

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061016\
 Data File : BF087894.D
 Acq On : 11 Jun 2016 1:29
 Operator : UM/SJ
 Sample : H3440-01
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W-09

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-BF060716.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	1.747	12	14	23	rVB	225337	475795	9.95%	1.807%
2	1.873	23	25	72	rVB	2165724	4782616	100.00%	18.162%
3	4.970	294	296	299	rBV	67002	85004	1.78%	0.323%
4	5.404	332	334	336	rVB	1276099	1248033	26.10%	4.739%
5	6.445	422	425	427	rBV	930664	865740	18.10%	3.288%
6	6.536	429	433	435	rBV2	45493	114529	2.39%	0.435%
7	6.582	435	437	439	rVV	2206302	2273951	47.55%	8.635%
8	6.662	443	444	447	rVB	216044	187432	3.92%	0.712%
9	6.810	455	457	459	rBV	749038	634806	13.27%	2.411%
10	6.890	462	464	467	rVB2	42527	59862	1.25%	0.227%
11	6.970	469	471	473	rVB	2400199	2098758	43.88%	7.970%
12	7.382	504	507	509	rBV	1716854	1815618	37.96%	6.895%
13	7.462	512	514	516	rBV2	39326	48926	1.02%	0.186%
14	8.102	567	570	572	rBV	743941	712115	14.89%	2.704%
15	8.365	589	593	596	rBV2	28055	48962	1.02%	0.186%
16	8.559	607	610	613	rVB4	22665	58567	1.22%	0.222%
17	8.673	618	620	622	rBV2	38729	76909	1.61%	0.292%
18	8.765	626	628	631	rBV2	42020	58366	1.22%	0.222%
19	9.176	661	664	666	rBV	2961701	2657383	55.56%	10.091%
20	9.279	671	673	676	rVV2	41578	81085	1.70%	0.308%
21	9.565	696	698	704	rVB	126845	146438	3.06%	0.556%
22	9.851	720	723	725	rBV	848777	797379	16.67%	3.028%
23	9.988	734	735	741	rVB2	64360	69214	1.45%	0.263%
24	10.216	753	755	758	rBV2	31970	50975	1.07%	0.194%
25	10.639	789	792	794	rVB2	1063653	1424167	29.78%	5.408%
26	10.731	797	800	802	rBV2	47526	108703	2.27%	0.413%
27	10.776	802	804	806	rVV2	450952	395854	8.28%	1.503%
28	10.822	806	808	809	rVV	57649	60777	1.27%	0.231%
29	10.856	809	811	812	rVV2	77451	94421	1.97%	0.359%
30	10.879	812	813	817	rVV3	61285	120862	2.53%	0.459%
31	10.948	817	819	822	rVV2	45304	110623	2.31%	0.420%
32	10.994	822	823	825	rVV2	91486	107407	2.25%	0.408%
33	11.039	825	827	829	rVV2	54196	90122	1.88%	0.342%
34	11.074	829	830	832	rVB	60251	51533	1.08%	0.196%

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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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35	11.199	839	841	845	rVB2	48752	72380	1.51%	0.275%
36	11.325	850	852	855	rVB	767753	639298	13.37%	2.428%
37	12.914	988	991	993	rBV	2242778	2290693	47.90%	8.699%
38	13.474	1038	1040	1042	rVB	320287	270941	5.67%	1.029%
39	13.954	1079	1082	1084	rBV	634616	605035	12.65%	2.298%
40	15.360	1202	1205	1208	rBV	383122	442170	9.25%	1.679%

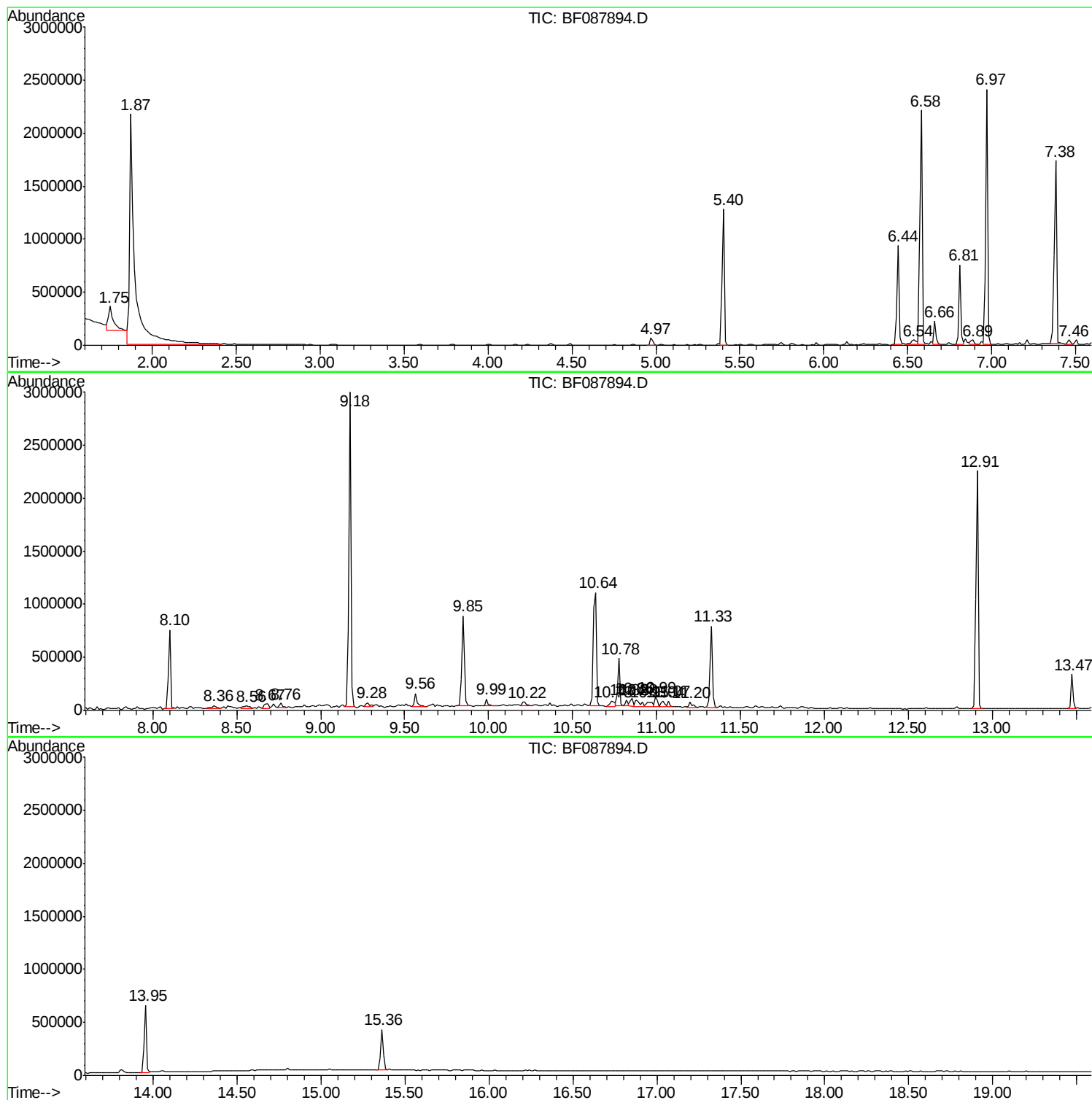
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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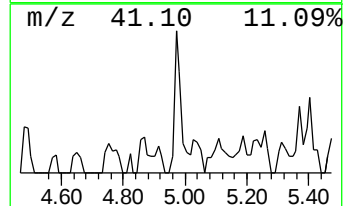
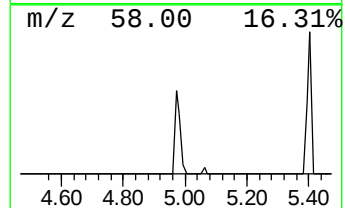
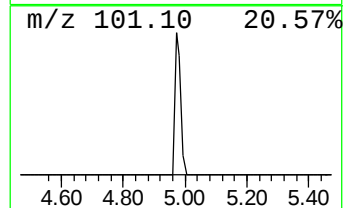
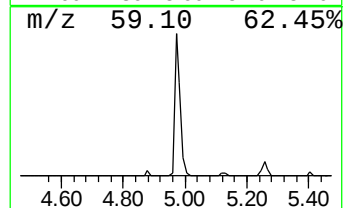
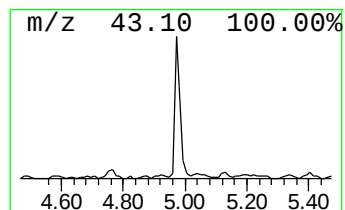
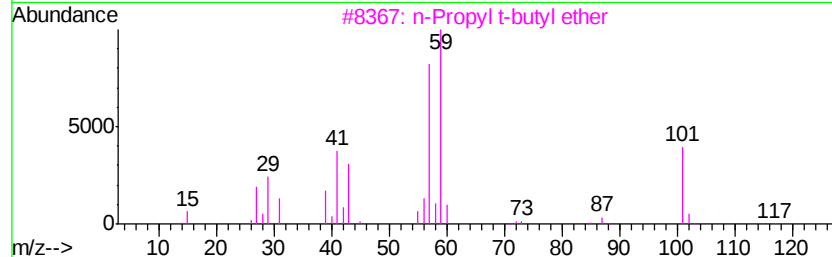
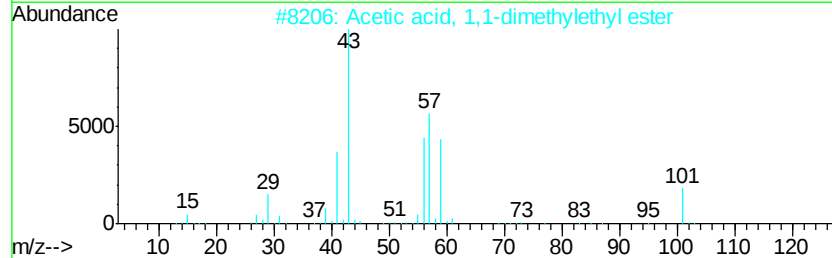
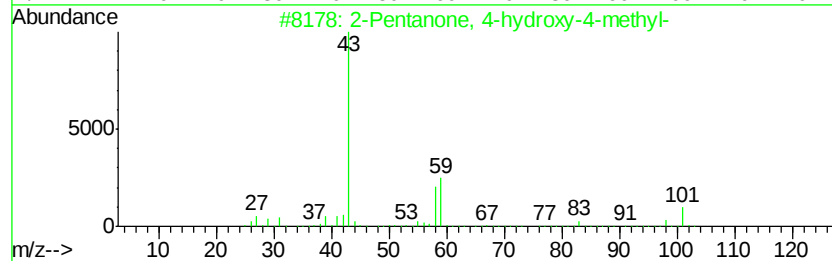
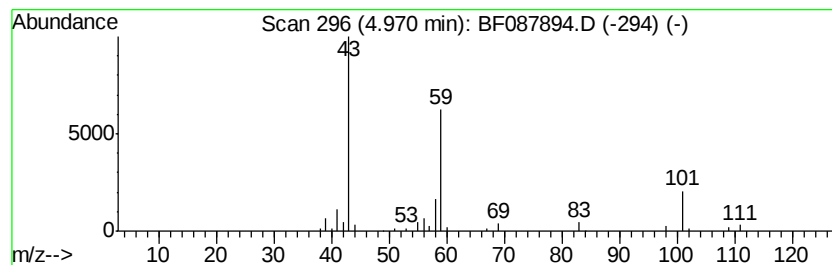
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.97	2.68 ng	85004	1,4-Dichlorobenzene-d4	6.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2		Acetic acid, 1,1-dimethylethyl ester	116	C6H12O2	000540-88-5	38
3		n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	28
4		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25



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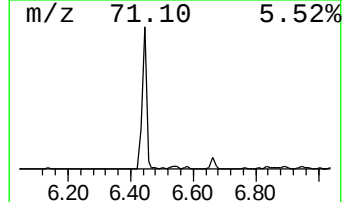
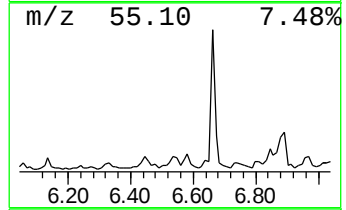
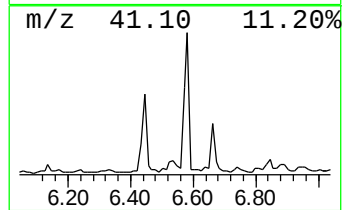
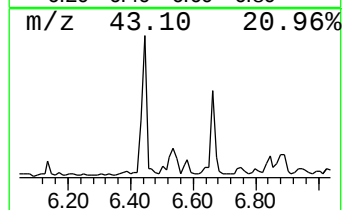
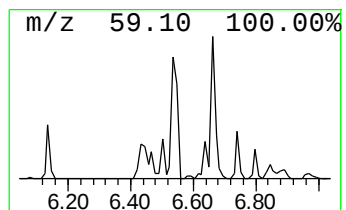
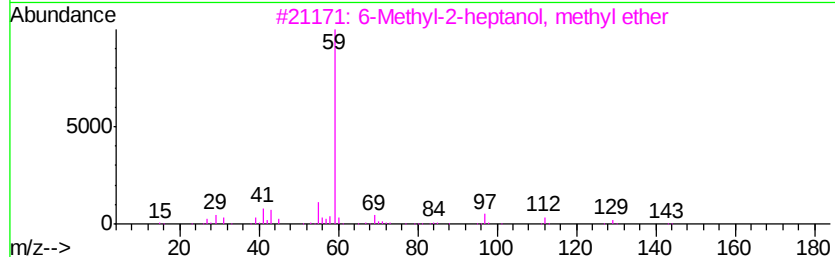
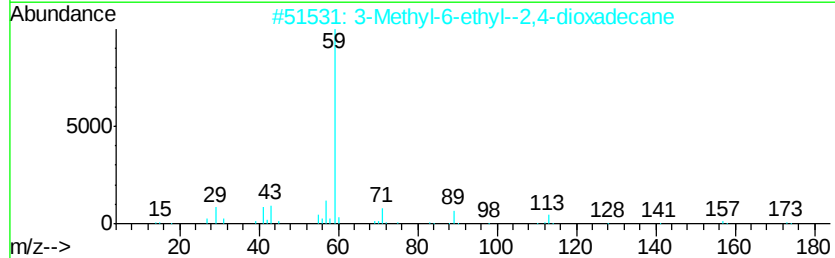
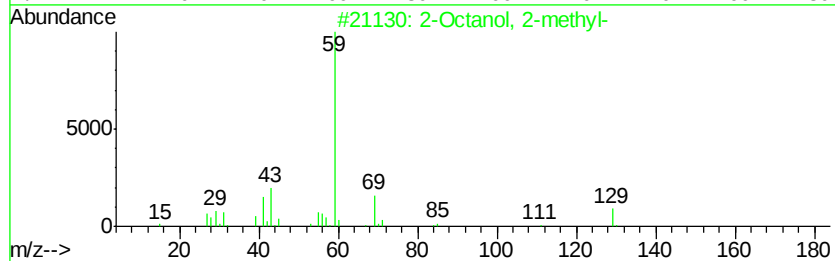
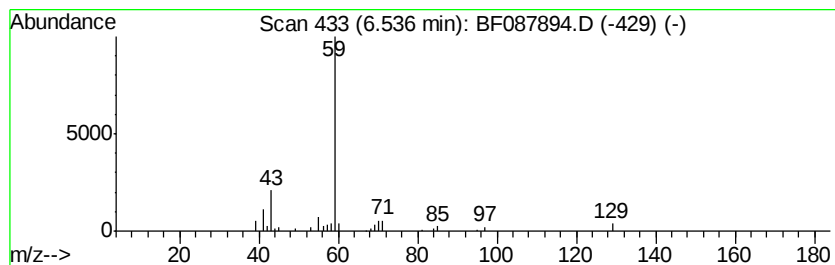
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TIC Library : C:\Database\NIST11.L
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Peak Number 4 2-Octanol, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.54	3.61 ng	114529	1,4-Dichlorobenzene-d4	6.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Octanol, 2-methyl-	144	C9H20O	000628-44-4	56
2			3-Methyl-6-ethyl--2,4-dioxadecane	188	C11H24O2	1000334-83-2	50
3			6-Methyl-2-heptanol, methyl ether	144	C9H20O	1000365-52-0	50
4			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	40
5			2,3,3-Trimethyl-2-pentanol	130	C8H18O	023171-85-9	40



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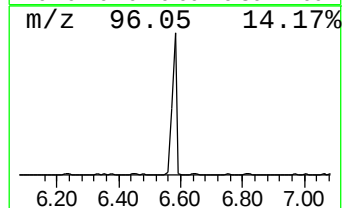
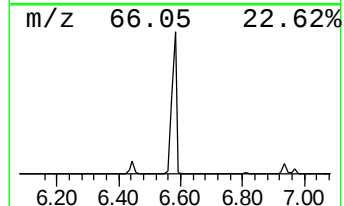
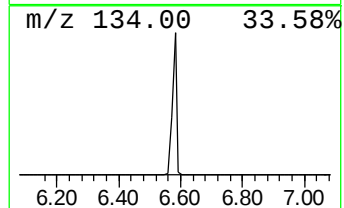
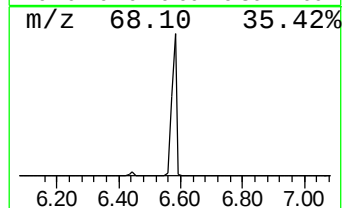
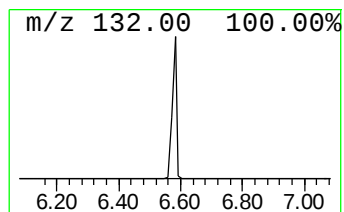
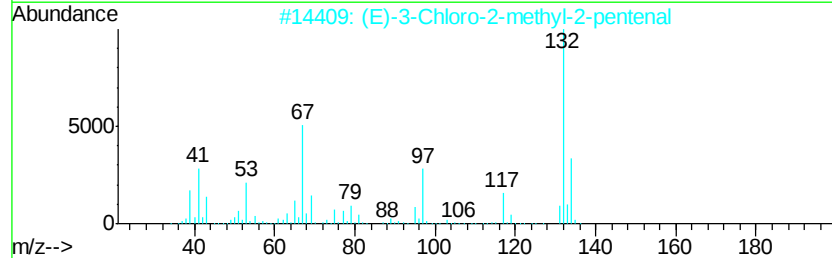
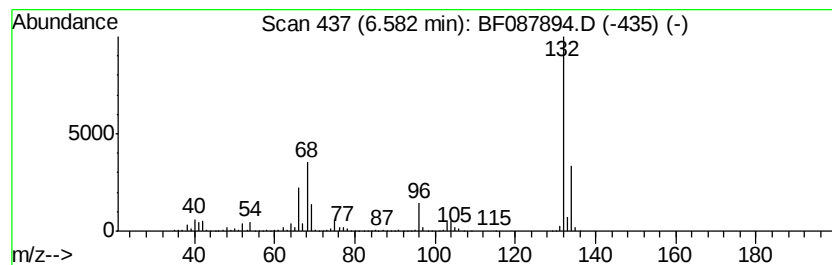
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown6.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	71.64 ng	2273950	1,4-Dichlorobenzene-d4	6.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	12
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		Xenon	132	Xe	007440-63-3	9



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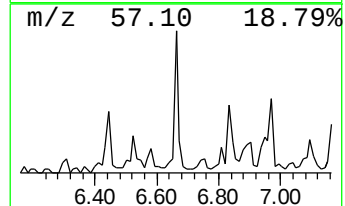
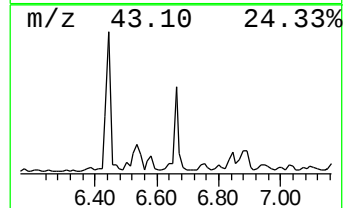
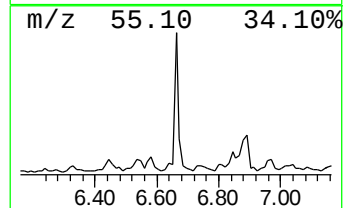
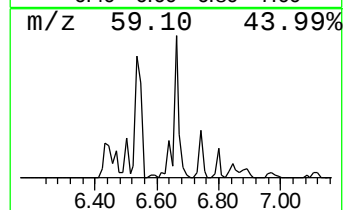
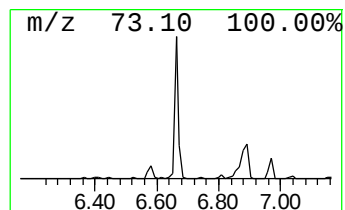
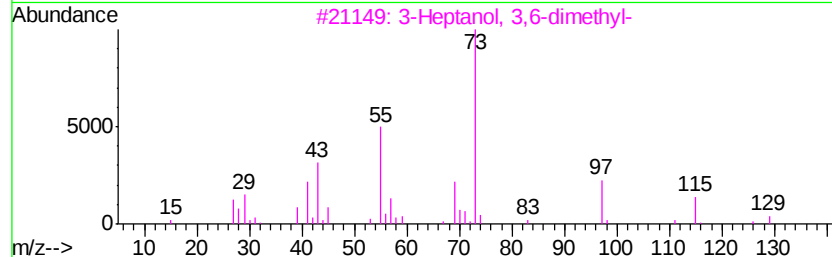
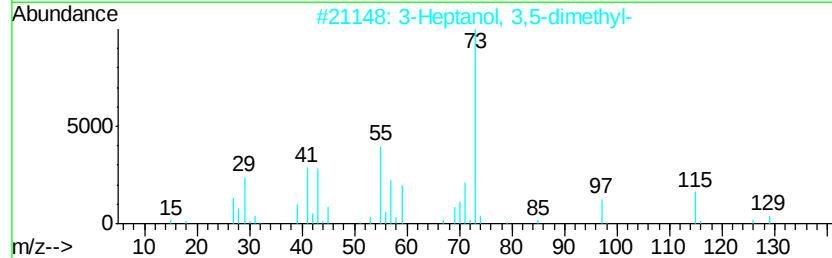
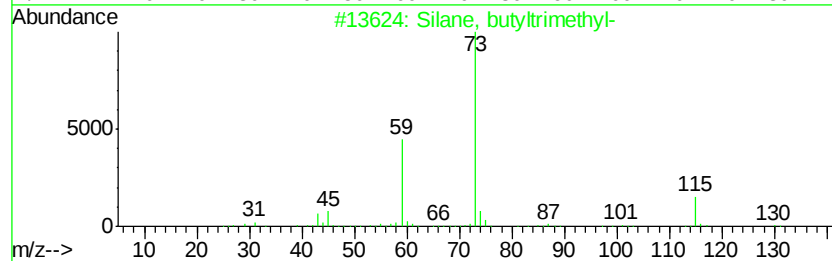
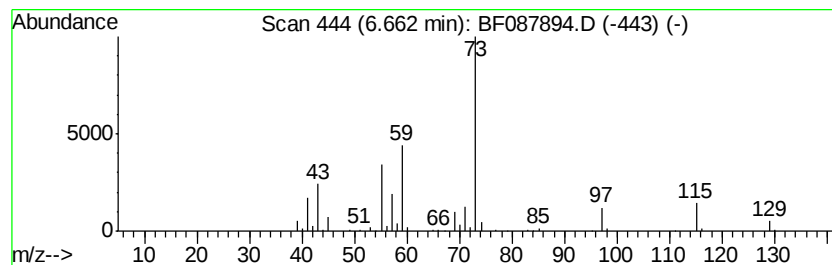
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown6.66 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	5.91 ng	187432	1,4-Dichlorobenzene-d4	6.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Silane, butyltrimethyl-	130	C7H18Si	001000-49-3	38
2			3-Heptanol, 3,5-dimethyl-	144	C9H20O	019549-74-7	37
3			3-Heptanol, 3,6-dimethyl-	144	C9H20O	001573-28-0	23
4			Silane, trimethylpropyl-	116	C6H16Si	003510-70-1	23
5			2,5-Dimethyl-4-hydroxy-3-hexanone	144	C8H16O2	000815-77-0	14



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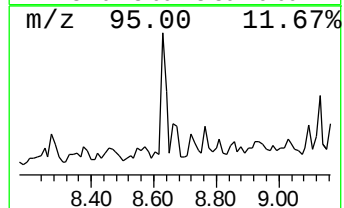
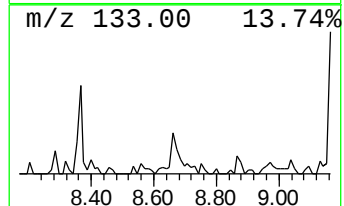
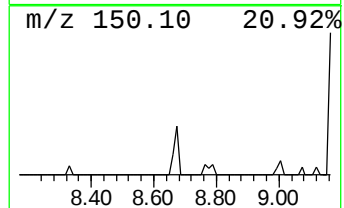
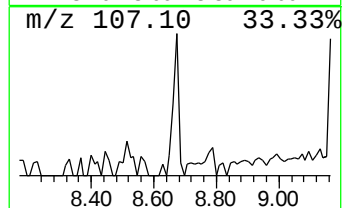
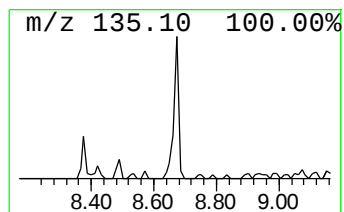
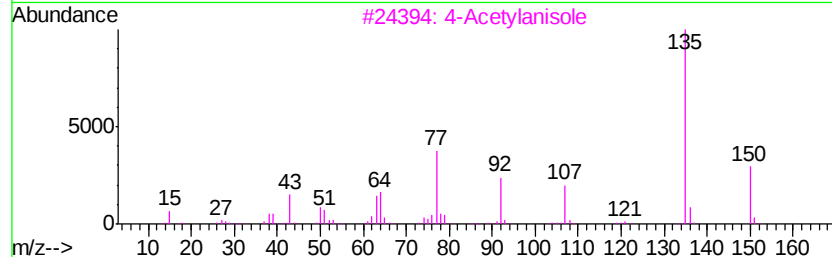
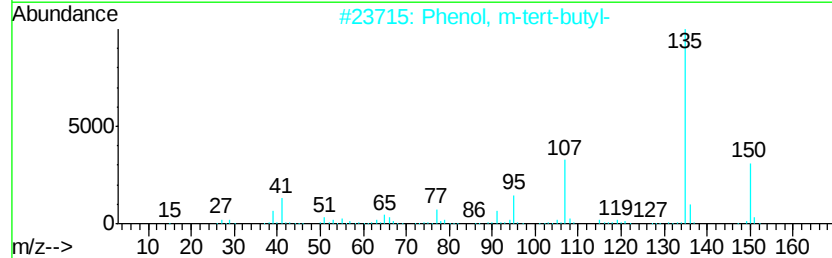
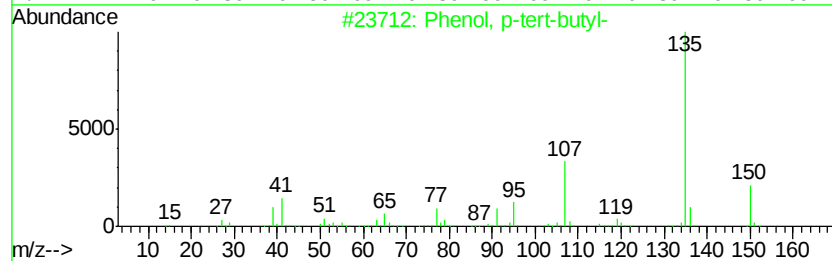
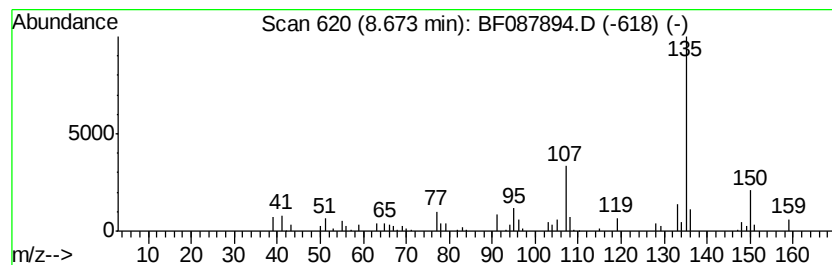
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Peak Number 8 Phenol, p-tert-butyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.67	2.16 ng	76909	Napthalene-d8	8.10
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Phenol, p-tert-butyl-	150 C10H14O	000098-54-4	81
2	Phenol, m-tert-butyl-	150 C10H14O	000585-34-2	74
3	4-Acetylanisole	150 C9H10O2	000100-06-1	58
4	Benzene, 1-methoxy-4-(1-methylet...	150 C10H14O	004132-48-3	53
5	Thymol	150 C10H14O	000089-83-8	53



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ALS Vial : 25 Sample Multiplier: 1

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ClientSampleId :
W-09

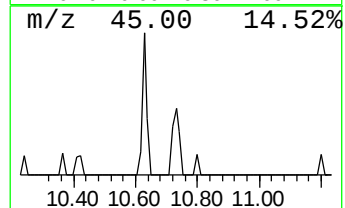
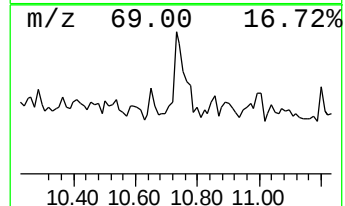
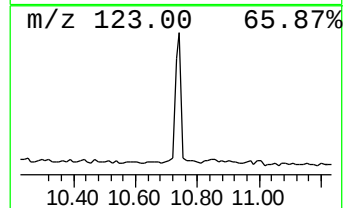
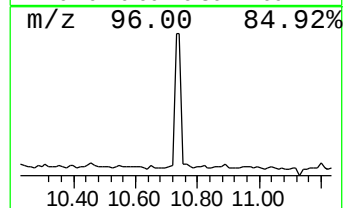
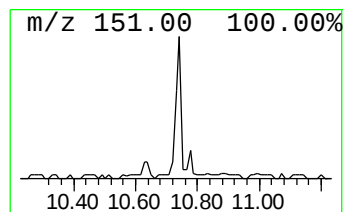
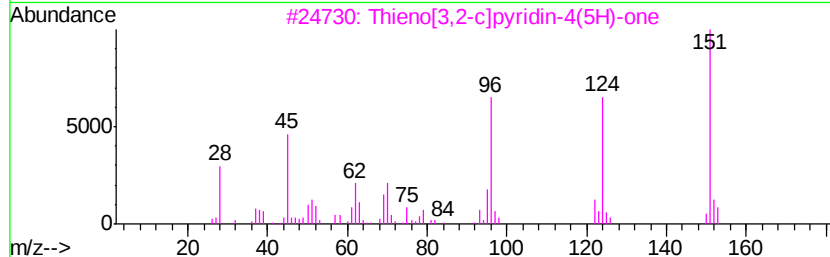
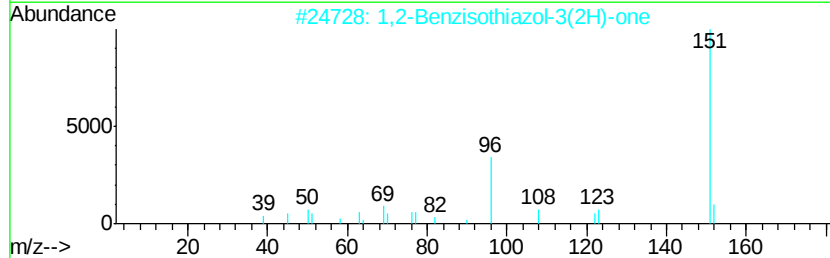
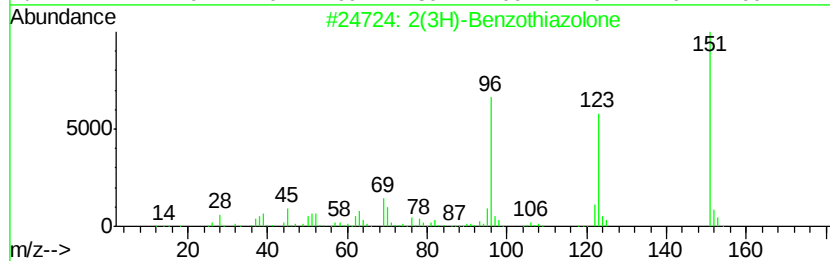
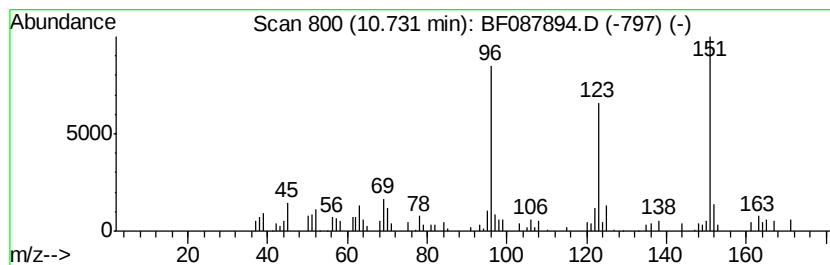
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 2(3H)-Benzothiazolone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.73	3.40 ng	108703	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	90
2		1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	76
3		Thieno[3,2-c]pyridin-4(5H)-one	151	C7H5NOS	027685-92-3	59
4		5-Ethoxycarbonylpiperidine-2-carbo...	195	C9H9NO4	017880-34-1	38
5		4-Methoxyformanilide	151	C8H9NO2	005470-34-8	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061016\
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Acq On : 11 Jun 2016 1:29
Operator : UM/SJ
Sample : H3440-01
Misc :
ALS Vial : 25 Sample Multiplier: 1

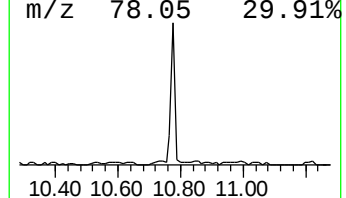
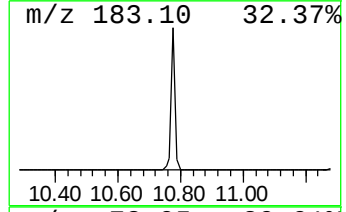
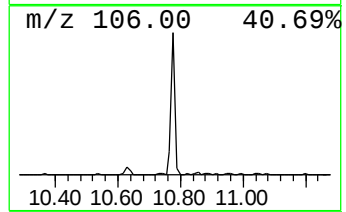
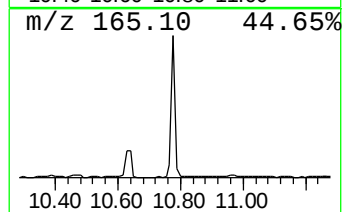
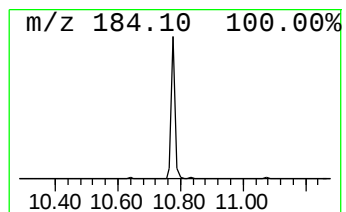
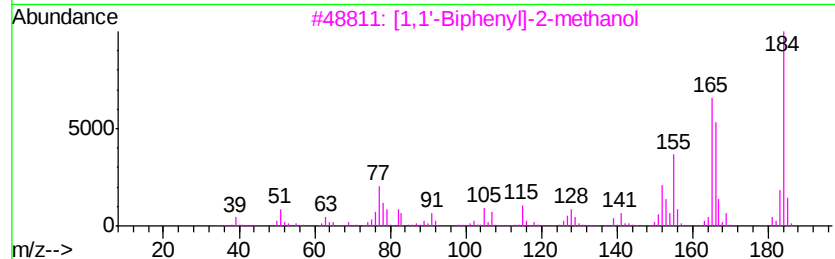
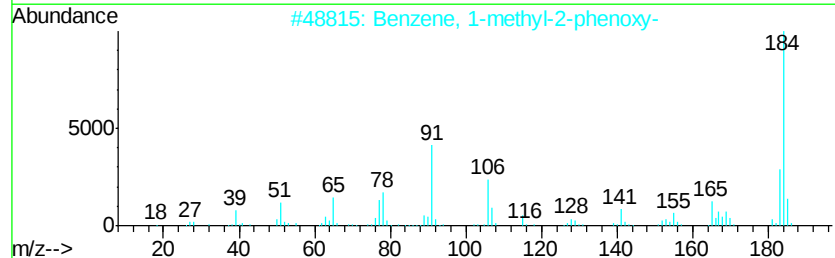
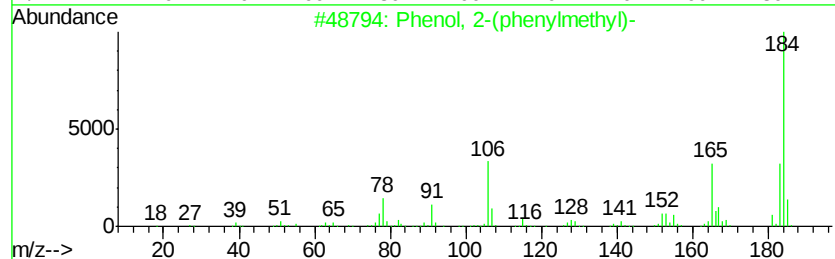
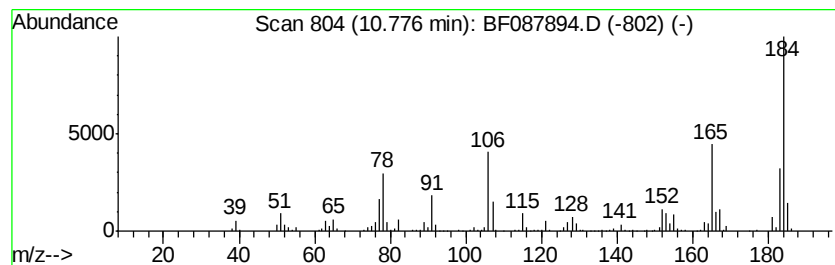
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Phenol, 2-(phenylmethyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.78	12.38 ng	395854	Phenanthrene-d10	11.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(phenylmethyl)-	184	C13H12O	028994-41-4	95
2	Benzene, 1-methyl-2-phenoxy-	184	C13H12O	003991-61-5	74
3	[1,1'-Biphenyl]-2-methanol	184	C13H12O	002928-43-0	70
4	p-Benzylphenol	184	C13H12O	000101-53-1	58
5	[1,1'-Biphenyl]-4-methanol	184	C13H12O	003597-91-9	58



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Data File : BF087894.D
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Instrument :
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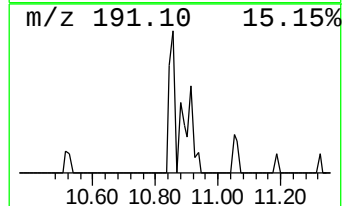
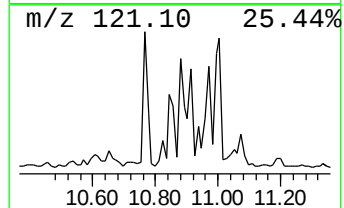
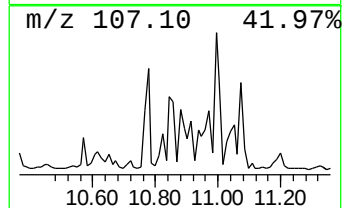
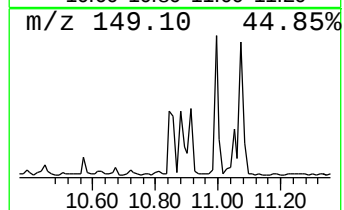
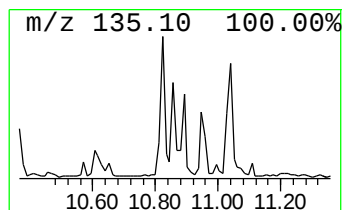
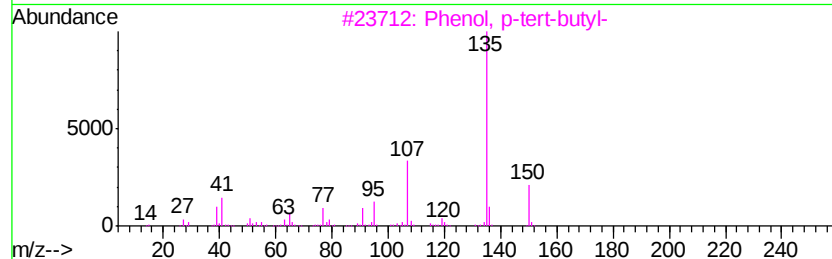
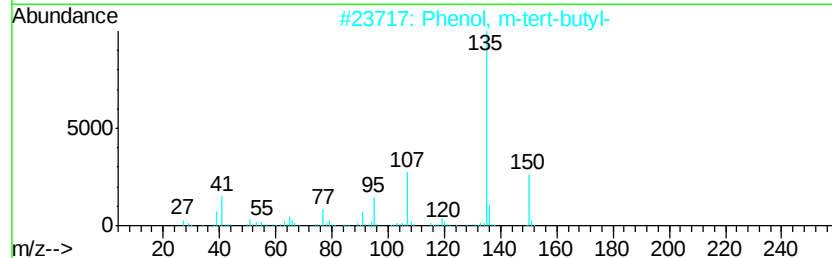
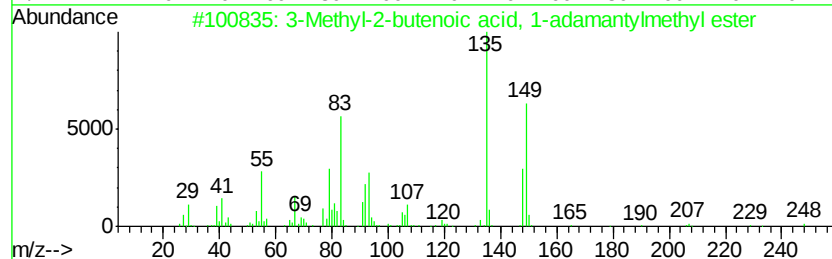
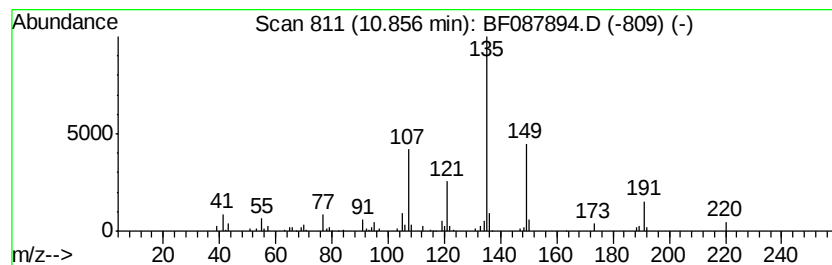
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 3-Methyl-2-butenic acid, 1... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.86	2.95 ng	94421	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Methyl-2-butenic acid, 1-adam...	248	C16H24O2	1000282-89-4	53
2		Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	52
3		Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	50
4		Hexestrol, 0-heptafluorobutyl-	466	C22H21F7O3	1000365-31-9	50
5		Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061016\
Data File : BF087894.D
Acq On : 11 Jun 2016 1:29
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Sample : H3440-01
Misc :
ALS Vial : 25 Sample Multiplier: 1

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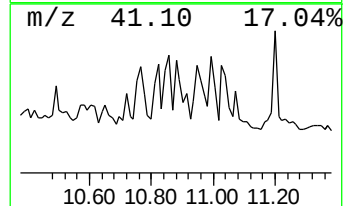
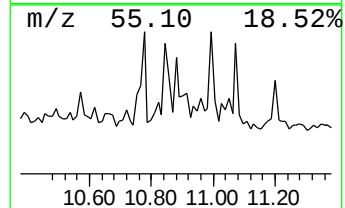
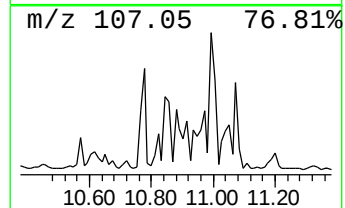
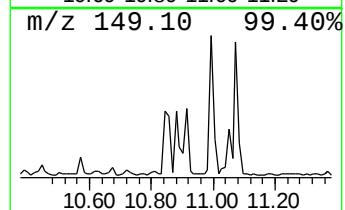
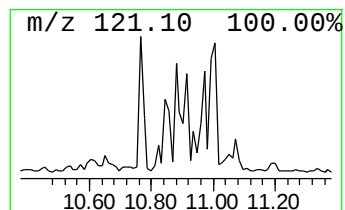
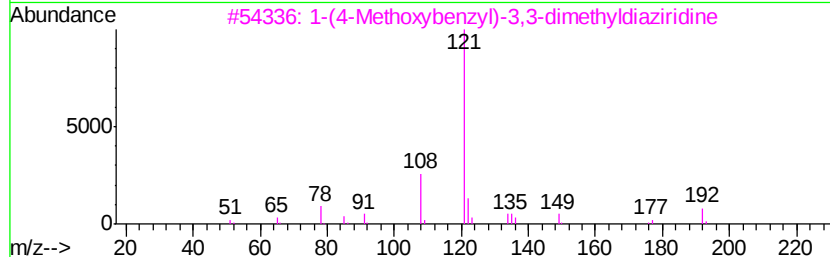
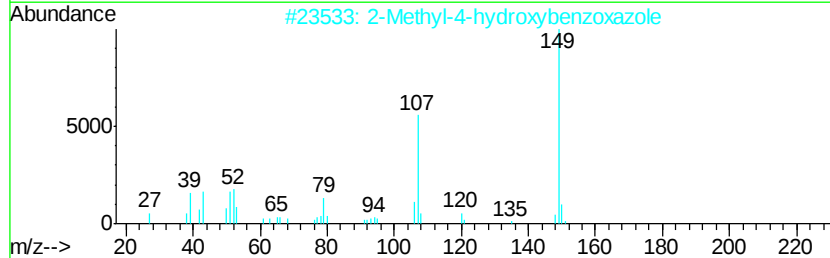
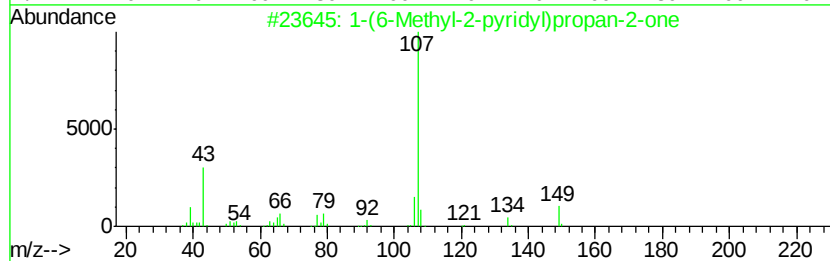
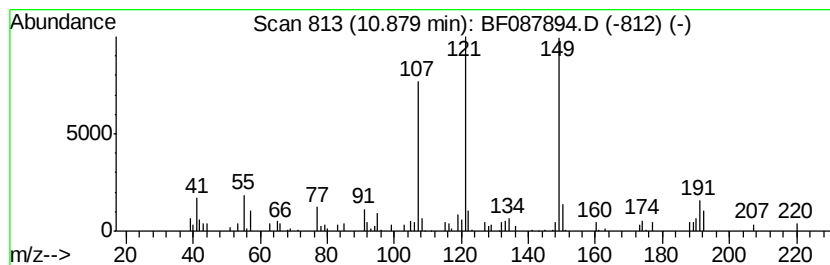
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown10.88 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.88	3.78 ng	120862	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-(6-Methyl-2-pyridyl)propan-2-one	149	C9H11NO	065702-08-1	46
2		2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	43
3		1-(4-Methoxybenzyl)-3,3-dimethyl...	192	C11H16N2O	022184-66-3	25
4		Isoxazole, 4,5-dihydro-5-[(4-met...	191	C11H13NO2	1000362-14-0	22
5		Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	22



Data Path : Z:\HPCHEM1\BNA F\DATA\BF061016\
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Misc :
ALS Vial : 25 Sample Multiplier: 1

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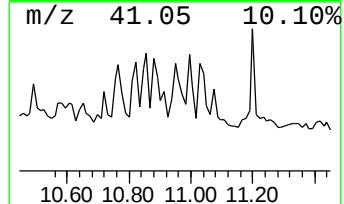
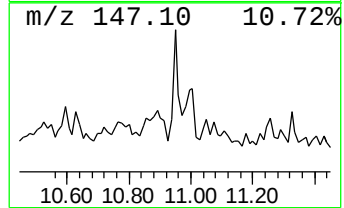
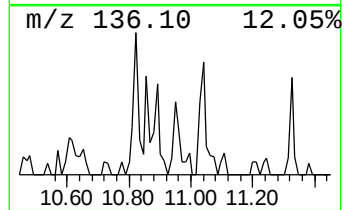
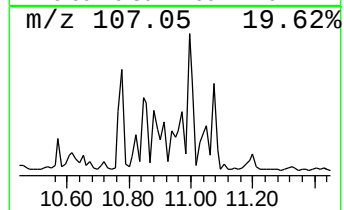
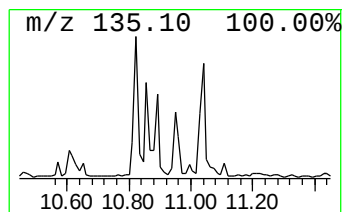
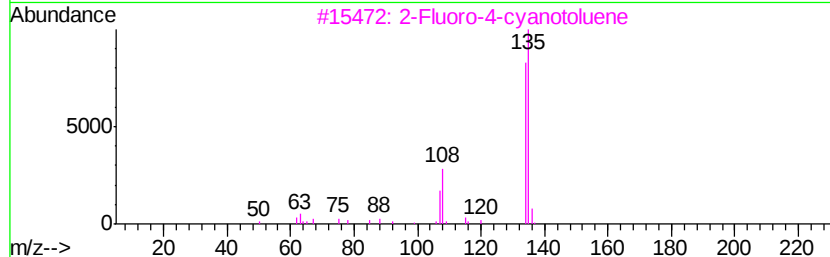
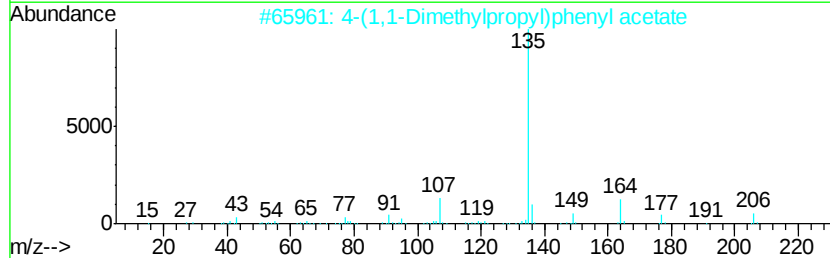
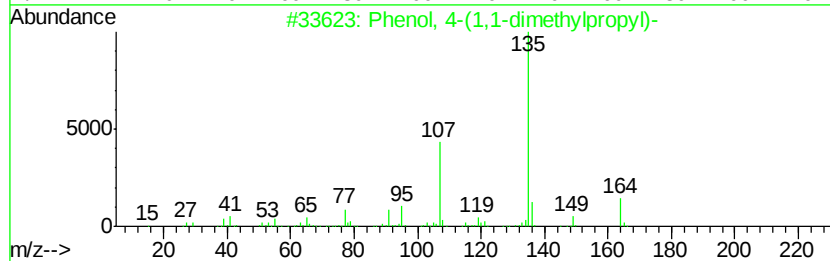
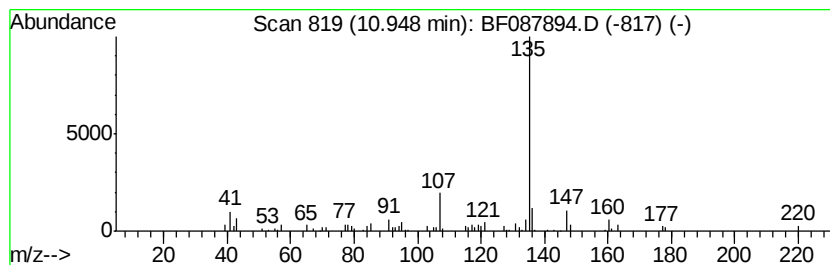
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Phenol, 4-(1,1-dimethylprop... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.95	3.46 ng	110623	Phenanthrene-d10	11.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	64
2			4-(1,1-Dimethylpropyl)phenyl ace...	206	C13H18O2	1000373-45-3	64
3			2-Fluoro-4-cyanotoluene	135	C8H6FN	170572-49-3	59
4			Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	59
5			1H-Pyrazolo[3,4-d]pyrimidin-4-amine	135	C5H5N5	002380-63-4	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061016\
Data File : BF087894.D
Acq On : 11 Jun 2016 1:29
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Misc :
ALS Vial : 25 Sample Multiplier: 1

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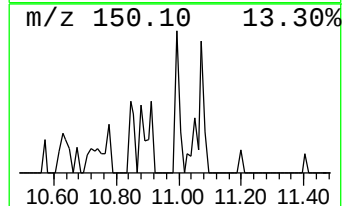
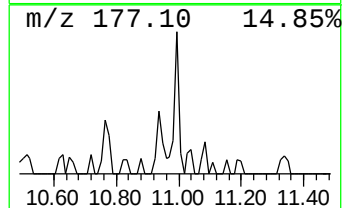
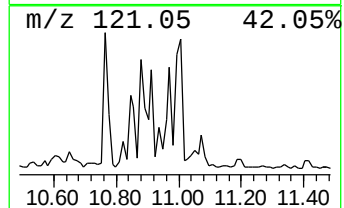
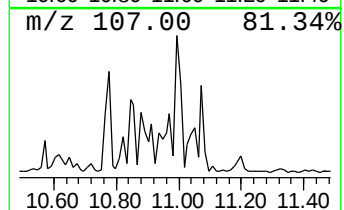
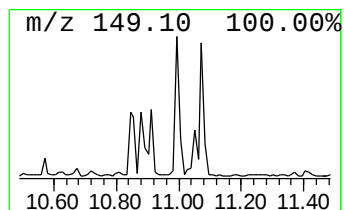
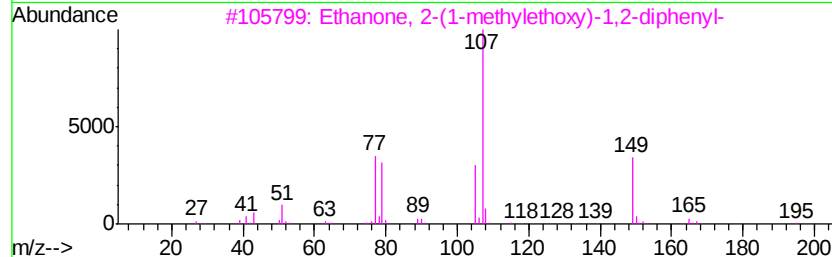
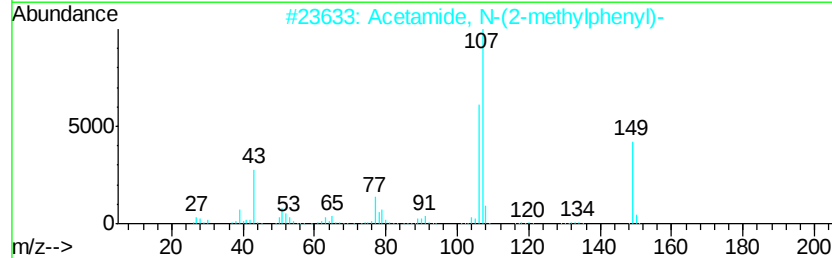
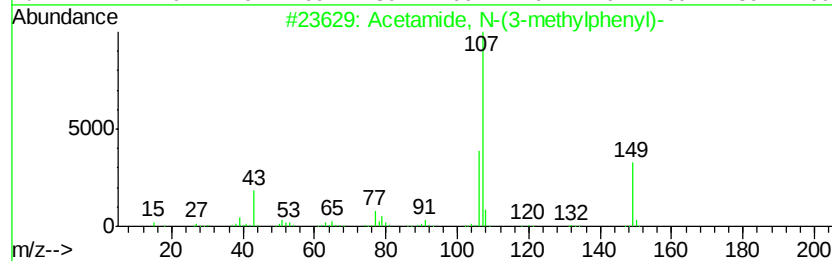
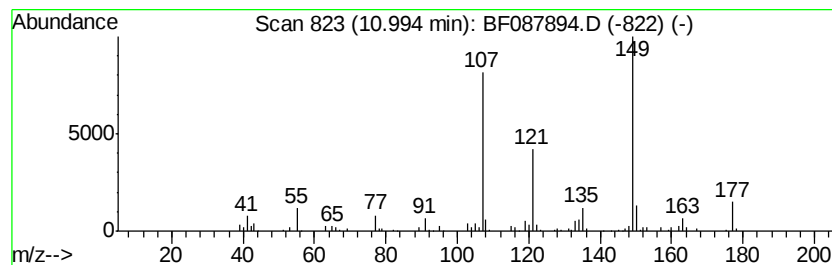
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Acetamide, N-(3-methylphenyl)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.99	3.36 ng	107407	Phenanthrene-d10	11.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	64
2			Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	59
3			Ethanone, 2-(1-methylethoxy)-1,2-diphenyl-	254	C17H18O2	006652-28-4	45
4			Benzene, [1-(1-methylethoxy)-3-methylphenyl]-	190	C13H18O	098088-47-2	45
5			Benzenemethanol, 4-(1,1-dimethylethoxy)-	164	C11H16O	000877-65-6	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061016\
Data File : BF087894.D
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Misc :
ALS Vial : 25 Sample Multiplier: 1

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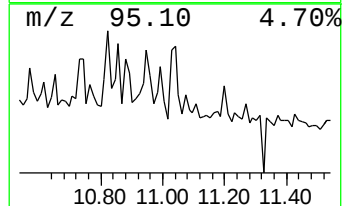
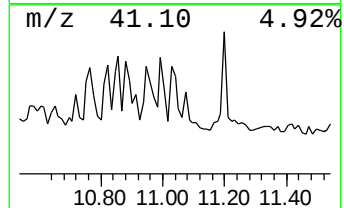
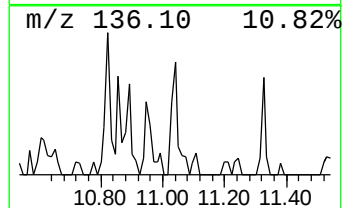
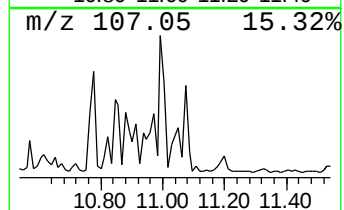
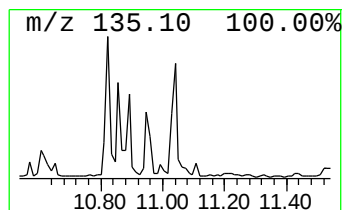
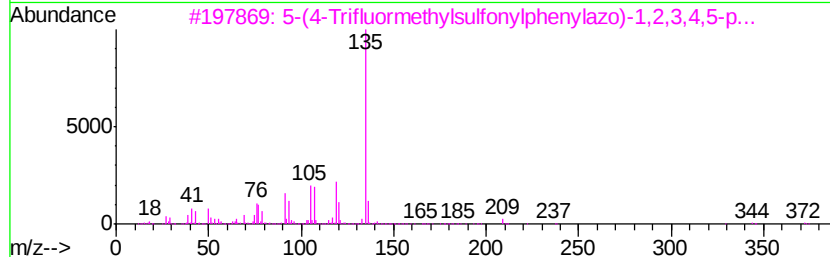
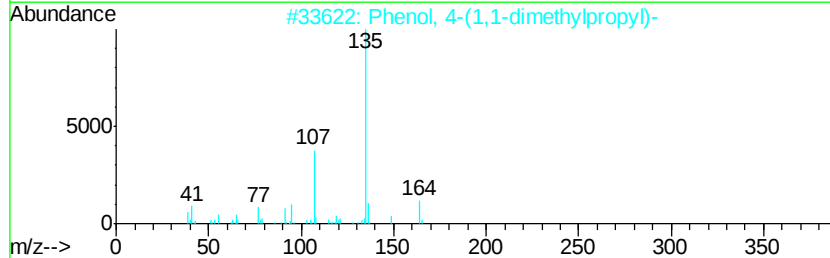
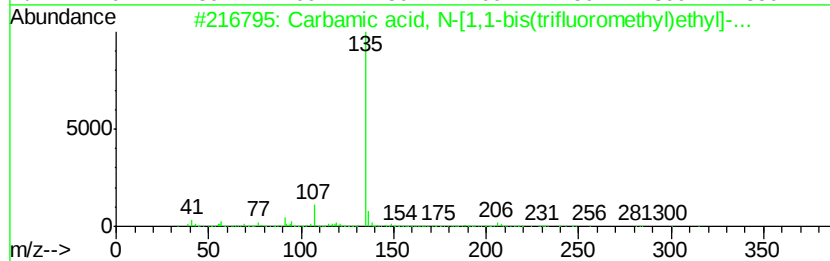
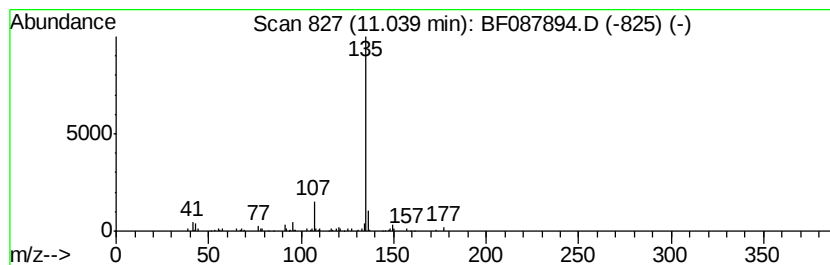
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Carbamic acid, N-[1,1-bis(t... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.04	2.82 ng	90122	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Carbamic acid, N-[1,1-bis(triflu...	413	C19H25F6N02	296242-69-8	72
2		Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	72
3		5-(4-Trifluormethylsulfonylphenyl)-	372	C17H19F3N2O2S	187338-96-1	64
4		2-Fluoro-4-cyanotoluene	135	C8H6FN	170572-49-3	64
5		Benzenepropanol, 4-methyl-	150	C10H14O	005406-39-3	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061016\
Data File : BF087894.D
Acq On : 11 Jun 2016 1:29
Operator : UM/SJ
Sample : H3440-01
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
W-09

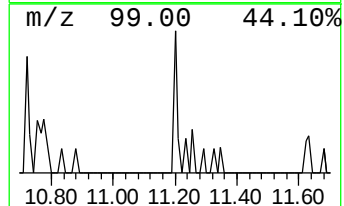
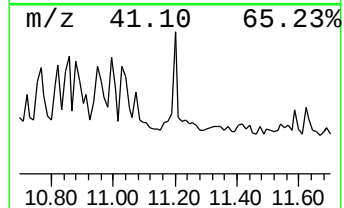
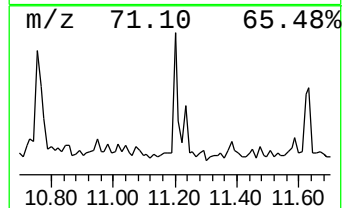
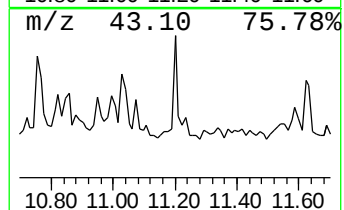
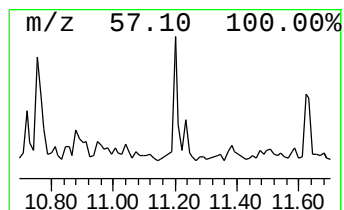
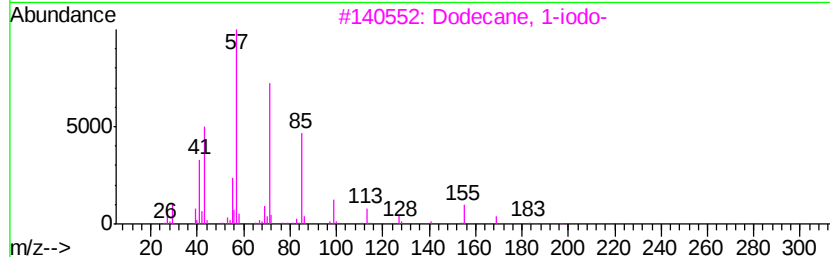
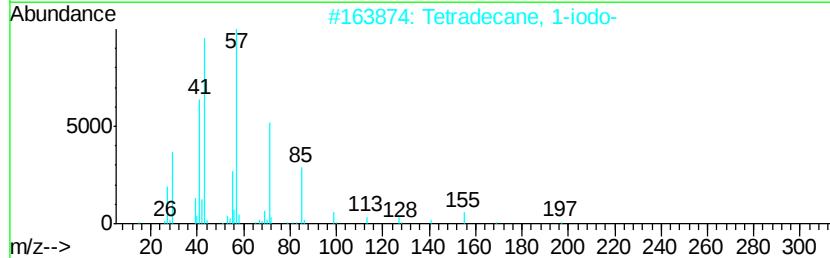
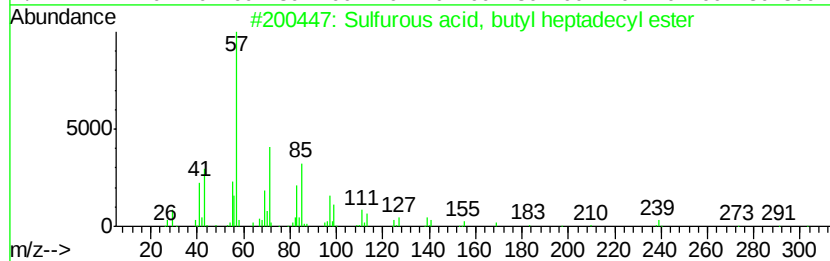
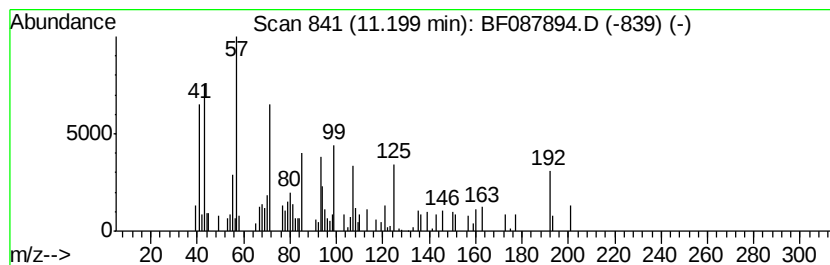
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 unknown11.20 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.20	2.26 ng	72380	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, butyl heptadecyl...	376	C21H44O3S	1000309-18-4	27
2		Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	22
3		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	22
4		Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	22
5		Tridecane, 2-methyl-	198	C14H30	001560-96-9	22



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Acq On : 11 Jun 2016 1:29
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Sample : H3440-01
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
W-09

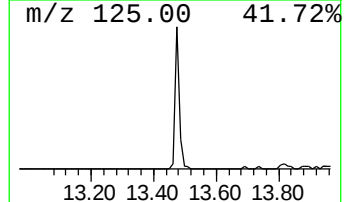
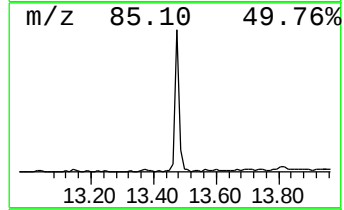
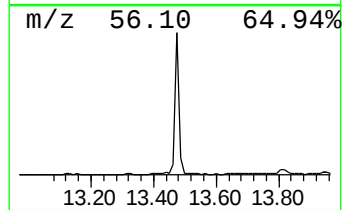
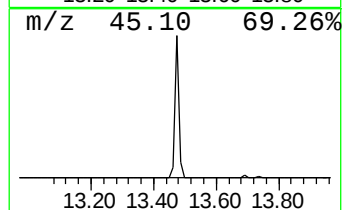
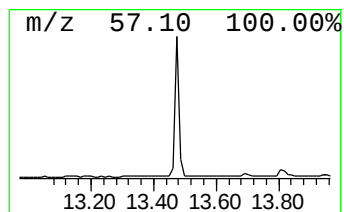
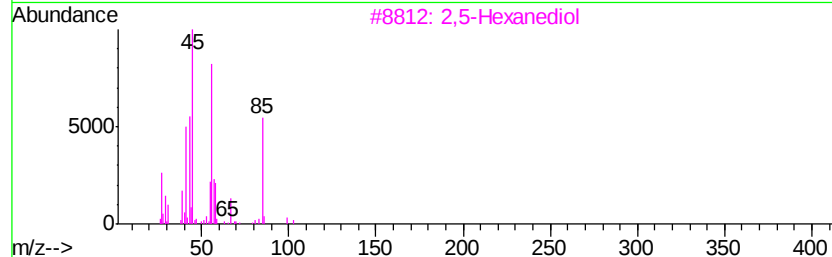
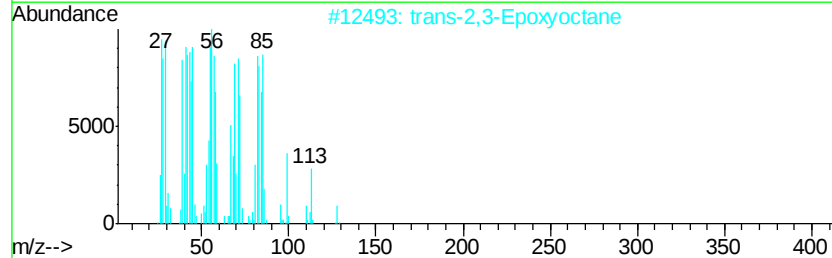
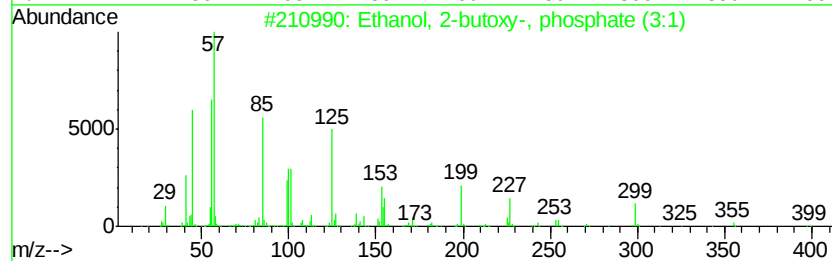
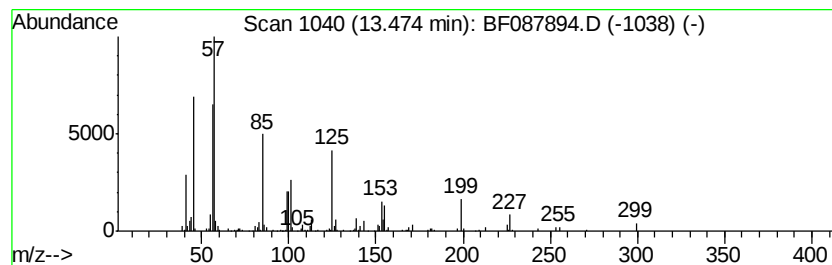
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 Ethanol, 2-butoxy-, phospho... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	8.96 ng	270941	Chrysene-d12	13.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-butoxy-, phosphate (3:1)	398	C18H39O7P	000078-51-3	52
2		trans-2,3-Epoxyoctane	128	C8H16O	028180-70-3	43
3		2,5-Hexanediol	118	C6H14O2	002935-44-6	35
4		Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	14
5		1-Methanesulfonylpiperazine	164	C5H12N2O2S	1000303-61-8	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061016\
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 Sample : H3440-01
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W-09

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.97	2.7 ng		85004	1	6.81	634806	20.0
2-Octanol, 2-methyl-	6.54	3.6 ng		114529	1	6.81	634806	20.0
unknown6.58	6.58	71.6 ng		2273950	1	6.81	634806	20.0
unknown6.66	6.66	5.9 ng		187432	1	6.81	634806	20.0
Phenol, p-tert-bu...	8.67	2.2 ng		76909	2	8.10	712115	20.0
2(3H)-Benzothiazo...	10.73	3.4 ng		108703	4	11.32	639298	20.0
Phenol, 2-(phenyl...	10.78	12.4 ng		395854	4	11.32	639298	20.0
3-Methyl-2-butenol...	10.86	3.0 ng		94421	4	11.32	639298	20.0
unknown10.88	10.88	3.8 ng		120862	4	11.32	639298	20.0
Phenol, 4-(1,1-di...	10.95	3.5 ng		110623	4	11.32	639298	20.0
Acetamide, N-(3-m...	10.99	3.4 ng		107407	4	11.32	639298	20.0
Carbamic acid, N-...	11.04	2.8 ng		90122	4	11.32	639298	20.0
unknown11.20	11.20	2.3 ng		72380	4	11.32	639298	20.0
Ethanol, 2-butoxy...	13.47	9.0 ng		270941	5	13.95	605035	20.0