

Data Path : Z:\HPCHEM1\BNA F\Data\BF110815\
 Data File : BF082794.D
 Acq On : 7 Nov 2015 7:10
 Operator : UM/IZ
 Sample : G4259-02
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-01-110315

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-BF110715.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	5.024	254	257	261	rVB	1284853	1569782	16.80%	2.213%
2	5.447	291	294	296	rBV	4155876	3782891	40.49%	5.334%
3	6.476	381	384	386	rBV	3207366	2791750	29.88%	3.936%
4	6.613	393	396	399	rBV	5978980	6069967	64.97%	8.559%
5	6.807	408	413	414	rBV2	78202	109657	1.17%	0.155%
6	6.842	414	416	418	rVV	1319357	1573993	16.85%	2.219%
7	7.002	428	430	433	rBV	4784463	5872959	62.86%	8.281%
8	7.413	462	466	468	rBV	5189866	5759813	61.65%	8.121%
9	7.505	471	474	475	rVB	134224	153871	1.65%	0.217%
10	7.550	475	478	481	rBV4	75545	113891	1.22%	0.161%
11	7.630	481	485	487	rBV2	109618	165573	1.77%	0.233%
12	7.790	496	499	501	rBV2	79301	128746	1.38%	0.182%
13	7.870	504	506	508	rBV2	218829	236846	2.53%	0.334%
14	8.053	519	522	523	rBV2	51120	99997	1.07%	0.141%
15	8.133	526	529	530	rVV	2624874	2330388	24.94%	3.286%
16	8.179	532	533	535	rVV	140939	197340	2.11%	0.278%
17	8.270	538	541	542	rVV2	91878	116093	1.24%	0.164%
18	8.328	544	546	548	rVV2	78995	115345	1.23%	0.163%
19	8.385	548	551	552	rVB	244834	237712	2.54%	0.335%
20	8.430	552	555	557	rBV3	45563	114000	1.22%	0.161%
21	8.522	561	563	565	rVB	132770	139030	1.49%	0.196%
22	8.613	566	571	572	rBV3	65203	177529	1.90%	0.250%
23	8.762	580	584	585	rBV2	75481	133863	1.43%	0.189%
24	8.796	585	587	589	rVB	292065	249295	2.67%	0.352%
25	8.979	598	603	606	rVB4	119482	308993	3.31%	0.436%
26	9.082	606	612	613	rBV4	90996	307857	3.29%	0.434%
27	9.219	621	624	625	rBV	8058783	9343181	100.00%	13.174%
28	9.368	635	637	638	rVB	112083	101349	1.08%	0.143%
29	9.608	654	658	659	rBV2	238633	345919	3.70%	0.488%
30	9.882	680	682	685	rVB	1771079	2350073	25.15%	3.314%
31	10.111	698	702	703	rBV3	80070	131109	1.40%	0.185%
32	10.191	707	709	711	rVV3	127020	182632	1.95%	0.258%
33	10.271	714	716	718	rBV	52382	102647	1.10%	0.145%
34	10.442	729	731	732	rBV2	48319	94310	1.01%	0.133%

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MW-01-110315

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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	10.556	739	741	744	rBV3	57863	131270	1.40%	0.185%
36	10.614	744	746	748	rBV	286622	349060	3.74%	0.492%
37	10.682	748	752	754	rBV	3831487	5754633	61.59%	8.114%
38	10.899	769	771	774	rVB2	118343	154269	1.65%	0.218%
39	11.242	799	801	806	rVB2	93860	197378	2.11%	0.278%
40	11.368	810	812	814	rVV	2781841	2304197	24.66%	3.249%
41	11.916	857	860	863	rBV2	435820	686291	7.35%	0.968%
42	12.077	871	874	878	rBV2	107772	180601	1.93%	0.255%
43	12.614	918	921	926	rBV3	65735	172091	1.84%	0.243%
44	12.694	926	928	934	rBV	618486	702789	7.52%	0.991%
45	12.785	934	936	939	rVB2	719782	872170	9.33%	1.230%
46	12.842	939	941	942	rBV	148420	152239	1.63%	0.215%
47	12.957	948	951	954	rBV	6054258	7681173	82.21%	10.830%
48	13.494	996	998	1001	rBV	759499	685264	7.33%	0.966%
49	13.540	1001	1002	1007	rVB2	87955	155499	1.66%	0.219%
50	13.871	1029	1031	1035	rBV	371828	429007	4.59%	0.605%
51	14.008	1040	1043	1045	rBV	1608863	2248886	24.07%	3.171%
52	14.191	1056	1059	1061	rBV	121417	145078	1.55%	0.205%
53	14.488	1084	1085	1090	rVB	101059	138959	1.49%	0.196%
54	14.797	1109	1112	1114	rVB	95529	133578	1.43%	0.188%
55	15.117	1138	1140	1142	rVB	100285	111869	1.20%	0.158%
56	15.425	1164	1167	1170	rBV	1325818	1687568	18.06%	2.379%
57	15.483	1170	1172	1175	rVB	86533	117039	1.25%	0.165%
58	16.386	1248	1251	1257	rVB2	48594	113476	1.21%	0.160%
59	17.403	1337	1340	1346	rVB4	40359	109128	1.17%	0.154%

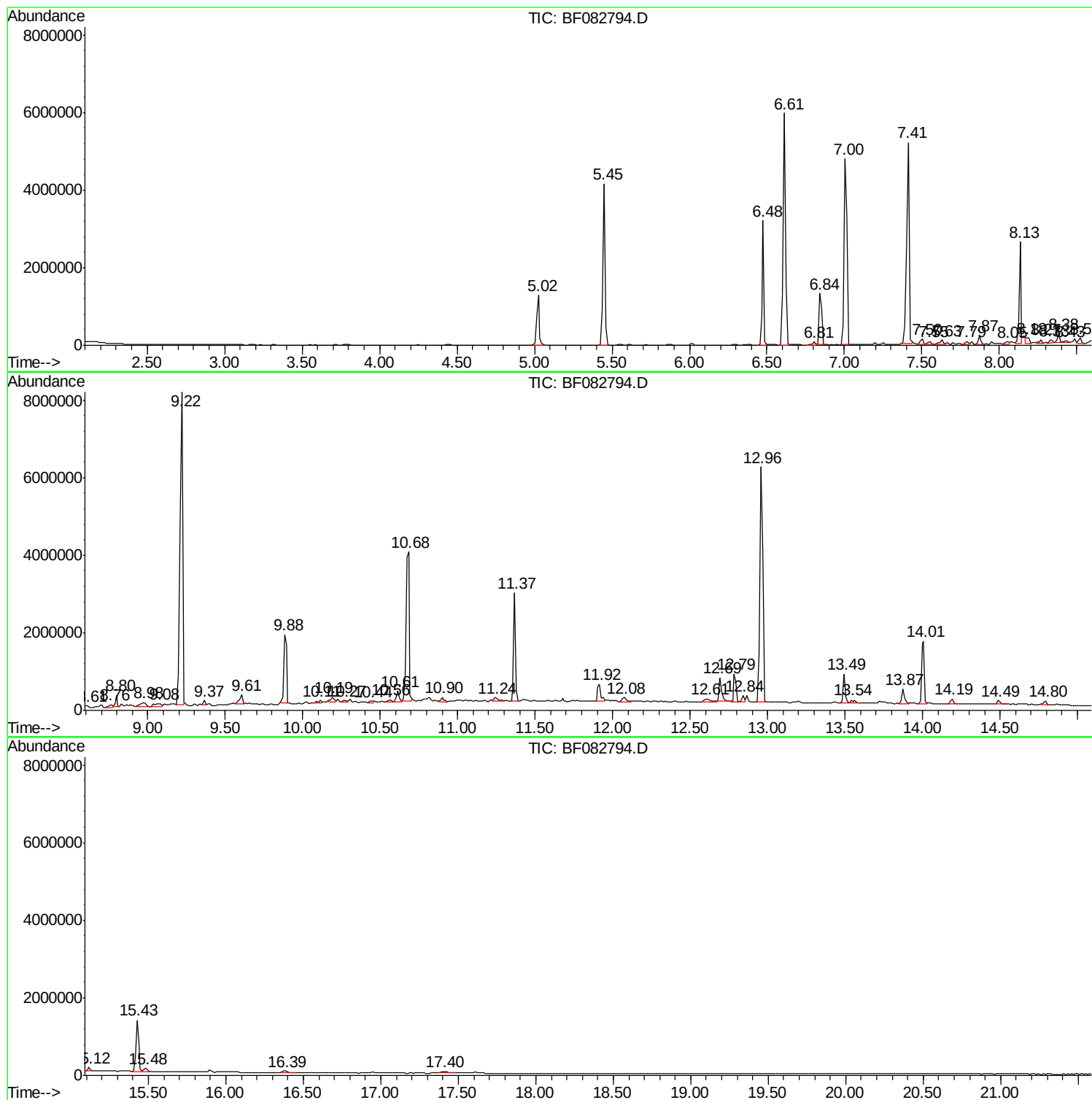
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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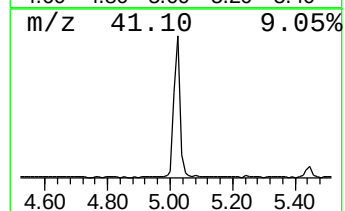
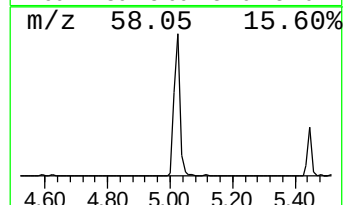
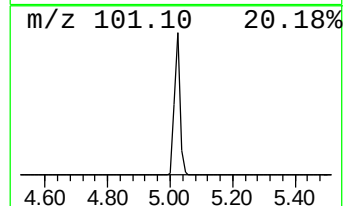
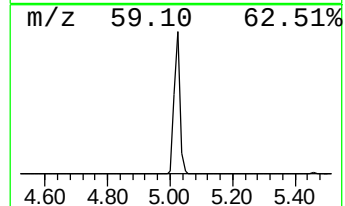
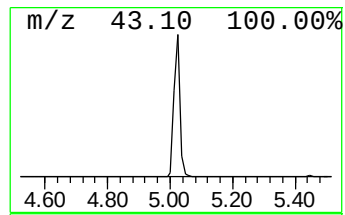
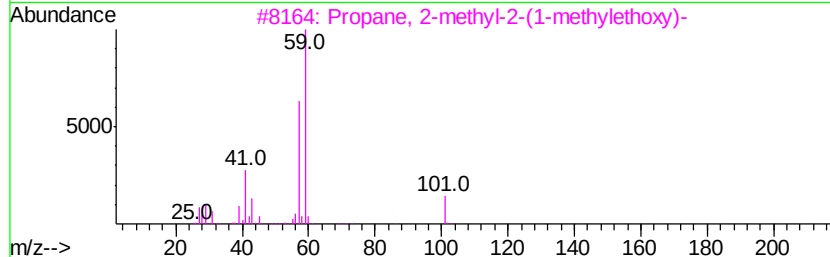
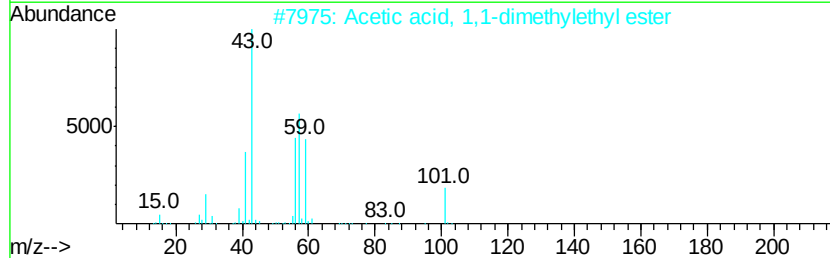
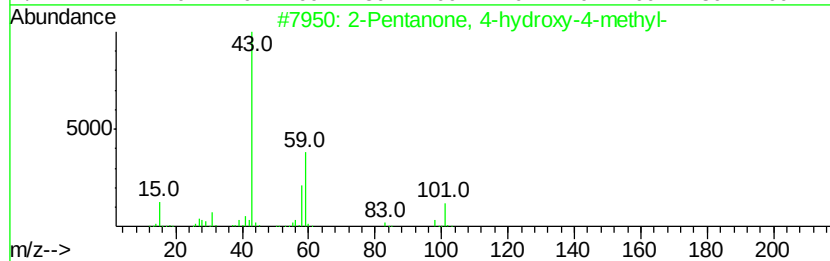
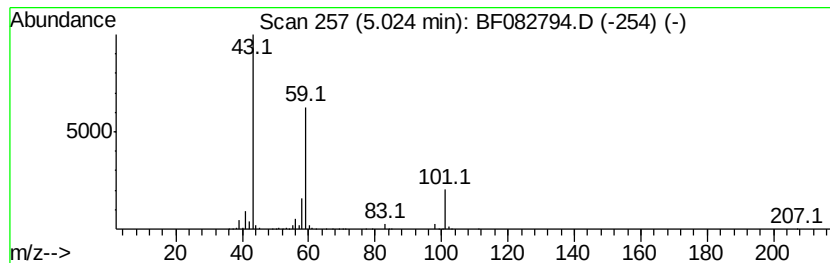
Instrument :
BNA_F
ClientSampleId :
MW-01-110315

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.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.
5.02	19.95 ng	1569780	1,4-Dichlorobenzene-d4	6.84
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	Propane, 2-methyl-2-(1-methyleth...	116 C7H16O	017348-59-3	28
4	Acetic acid, cyano-, 1,1-dimethy...	141 C7H11NO2	001116-98-9	17
5	Morpholine, 4-methyl-	101 C5H11NO	000109-02-4	9



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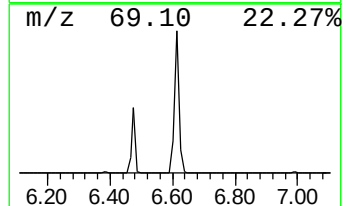
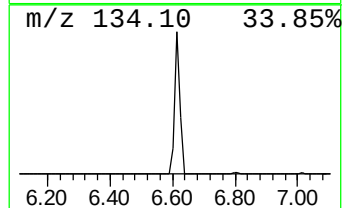
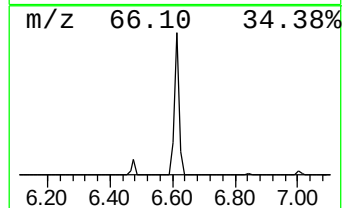
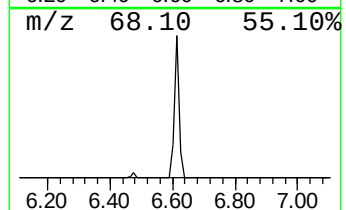
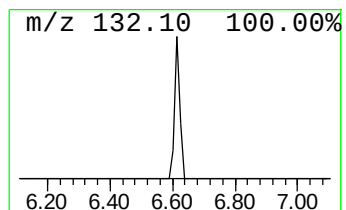
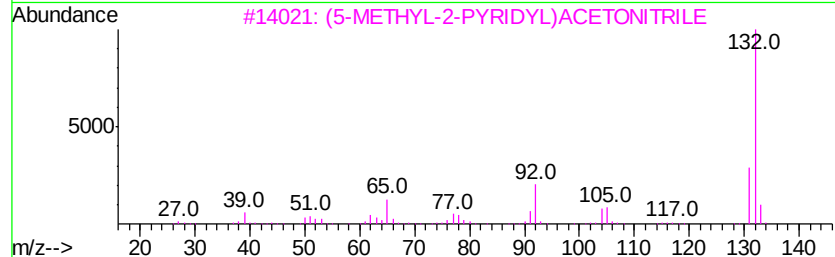
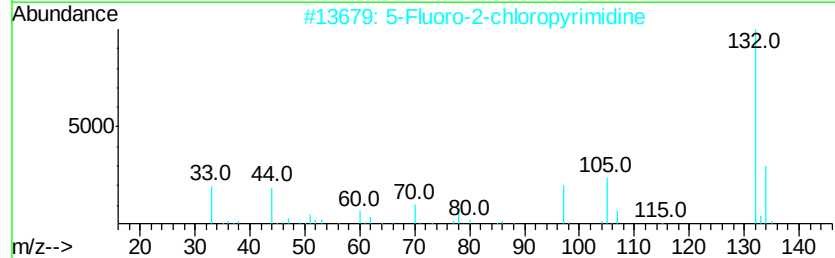
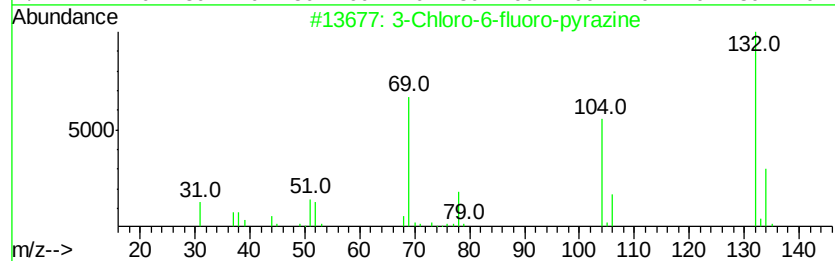
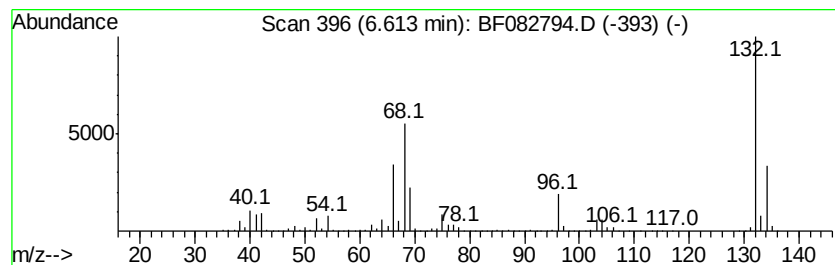
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.61	77.13 ng	6069970	1,4-Dichlorobenzene-d4	6.84

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	12
3		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
4		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10



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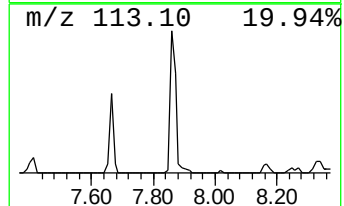
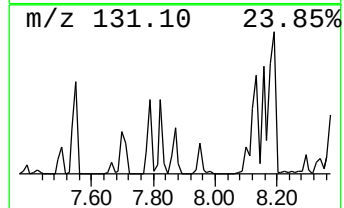
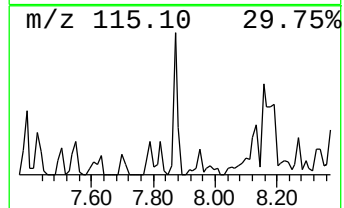
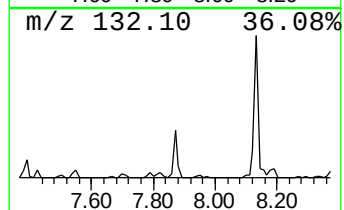
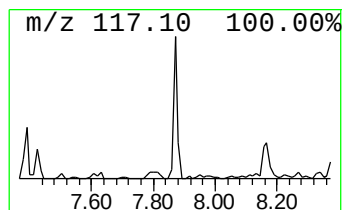
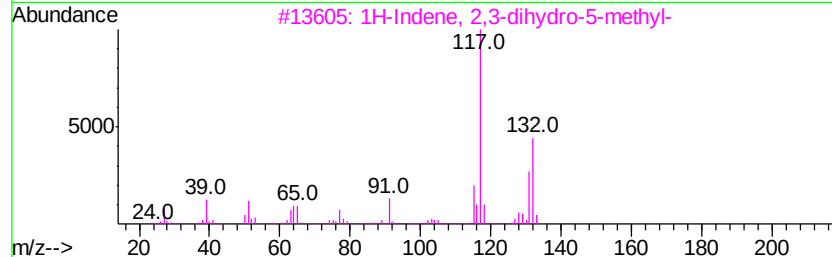
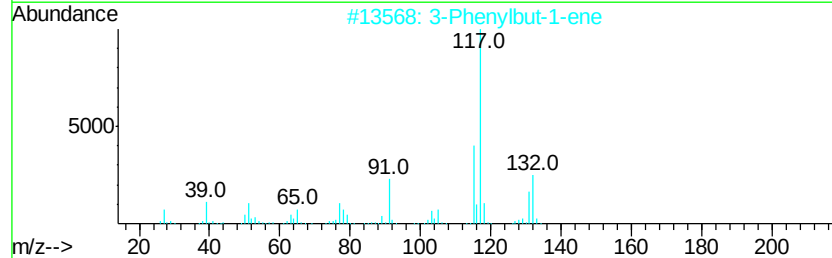
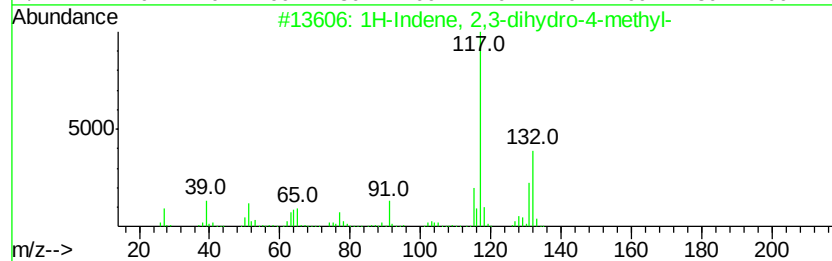
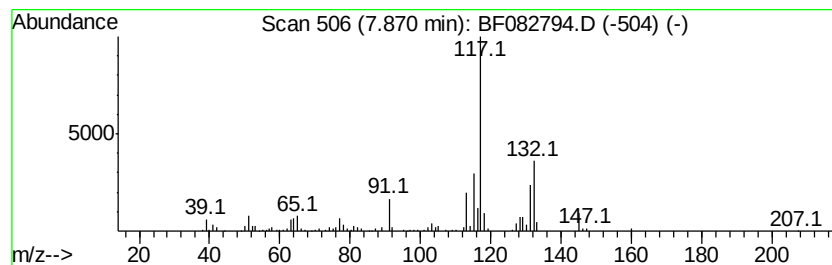
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.87	2.03 ng	236846	Napthalene-d8	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	81
4		Indan, 1-methyl-	132	C10H12	000767-58-8	70
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	68



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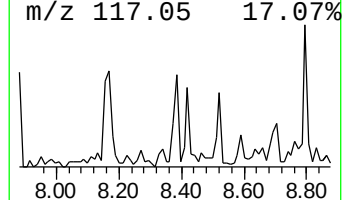
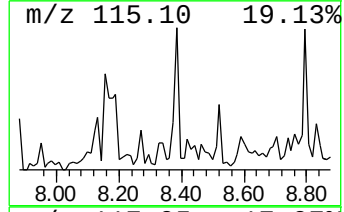
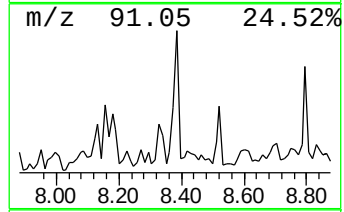
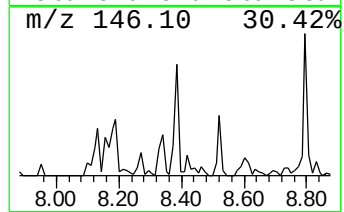
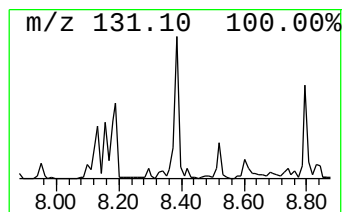
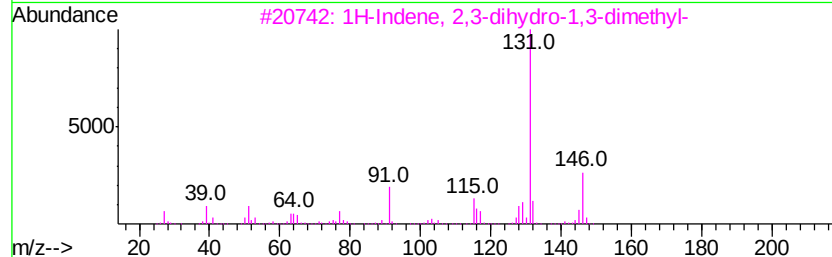
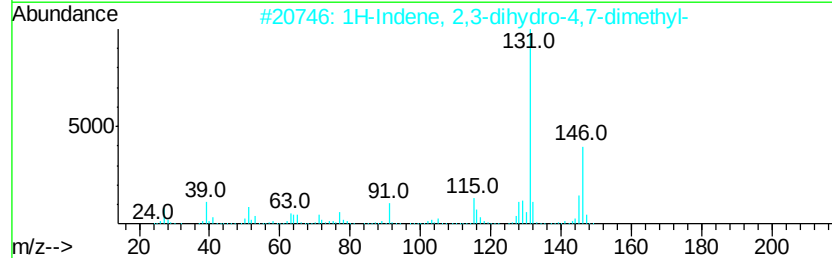
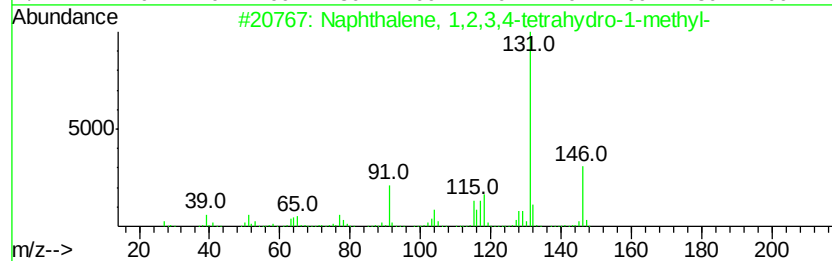
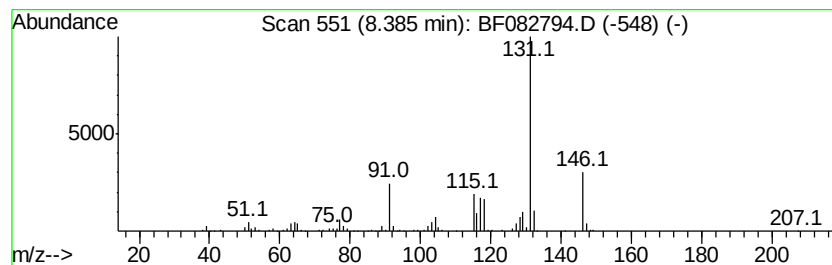
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.38	2.04 ng	237712	Naphthalene-d8	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro...	146	C11H14	001559-81-5	97
2		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87
3		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	83
4		Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	83
5		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	83



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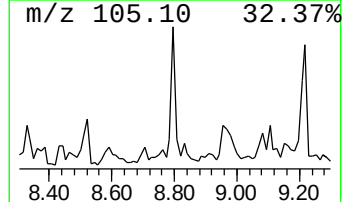
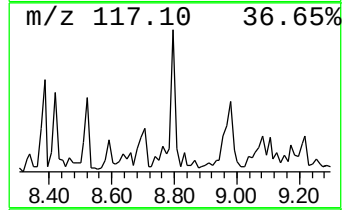
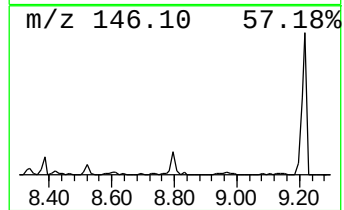
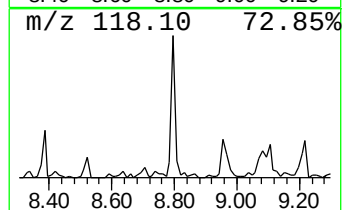
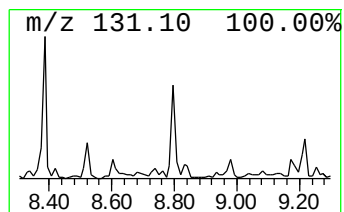
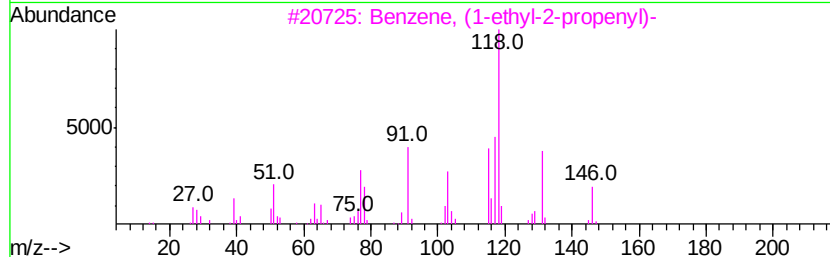
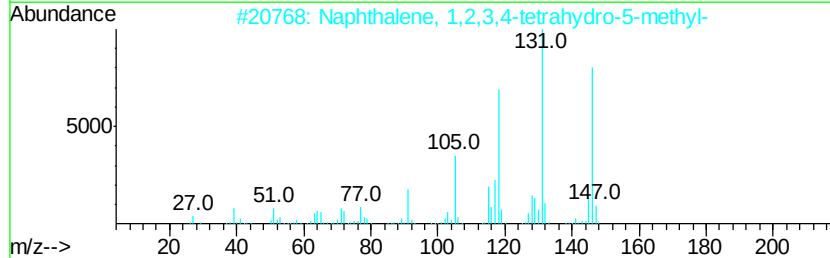
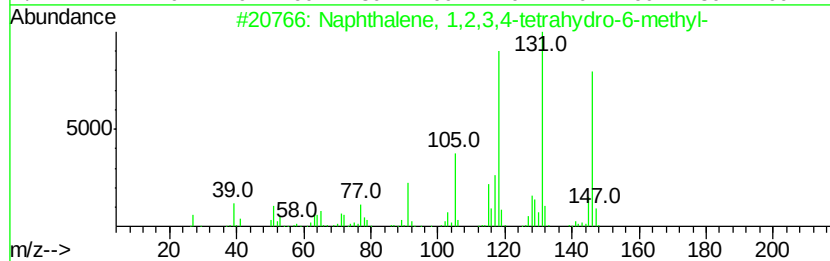
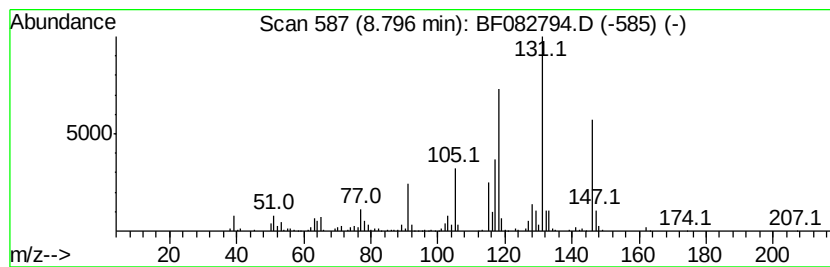
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Naphthalene, 1,2,3,4-tetra... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.80	2.14 ng	249295	Naphthalene-d8	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	94
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	93
3		Benzene, (1-ethyl-2-propenyl)-	146	C11H14	019947-22-9	76
4		1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	76
5		1H-Indene, 2,3-dihydro-2,2-dimethyl-	146	C11H14	020836-11-7	70



Data Path : Z:\HPCHEM1\BNA F\Data\BF110815\
Data File : BF082794.D
Acq On : 7 Nov 2015 7:10
Operator : UM/IZ
Sample : G4259-02
Misc :
ALS Vial : 32 Sample Multiplier: 1

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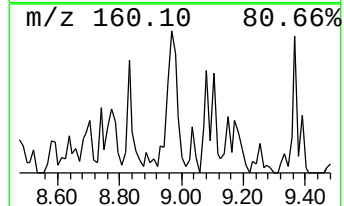
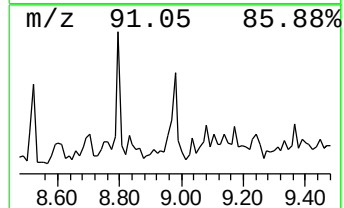
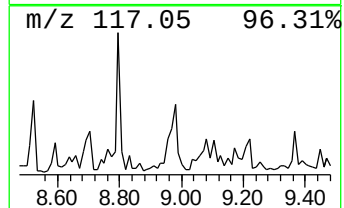
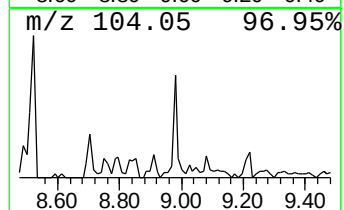
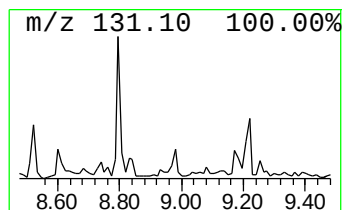
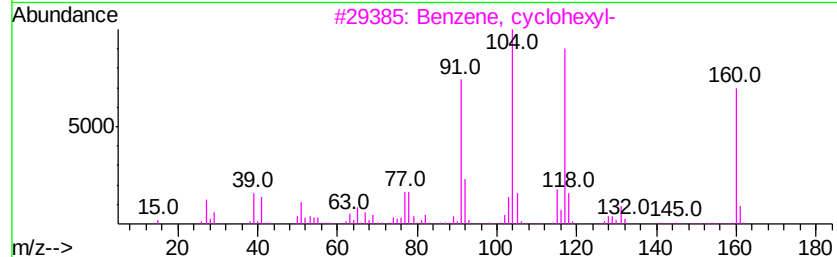
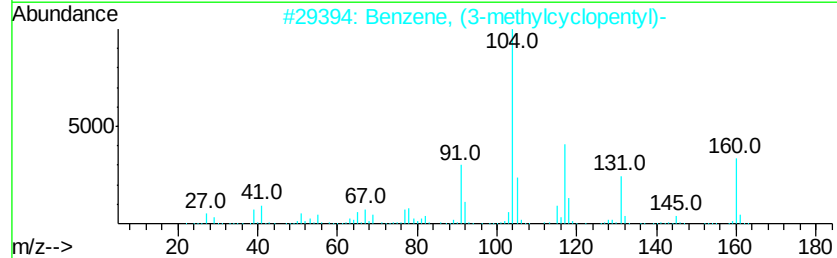
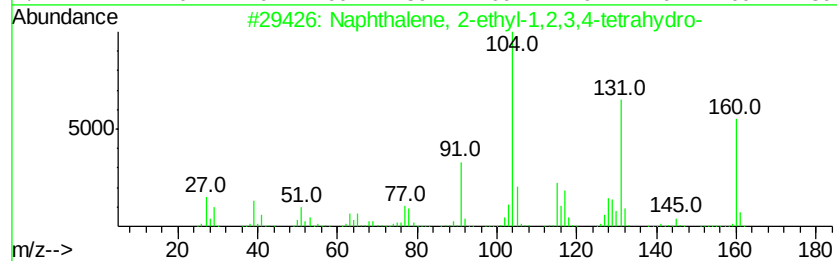
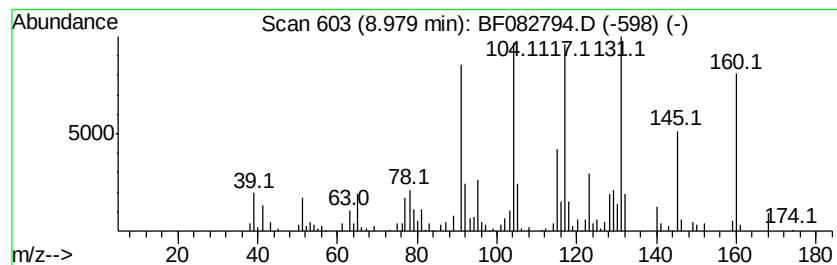
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Naphthalene, 2-ethyl-1,2,3,... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.98	2.65 ng	308993	Naphthalene-d8	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-ethyl-1,2,3,4-tet...	160	C12H16	032367-54-7	83
2		Benzene, (3-methylcyclopentyl)-	160	C12H16	005078-75-1	58
3		Benzene, cyclohexyl-	160	C12H16	000827-52-1	49
4		Quinoxaline, 2-methyl-, 4-oxide	160	C9H8N2O	018916-45-5	27
5		Naphthalene, 1,2-dihydro-6-methoxy-	160	C11H12O	060573-58-2	25



Data Path : Z:\HPCHEM1\BNA F\Data\BF110815\
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Operator : UM/IZ
Sample : G4259-02
Misc :
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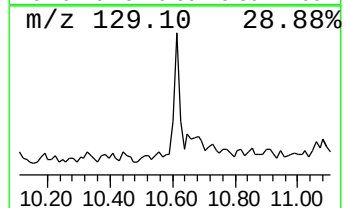
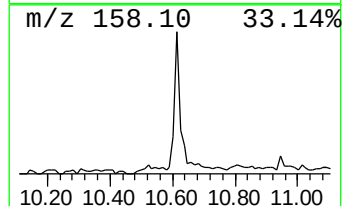
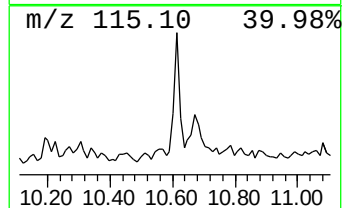
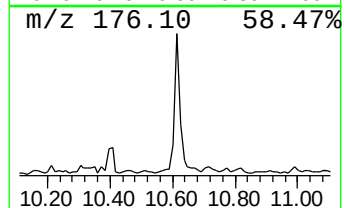
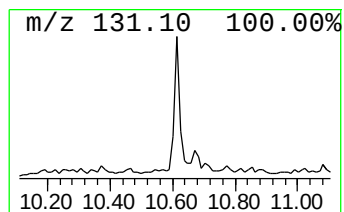
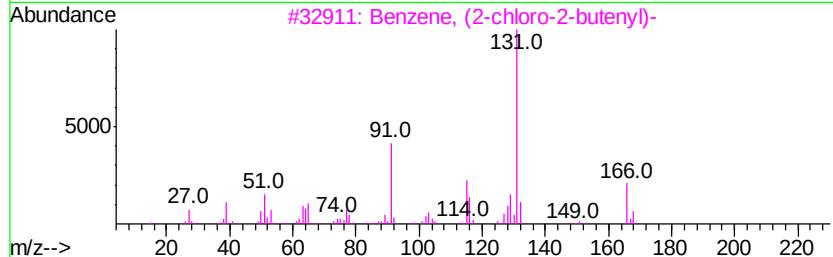
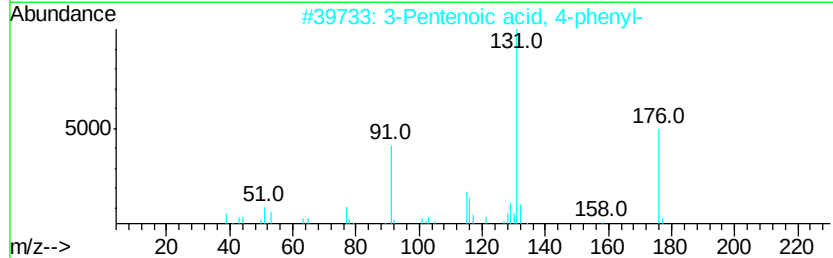
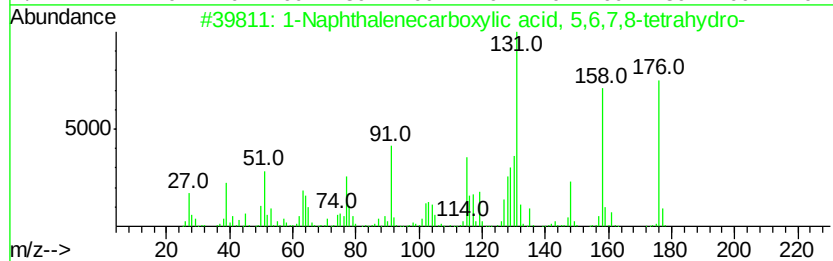
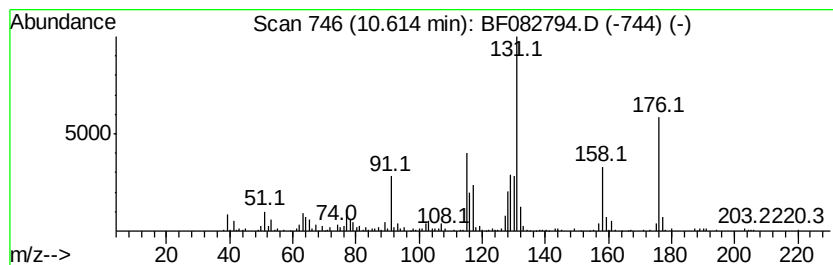
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Naphthalenecarboxylic aci... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.61	2.97 ng	349060	Acenaphthene-d10	9.88
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1-Naphthalenecarboxylic acid, 5,...	176 C11H12O2	004242-18-6 72
2		3-Pentenoic acid, 4-phenyl-	176 C11H12O2	053774-19-9 47
3		Benzene, (2-chloro-2-butenyl)-	166 C10H11Cl	054411-12-0 38
4		Benzene, (1,2-dicyclopropyl-2-ph...	262 C20H22	110330-90-0 38
5		Cinnamic acid, .alpha.,.beta.-di...	176 C11H12O2	1000151-56-4 37



Data Path : Z:\HPCHEM1\BNA F\Data\BF110815\
Data File : BF082794.D
Acq On : 7 Nov 2015 7:10
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Sample : G4259-02
Misc :
ALS Vial : 32 Sample Multiplier: 1

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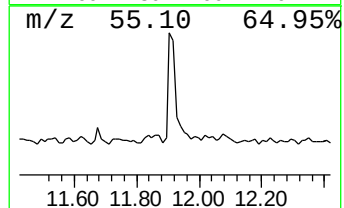
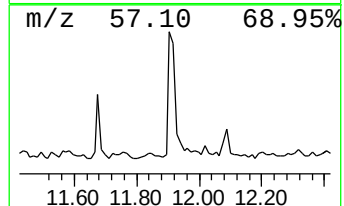
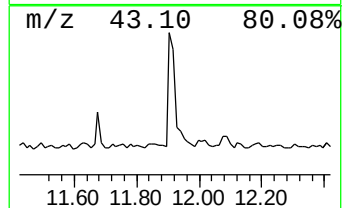
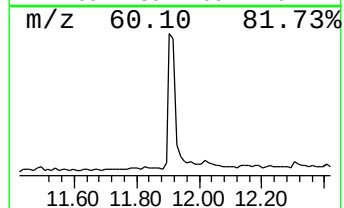
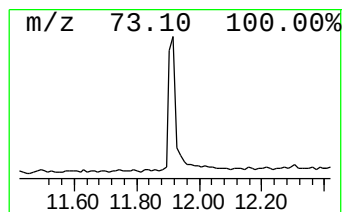
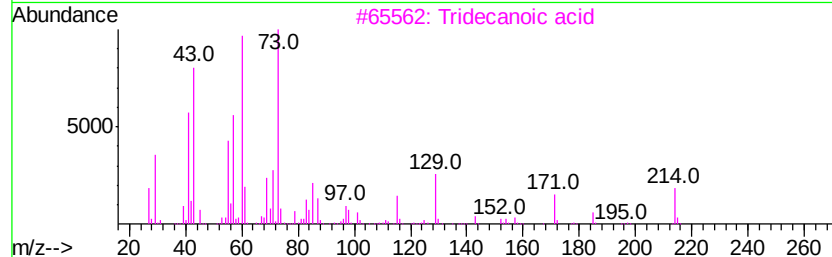
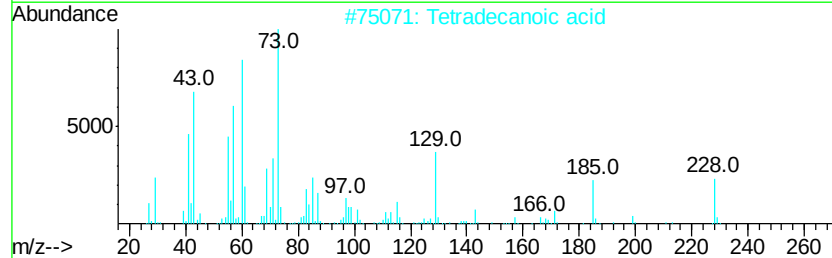
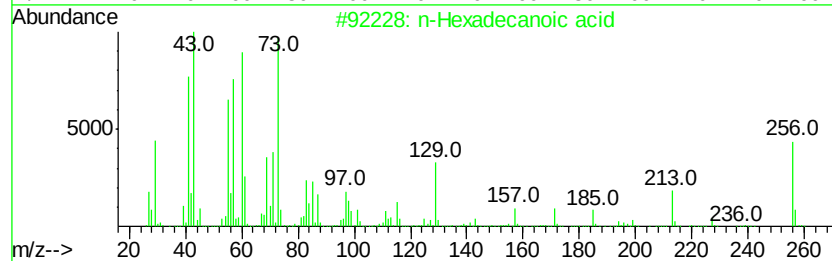
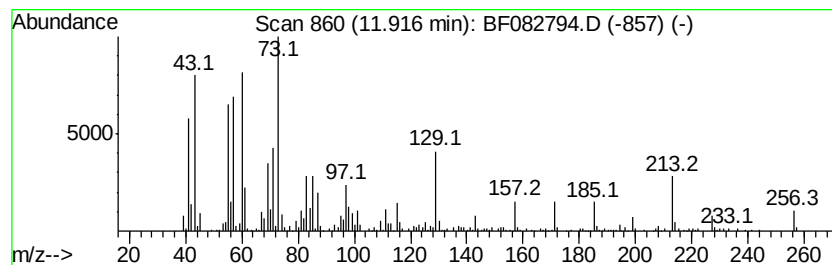
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.92	5.96 ng	686291	Phenanthrene-d10	11.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	94
3		Tridecanoic acid	214	C13H26O2	000638-53-9	90
4		Undecanoic acid	186	C11H22O2	000112-37-8	83
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	81



Data Path : Z:\HPCHEM1\BNA F\Data\BF110815\
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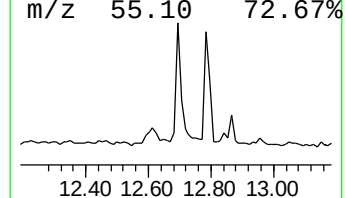
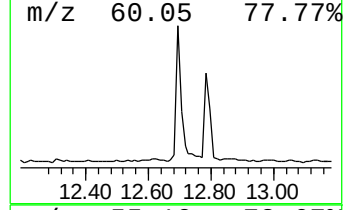
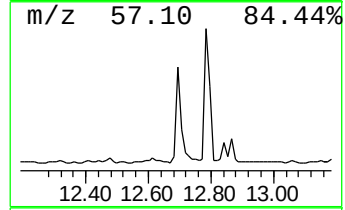
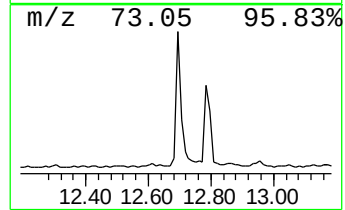
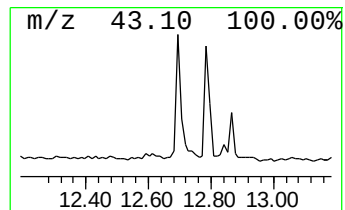
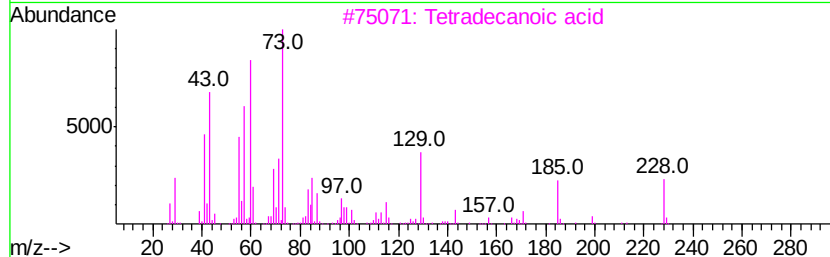
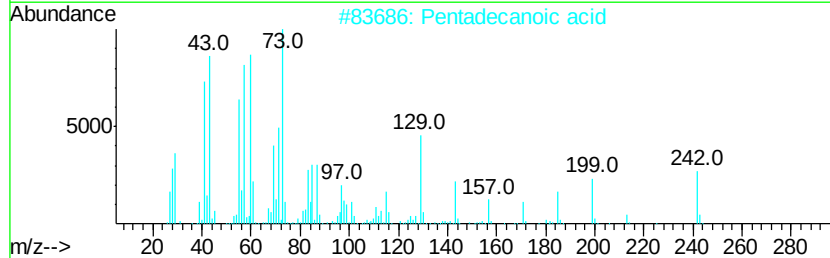
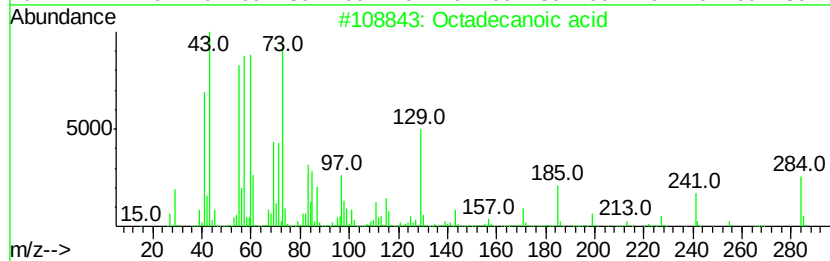
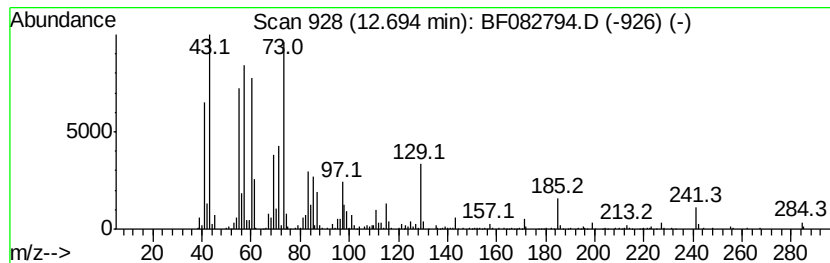
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.69	6.25 ng	702789	Chrysene-d12	14.01	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	98
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	58
4	Undecane	156	C11H24	001120-21-4	11
5	Tetradecane, 2-methyl-	212	C15H32	001560-95-8	11



Data Path : Z:\HPCHEM1\BNA F\Data\BF110815\
Data File : BF082794.D
Acq On : 7 Nov 2015 7:10
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Misc :
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ClientSampleId :
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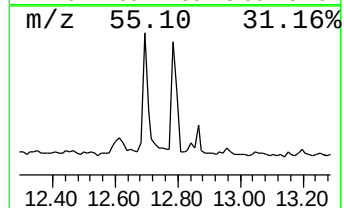
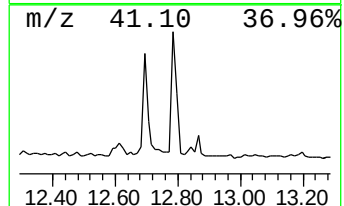
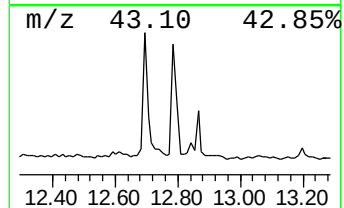
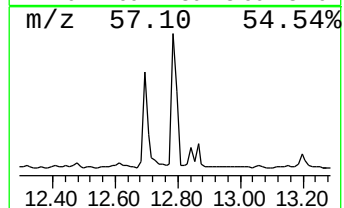
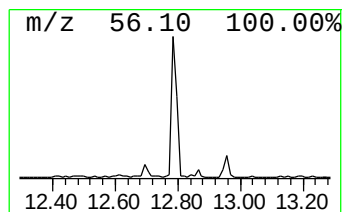
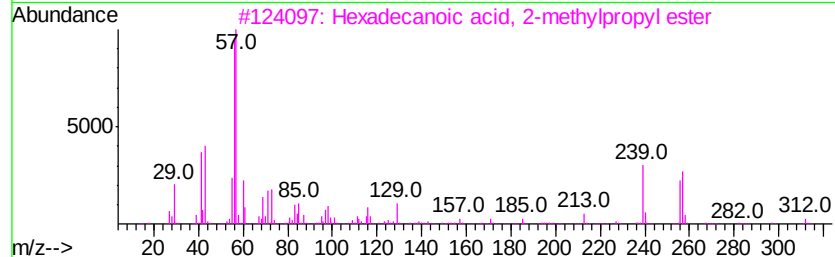
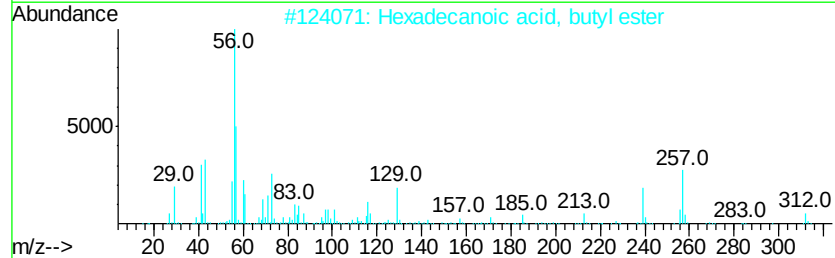
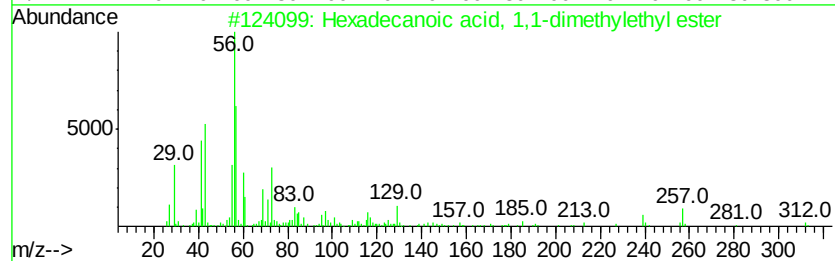
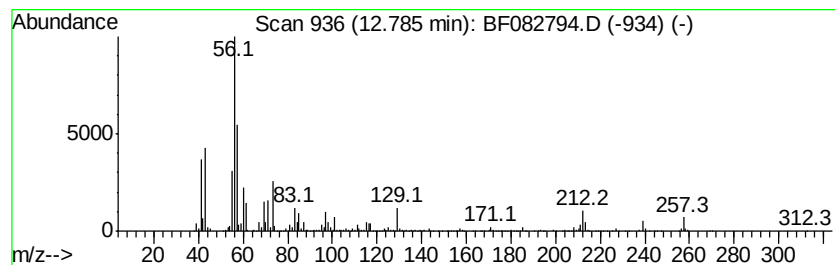
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexadecanoic acid, 1,1-dime... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.79	7.76 ng	872170	Chrysene-d12	14.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	91
2		Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	90
3		Hexadecanoic acid, 2-methylpropyl...	312	C20H40O2	000110-34-9	64
4		Pentadecane, 8-methylene-	224	C16H32	055668-09-2	43
5		2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	40



Data Path : Z:\HPCHEM1\BNA_F\Data\BF110815\
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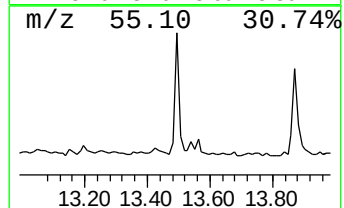
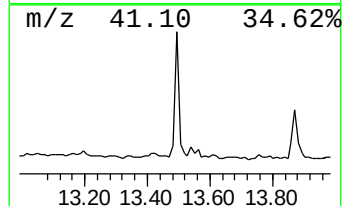
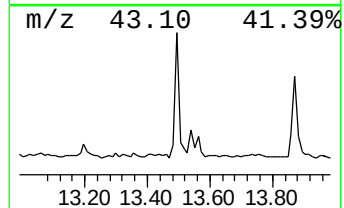
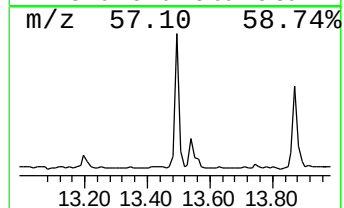
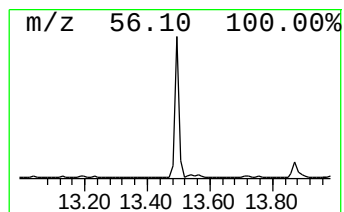
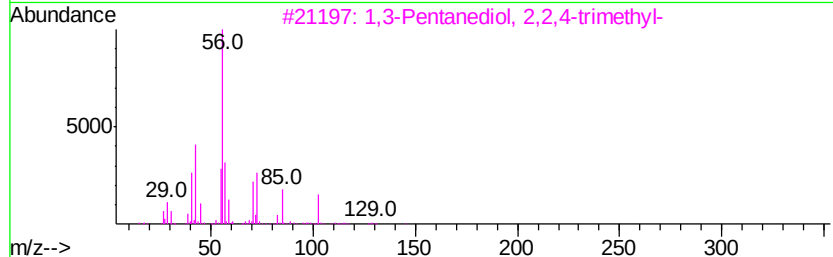
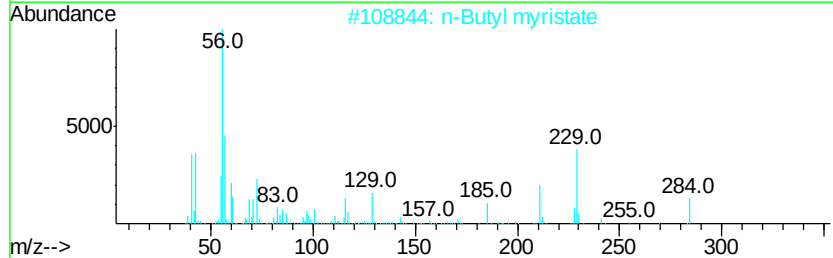
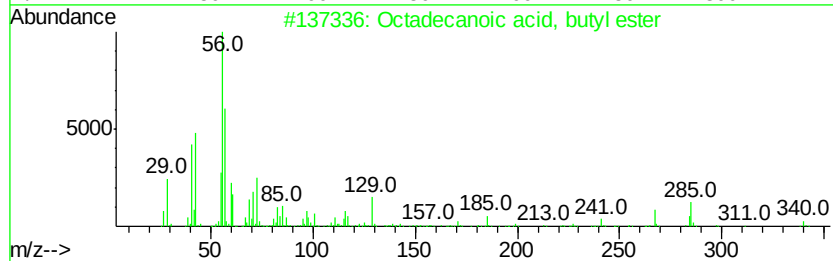
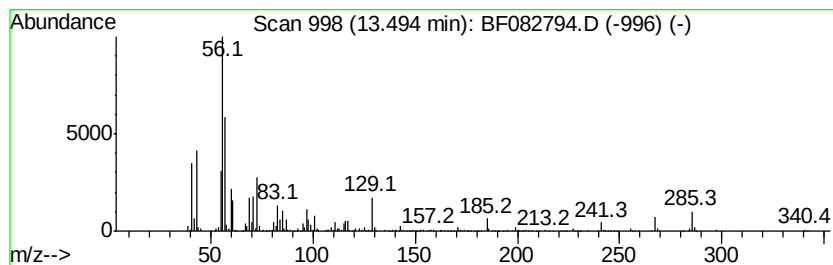
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octadecanoic acid, butyl ester Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	6.09 ng	685264	Chrysene-d12	14.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	93
2		n-Butyl myristate	284	C18H36O2	000110-36-1	86
3		1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	50
4		Cyclobutane, 1,1-dimethyl-2-octyl-	196	C14H28	062338-30-1	38
5		1-Pentadecene, 2-methyl-	224	C16H32	029833-69-0	37



Data Path : Z:\HPCHEM1\BNA F\Data\BF110815\
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Misc :
ALS Vial : 32 Sample Multiplier: 1

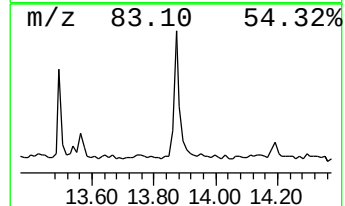
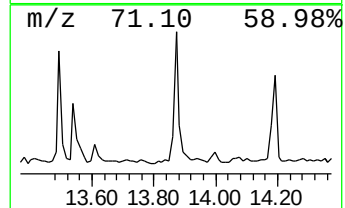
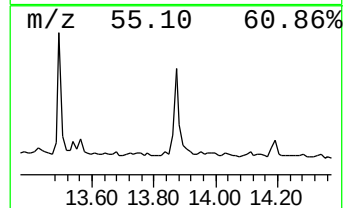
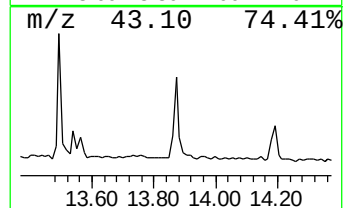
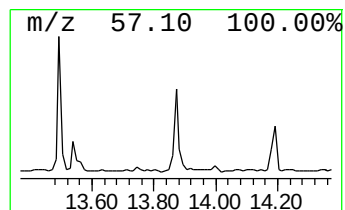
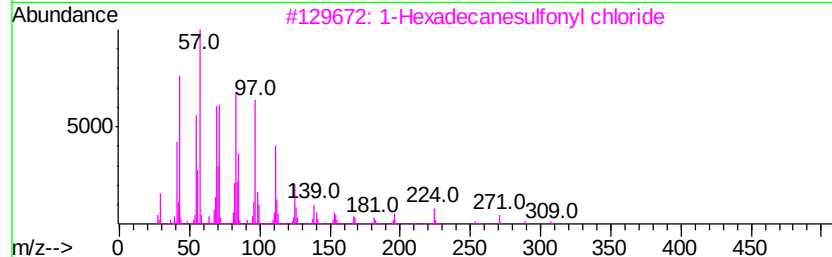
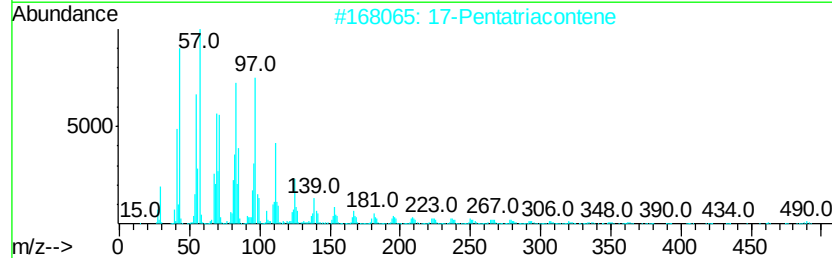
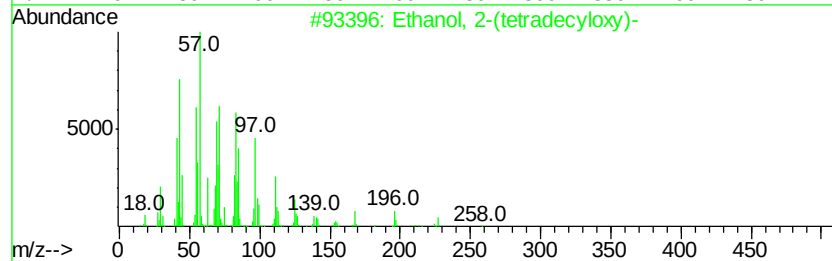
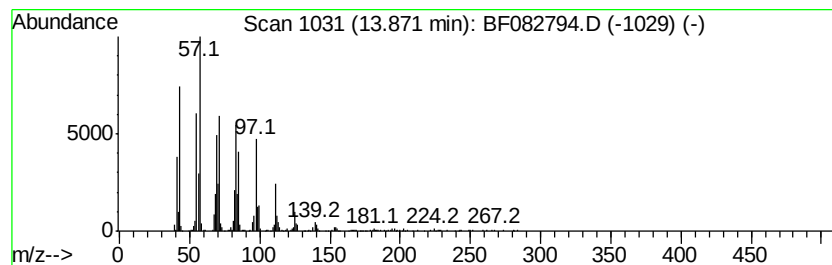
Instrument :
BNA_F
ClientSampleId :
MW-01-110315

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Ethanol, 2-(tetradecyloxy)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.87	3.82 ng	429007	Chrysene-d12	14.01		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
2		17-Pentatriacontene	491	C35H70	006971-40-0	87
3		1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	87
4		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	81
5		Cyclohexadecane	224	C16H32	000295-65-8	78



Data Path : Z:\HPCHEM1\BNA_F\Data\BF110815\
Data File : BF082794.D
Acq On : 7 Nov 2015 7:10
Operator : UM/IZ
Sample : G4259-02
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-01-110315

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.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...	5.02	19.9	ng	1569780	1	6.84	1573990	20.0
unknown6.61	6.61	77.1	ng	6069970	1	6.84	1573990	20.0
1H-Indene, 2,3-di...	7.87	2.0	ng	236846	2	8.13	2330390	20.0
Naphthalene, 1,2,...	8.38	2.0	ng	237712	2	8.13	2330390	20.0
Naphthalene, 1,2,...	8.80	2.1	ng	249295	2	8.13	2330390	20.0
Naphthalene, 2-et...	8.98	2.6	ng	308993	2	8.13	2330390	20.0
1-Naphthalenecarb...	10.61	3.0	ng	349060	3	9.88	2350070	20.0
n-Hexadecanoic acid	11.92	6.0	ng	686291	4	11.37	2304200	20.0
Octadecanoic acid	12.69	6.3	ng	702789	5	14.01	2248890	20.0
Hexadecanoic acid...	12.79	7.8	ng	872170	5	14.01	2248890	20.0
Octadecanoic acid...	13.49	6.1	ng	685264	5	14.01	2248890	20.0
Ethanol, 2-(tetra...	13.87	3.8	ng	429007	5	14.01	2248890	20.0