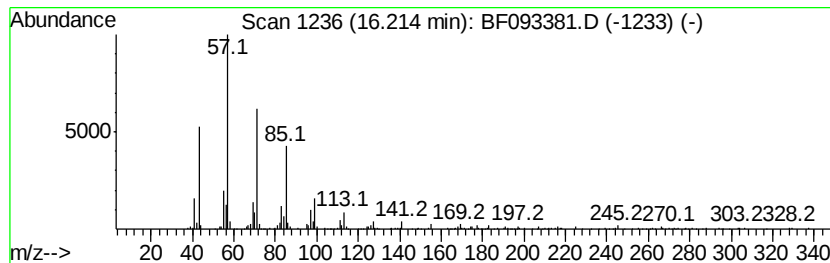
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Octadecane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.21	6.12 ng	357169	Perylene-d12	15.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane, 2-methyl-	268	C19H40	001560-88-9	93
2			2-Bromo dodecane	248	C12H25Br	013187-99-0	91
3			Heneicosane	296	C21H44	000629-94-7	91
4			Tetratetracontane	619	C44H90	007098-22-8	90
5			Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030317\
Data File : BF093381.D
Acq On : 4 Mar 2017 1:35
Operator : SJ/MA
Sample : I2078-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
C-4

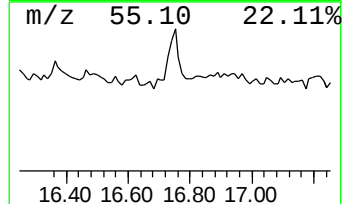
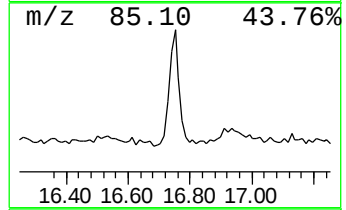
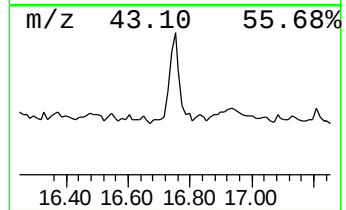
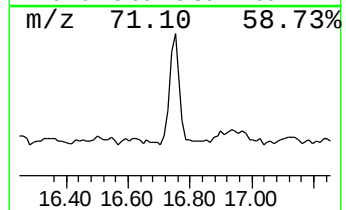
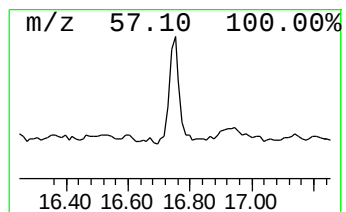
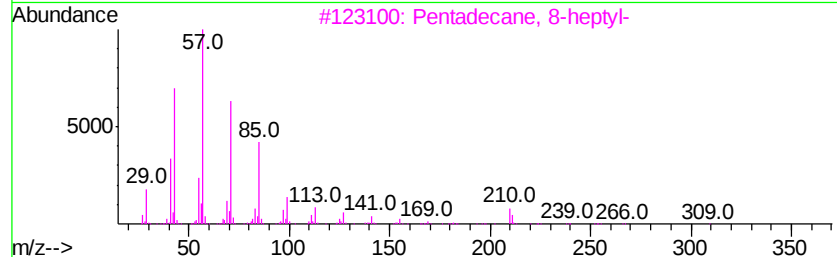
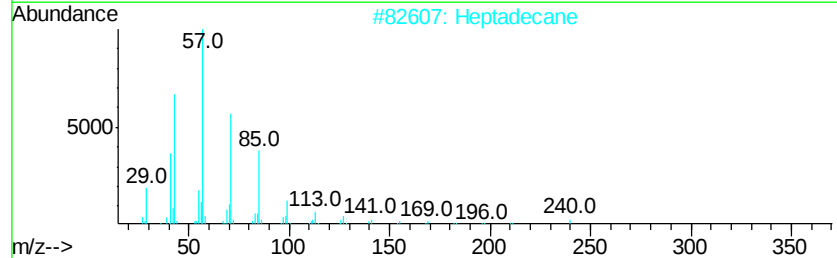
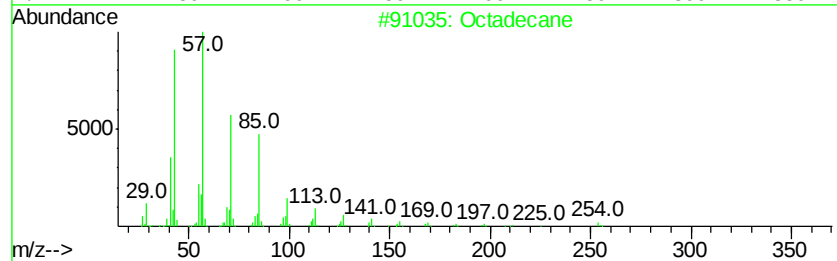
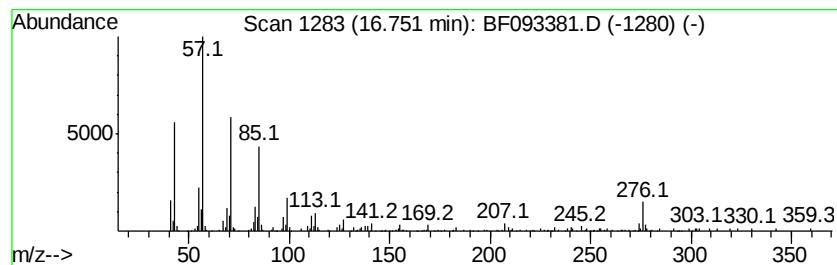
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Octadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.75	4.90 ng	285734	Perylene-d12	15.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	89
2		Heptadecane	240	C17H36	000629-78-7	87
3		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	83
4		Docosane	310	C22H46	000629-97-0	83
5		Heptacosane	380	C27H56	000593-49-7	83



Data Path : Z:\HPCHEM1\BNA F\DATA\BF030317\
Data File : BF093381.D
Acq On : 4 Mar 2017 1:35
Operator : SJ/MA
Sample : I2078-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
C-4

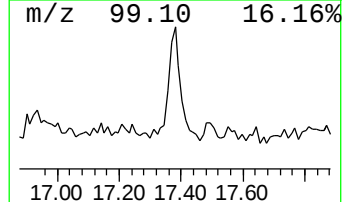
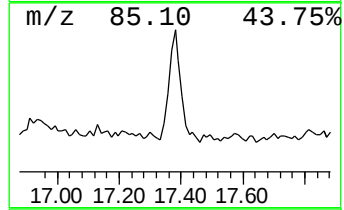
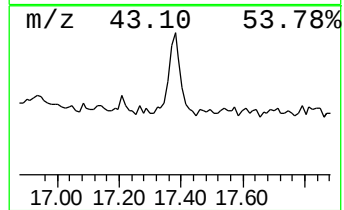
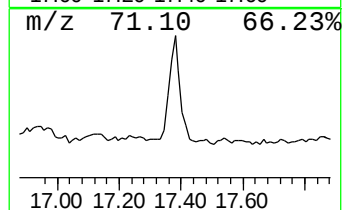
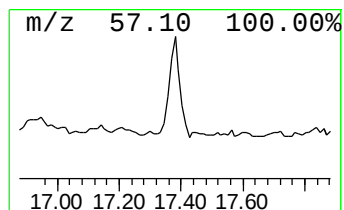
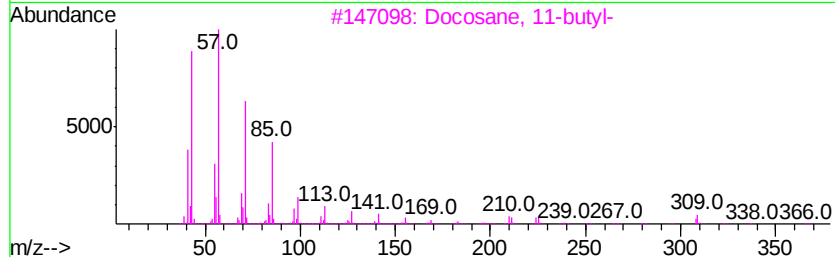
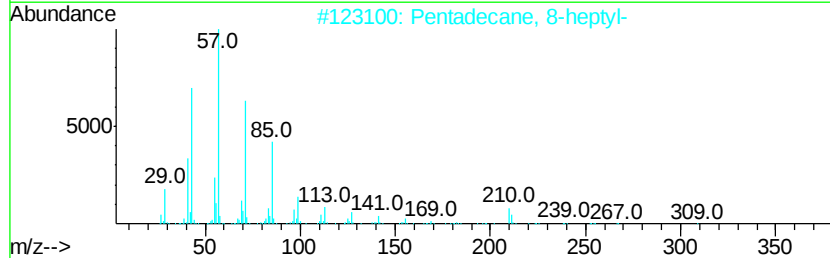
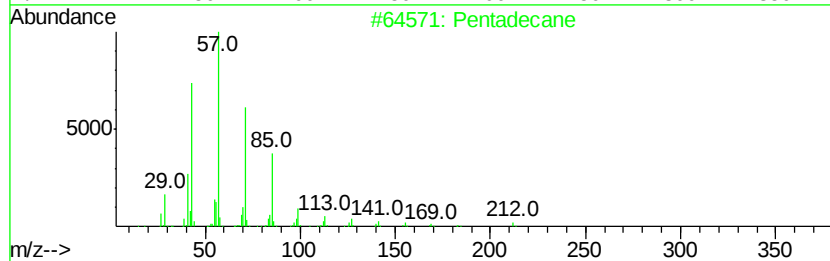
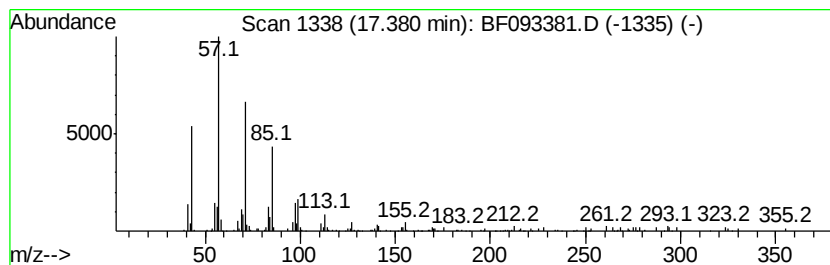
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Pentadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.38	4.37 ng	255227	Perylene-d12	15.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane	212	C15H32	000629-62-9	91
2		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	86
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	86
4		Heptacosane	380	C27H56	000593-49-7	83
5		Octadecane, 2-methyl-	268	C19H40	001560-88-9	80



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 Data File : BF093381.D
 Acq On : 4 Mar 2017 1:35
 Operator : SJ/MA
 Sample : I2078-07
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.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
4-Penten-2-one, 4...	3.21	3.0	ng	190932	1	6.78	1276620	20.0
3-Penten-2-one, 4...	4.26	116.8	ng	7454050	1	6.78	1276620	20.0
2-Pentanone, 4-hy...	4.96	128.4	ng	8193110	1	6.78	1276620	20.0
unknown6.56	6.56	13.5	ng	861878	1	6.78	1276620	20.0
Hexadecane	10.72	3.3	ng	282081	4	11.30	1689370	20.0
1,2-Benzenedicarb...	11.93	3.2	ng	272112	4	11.30	1689370	20.0
Trifluoroacetic a...	13.78	6.3	ng	317569	5	13.94	1001630	20.0
Heptadecane	14.41	3.0	ng	151374	5	13.94	1001630	20.0
Tetradecane	14.69	2.1	ng	124886	6	15.36	1166850	20.0
Pentadecane, 8-he...	15.00	3.3	ng	194312	6	15.36	1166850	20.0
Tetracosane	15.76	4.9	ng	288265	6	15.36	1166850	20.0
Octadecane, 2-met...	16.21	6.1	ng	357169	6	15.36	1166850	20.0
Octadecane	16.75	4.9	ng	285734	6	15.36	1166850	20.0
Pentadecane	17.38	4.4	ng	255227	6	15.36	1166850	20.0