

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042716\
 Data File : BF086459.D
 Acq On : 28 Apr 2016 8:05
 Operator : UM/SJ
 Sample : H2736-02
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RICHMOND-AMBOY-SB02(0.5-8.0)

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-BF042716.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.990	251	254	260	rBV	63928	94564	2.72%	0.336%
2	5.402	287	290	292	rBV	3247747	3470268	100.00%	12.343%
3	6.442	378	381	383	rBV	3129929	3369386	97.09%	11.984%
4	6.567	390	392	395	rBV	2588071	3230322	93.09%	11.489%
5	6.807	411	413	415	rBV	717798	768653	22.15%	2.734%
6	6.956	424	426	429	rBV	2014681	2465979	71.06%	8.771%
7	7.367	459	462	469	rBV	1829241	1964918	56.62%	6.989%
8	7.470	469	471	478	rVB	41614	90196	2.60%	0.321%
9	8.088	523	525	528	rBV	924912	901010	25.96%	3.205%
10	9.162	617	619	622	rBV	1879907	2737832	78.89%	9.738%
11	9.562	652	654	656	rBV	423307	362030	10.43%	1.288%
12	9.779	671	673	675	rBV	69917	60962	1.76%	0.217%
13	9.836	676	678	681	rVB	542504	766912	22.10%	2.728%
14	10.282	715	717	718	rVB	126345	96439	2.78%	0.343%
15	10.499	733	736	739	rBV	44398	72958	2.10%	0.259%
16	10.625	745	747	750	rBV2	1279533	1482665	42.72%	5.273%
17	10.751	757	758	765	rVB	150930	185525	5.35%	0.660%
18	11.196	796	797	799	rBV	72085	71689	2.07%	0.255%
19	11.322	806	808	811	rBV	677425	671024	19.34%	2.387%
20	11.562	826	829	833	rBV	98036	178361	5.14%	0.634%
21	11.631	833	835	838	rVB	27148	41810	1.20%	0.149%
22	12.385	898	901	909	rBV2	88276	208679	6.01%	0.742%
23	12.739	929	932	935	rBV2	23030	40340	1.16%	0.143%
24	12.911	944	947	949	rBV	2243316	2251774	64.89%	8.009%
25	13.151	965	968	970	rBV	32103	38187	1.10%	0.136%
26	13.677	1011	1014	1018	rBV	22048	40985	1.18%	0.146%
27	13.814	1024	1026	1034	rBV	119937	228546	6.59%	0.813%
28	13.951	1036	1038	1046	rVV	632310	894742	25.78%	3.182%
29	14.065	1046	1048	1050	rVB	69430	83135	2.40%	0.296%
30	14.431	1078	1080	1082	rBV	31813	51654	1.49%	0.184%
31	14.797	1110	1112	1115	rVB	47585	57150	1.65%	0.203%
32	14.968	1125	1127	1132	rBV	22880	53414	1.54%	0.190%
33	15.048	1132	1134	1137	rBV	97695	110552	3.19%	0.393%
34	15.357	1159	1161	1169	rBV	328699	587392	16.93%	2.089%

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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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042716.M
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35	15.814	1198	1201	1205	rBV	61923	105916	3.05%	0.377%
36	16.054	1217	1222	1227	rBV4	23887	77325	2.23%	0.275%
37	17.300	1328	1331	1337	rVB7	20803	65868	1.90%	0.234%
38	17.791	1371	1374	1380	rVB7	24059	66521	1.92%	0.237%
39	18.214	1408	1411	1419	rVB7	21202	70448	2.03%	0.251%

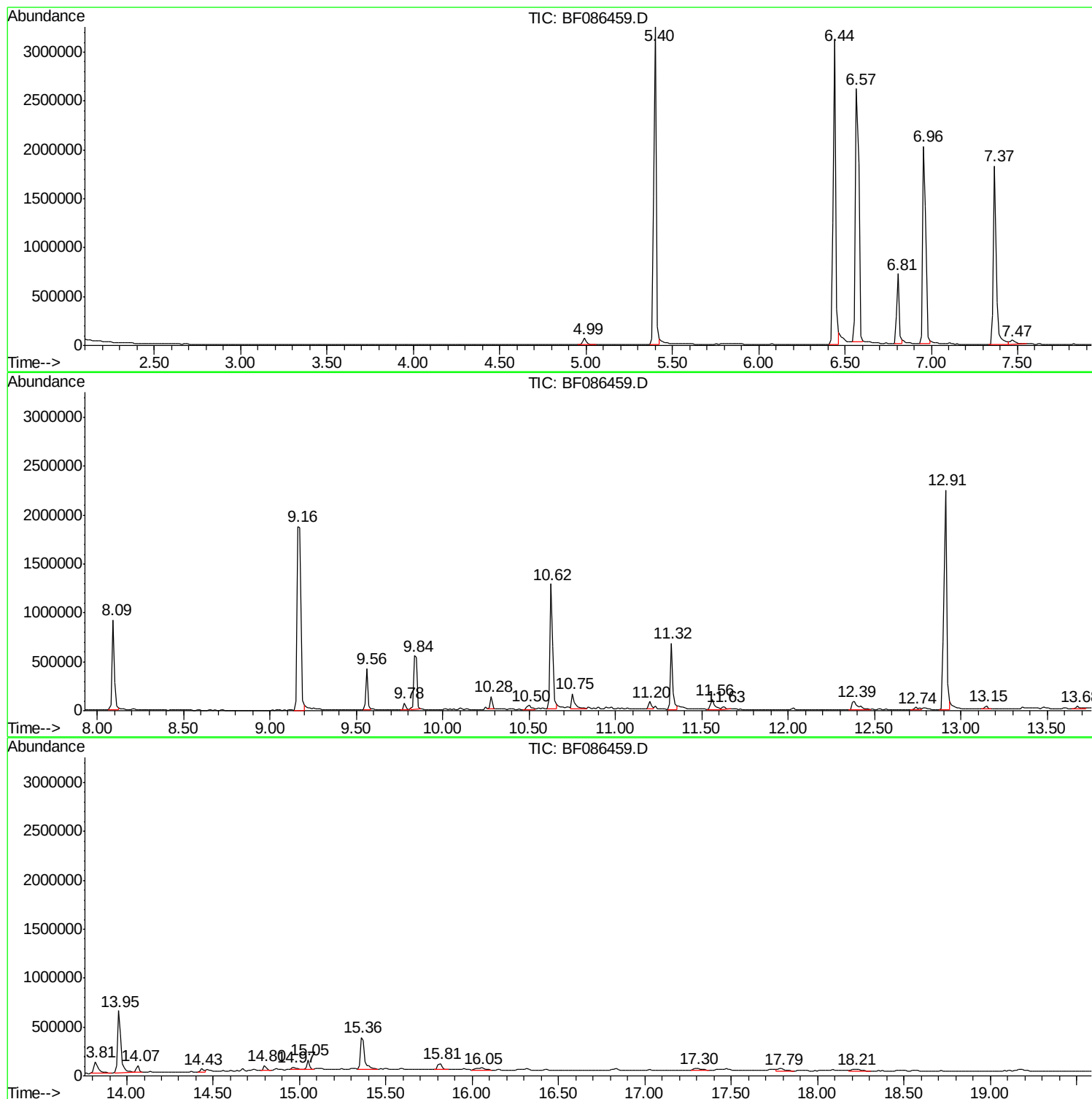
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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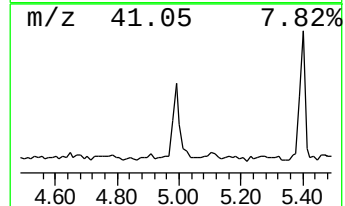
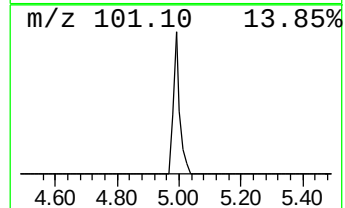
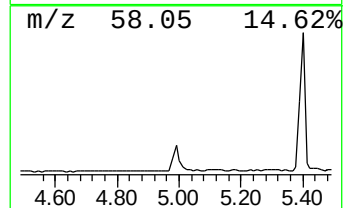
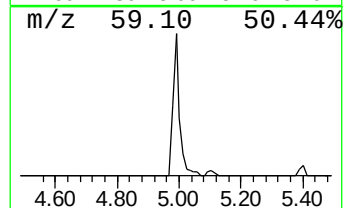
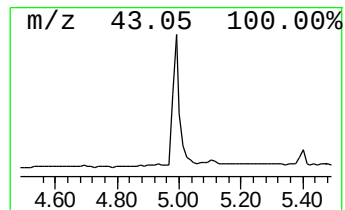
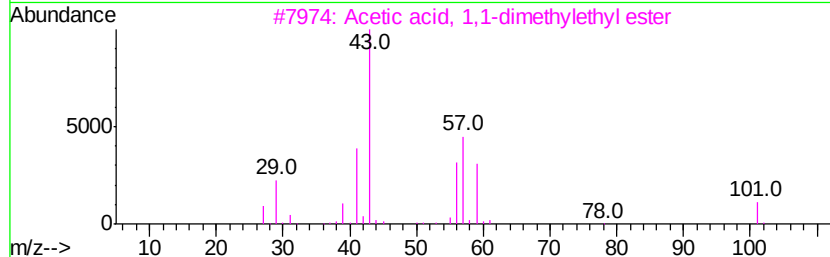
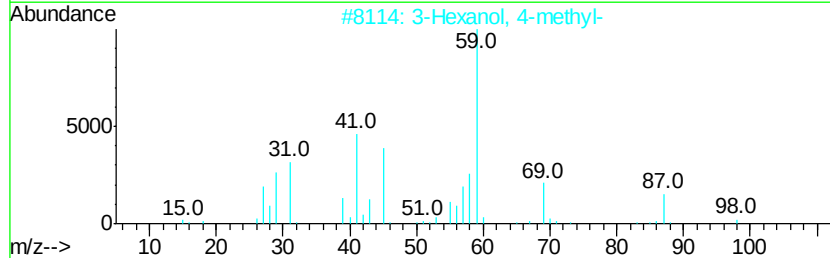
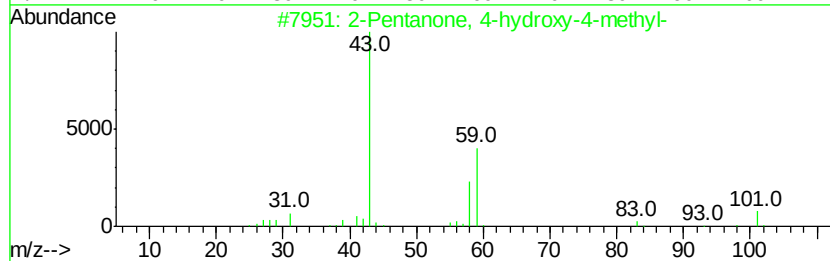
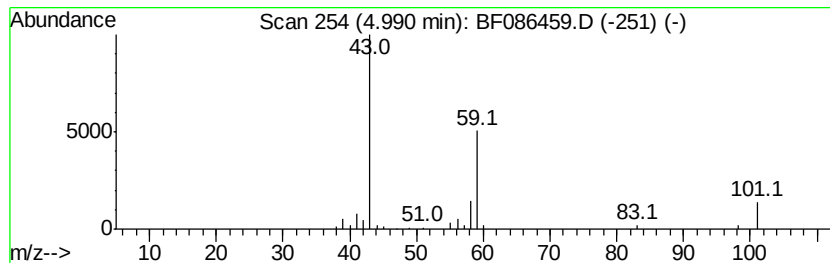
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	2.46 ng	94564	1,4-Dichlorobenzene-d4	6.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
3			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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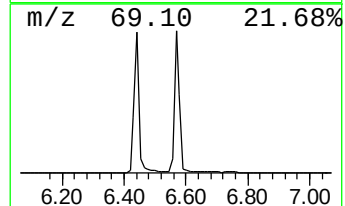
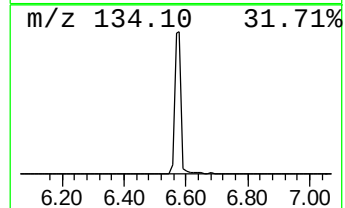
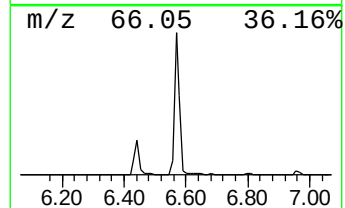
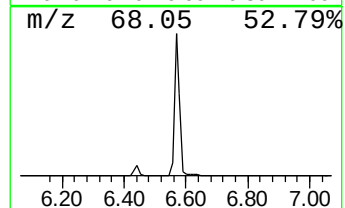
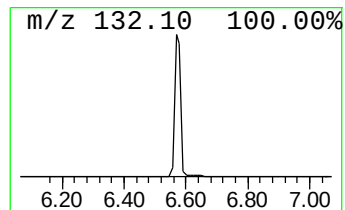
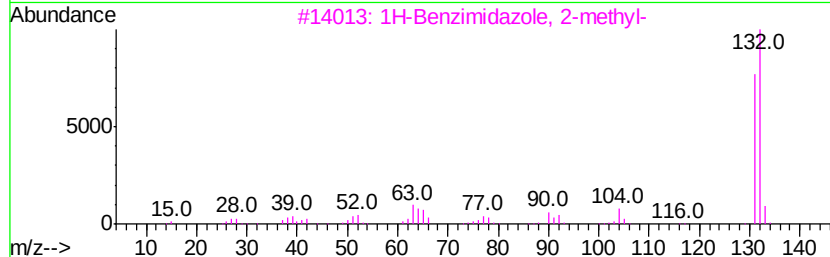
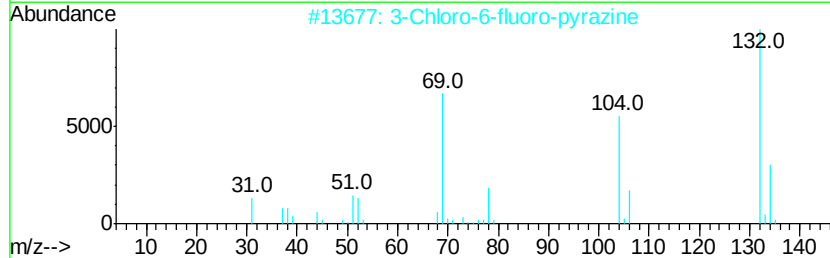
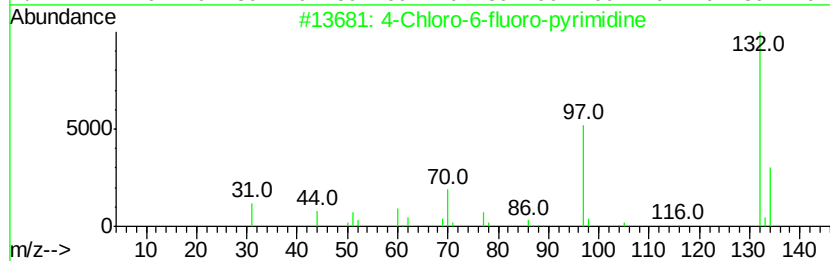
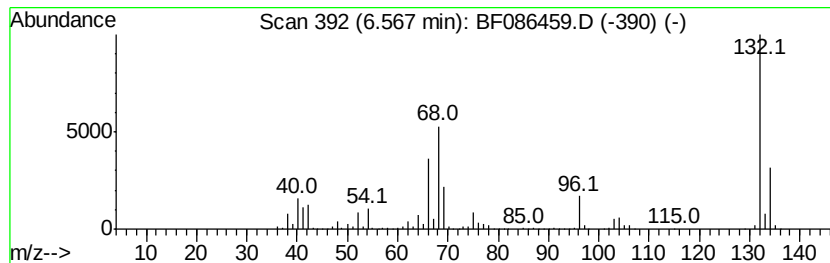
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown6.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	84.05 ng	3230320	1,4-Dichlorobenzene-d4	6.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12



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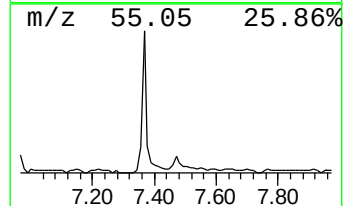
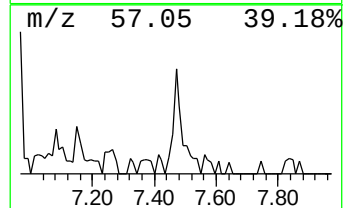
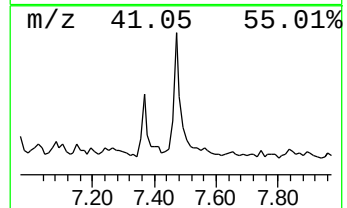
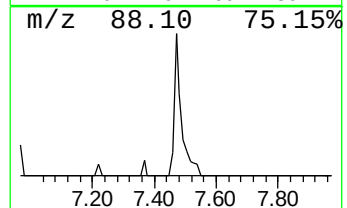
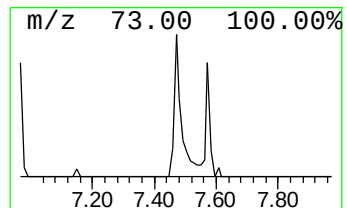
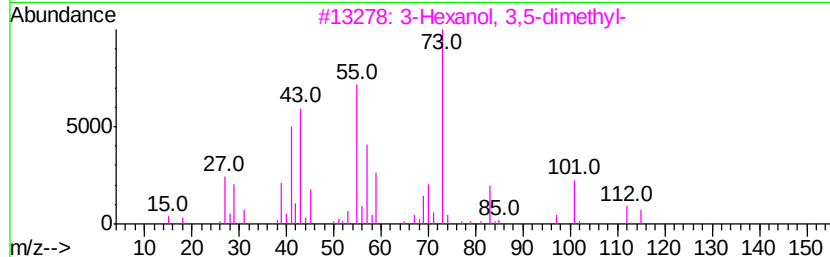
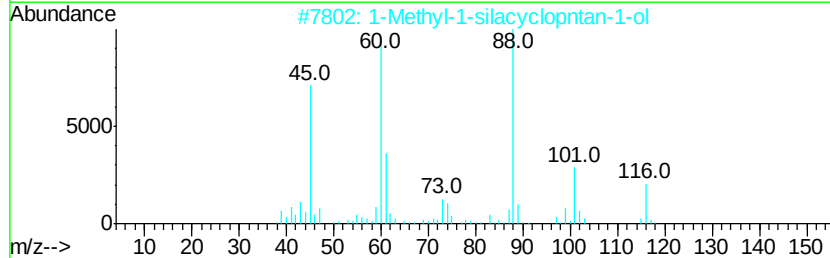
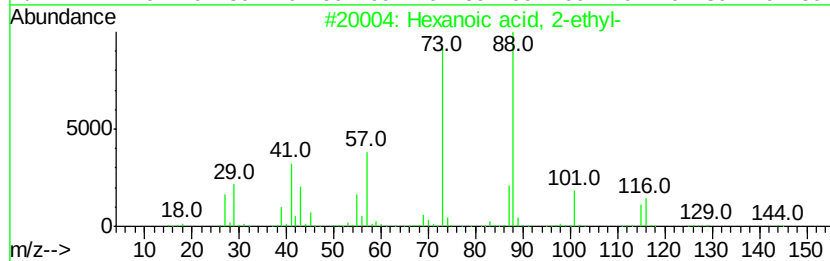
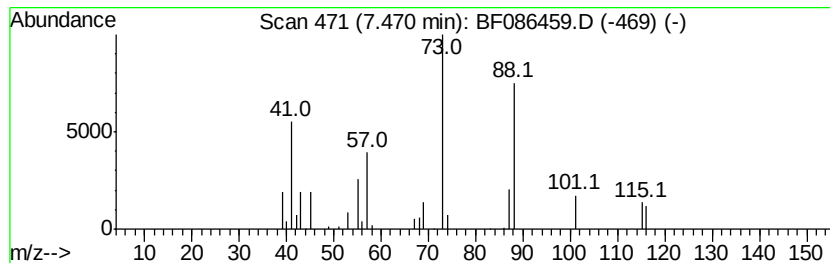
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Hexanoic acid, 2-ethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.47	2.00 ng	90196	Naphthalene-d8	8.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78
2			1-Methyl-1-silacyclopentan-1-ol	116	C5H12OSi	1000216-84-4	28
3			3-Hexanol, 3,5-dimethyl-	130	C8H18O	004209-91-0	10
4			Hexanamide, 6-(heptyloxy)-	229	C13H27NO2	055191-08-7	9
5			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	9



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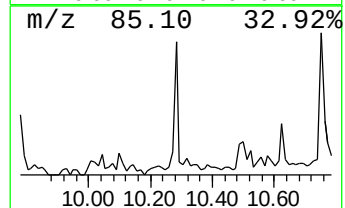
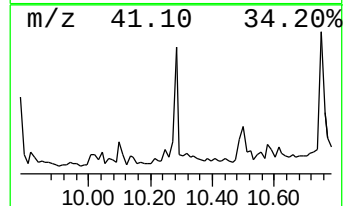
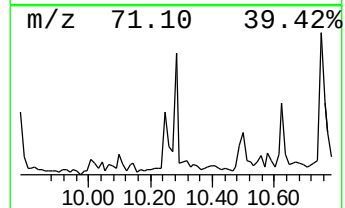
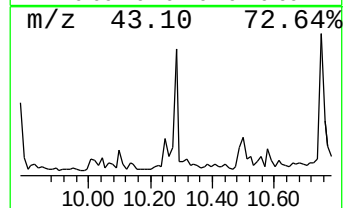
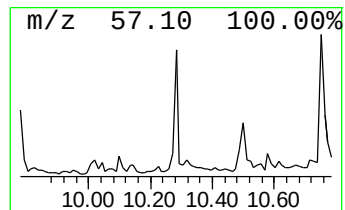
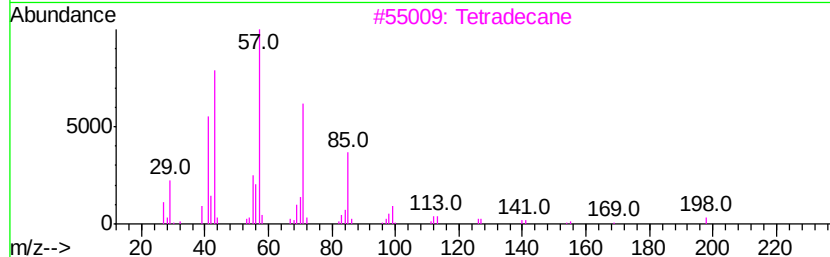
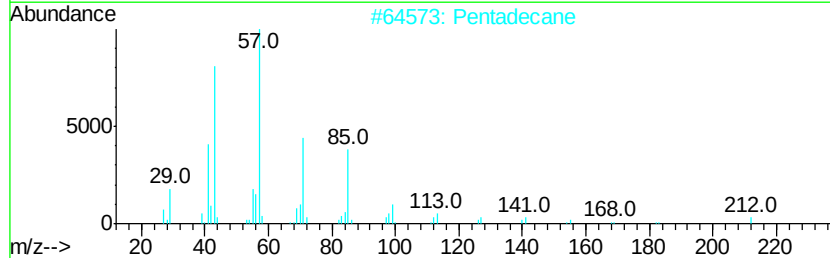
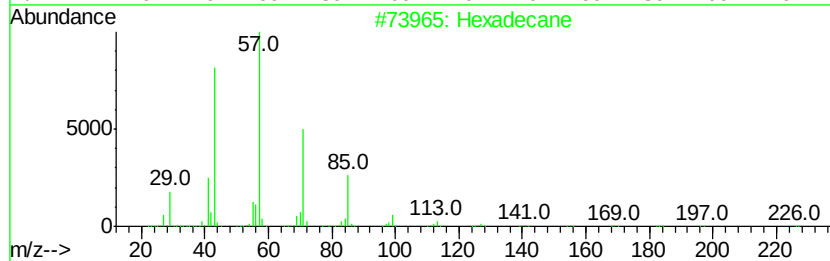
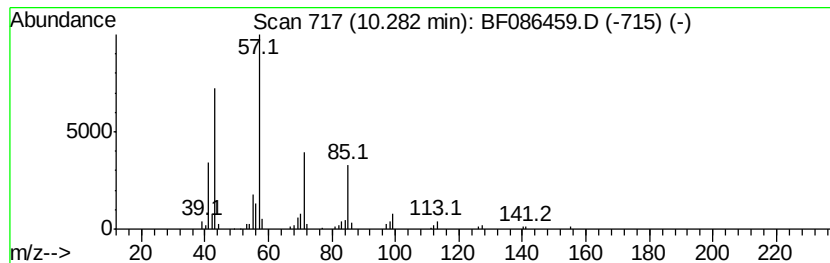
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.28	2.52 ng	96439	Acenaphthene-d10	9.85	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	90
2	Pentadecane	212	C15H32	000629-62-9	90
3	Tetradecane	198	C14H30	000629-59-4	86
4	Octadecane	254	C18H38	000593-45-3	86
5	Eicosane	282	C20H42	000112-95-8	83



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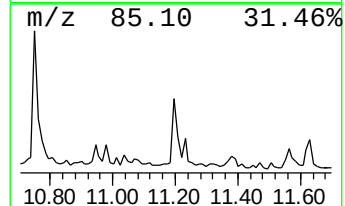
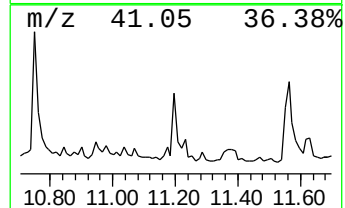
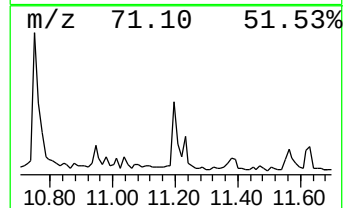
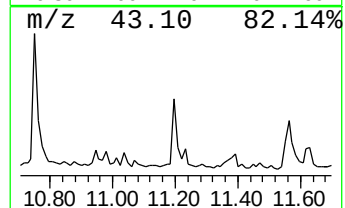
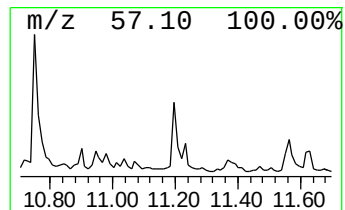
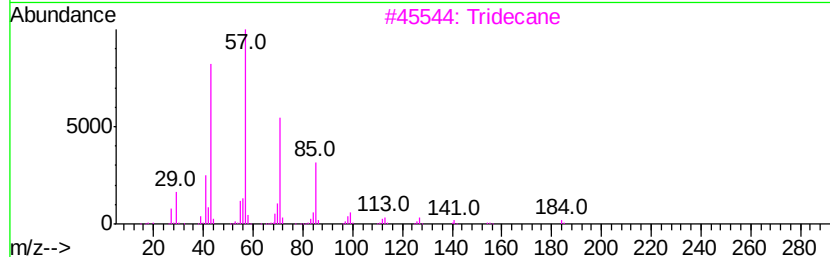
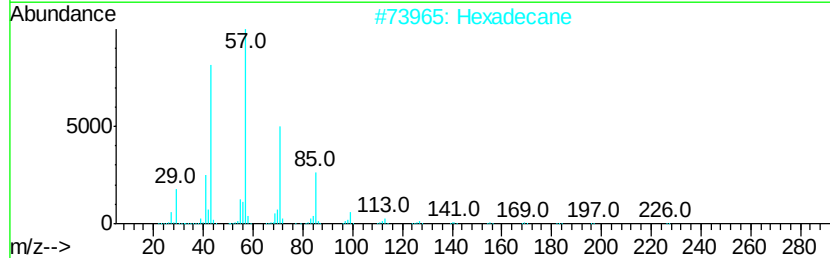
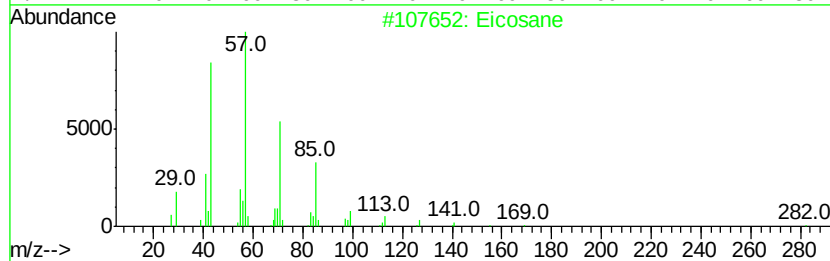
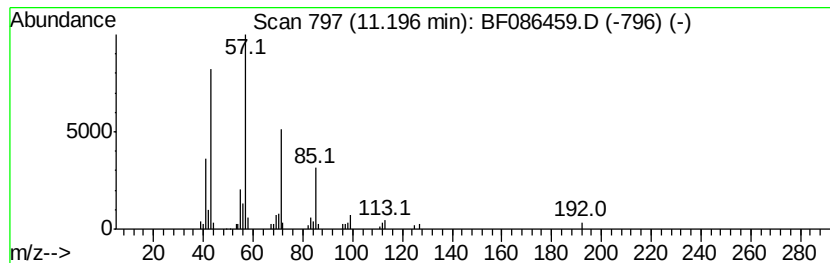
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Peak Number 7 Eicosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.20	2.14 ng	71689	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Hexadecane	226	C16H34	000544-76-3	91
3		Tridecane	184	C13H28	000629-50-5	86
4		Tridecane, 7-propyl-	226	C16H34	055045-09-5	80
5		Octadecane	254	C18H38	000593-45-3	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042716\
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Acq On : 28 Apr 2016 8:05
Operator : UM/SJ
Sample : H2736-02
Misc :
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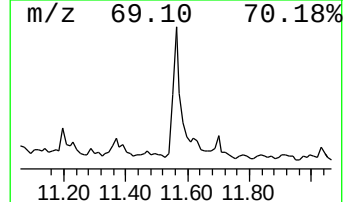
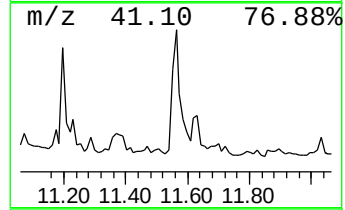
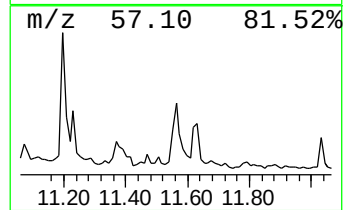
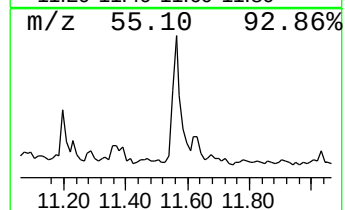
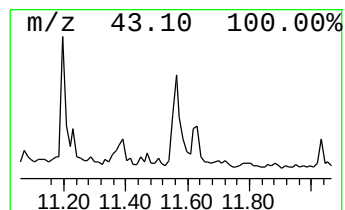
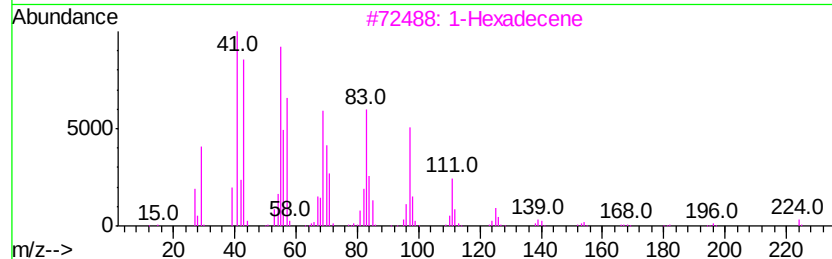
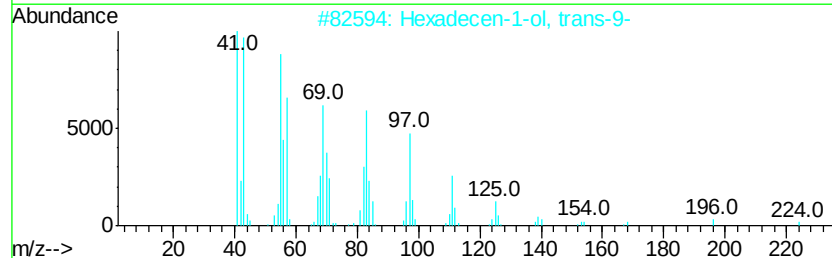
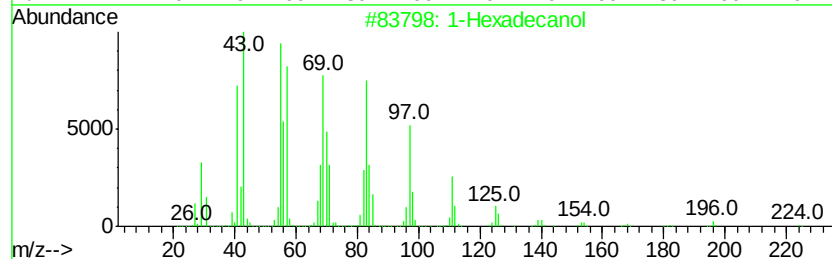
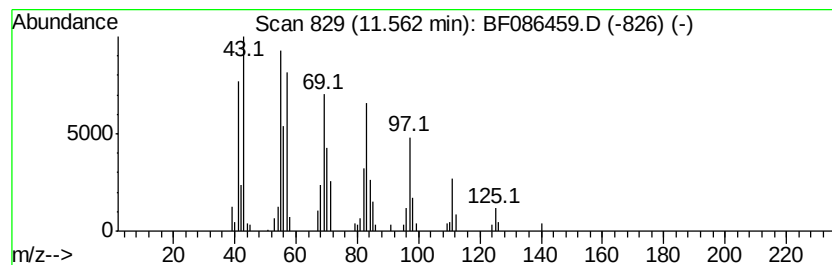
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1-Hexadecanol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.56	5.32 ng	178361	Phenanthrene-d10	11.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexadecanol	242	C16H34O	036653-82-4	95
2	Hexadecen-1-ol, trans-9-	240	C16H32O	064437-47-4	91
3	1-Hexadecene	224	C16H32	000629-73-2	91
4	5-Eicosene, (E)-	280	C20H40	074685-30-6	91
5	Trifluoroacetic acid, n-heptadec...	324	C17H31F3O2	1000216-79-2	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042716\
Data File : BF086459.D
Acq On : 28 Apr 2016 8:05
Operator : UM/SJ
Sample : H2736-02
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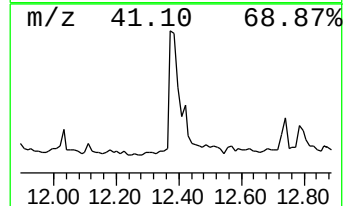
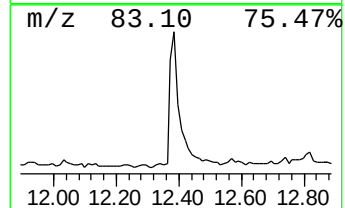
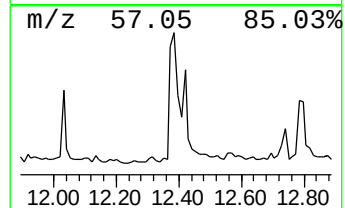
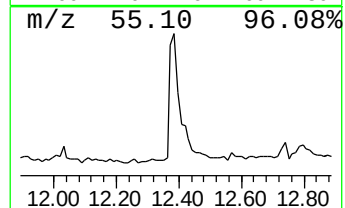
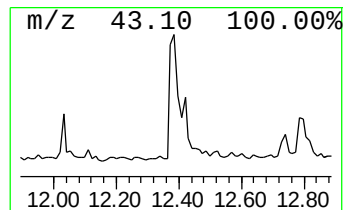
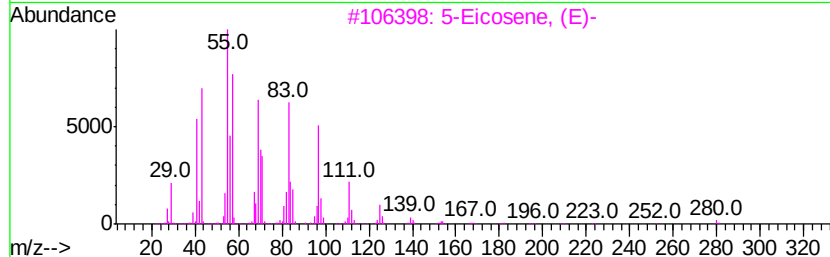
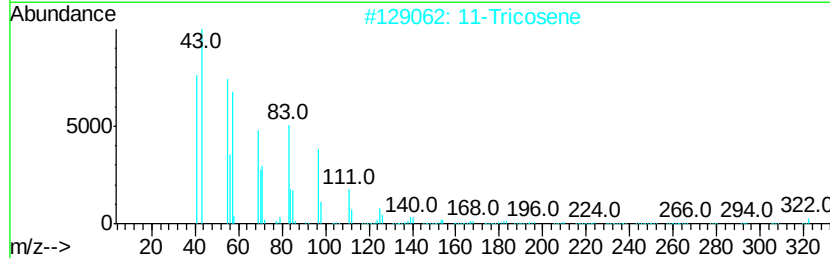
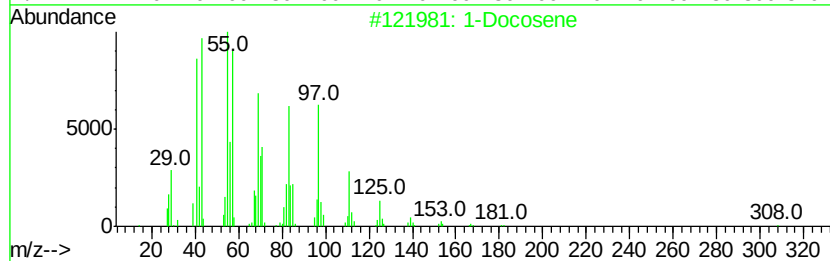
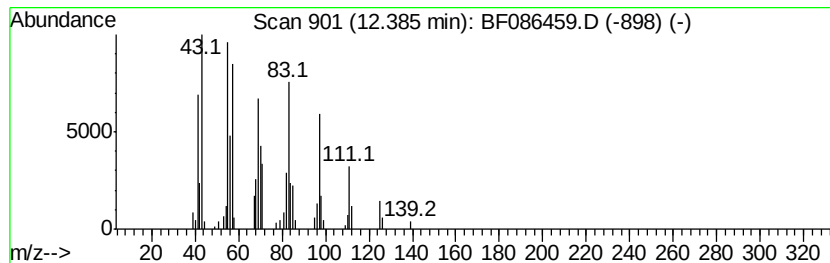
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Docosene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.39	6.22 ng	208679	Phenanthrene-d10	11.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Docosene	308 C22H44	001599-67-3	94
2	11-Tricosene	322 C23H46	052078-56-5	91
3	5-Eicosene, (E)-	280 C20H40	074685-30-6	91
4	3-Eicosene, (E)-	280 C20H40	074685-33-9	91
5	Hexadecen-1-ol, trans-9-	240 C16H32O	064437-47-4	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042716\
Data File : BF086459.D
Acq On : 28 Apr 2016 8:05
Operator : UM/SJ
Sample : H2736-02
Misc :
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Instrument :
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ClientSampleId :
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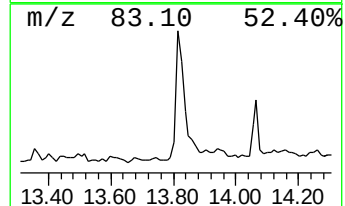
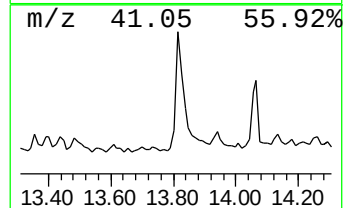
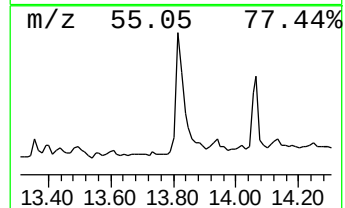
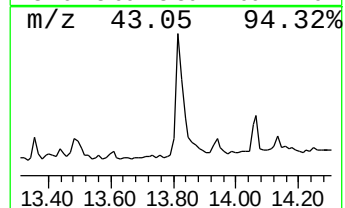
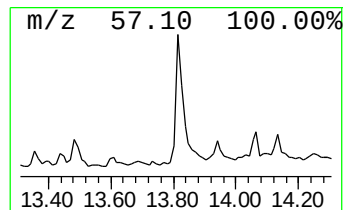
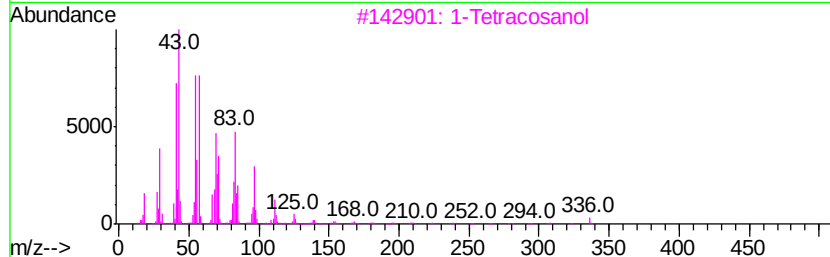
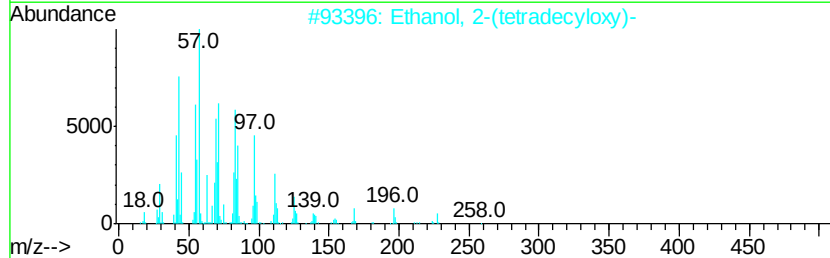
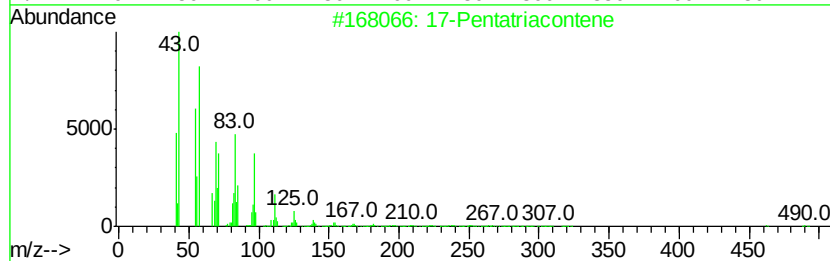
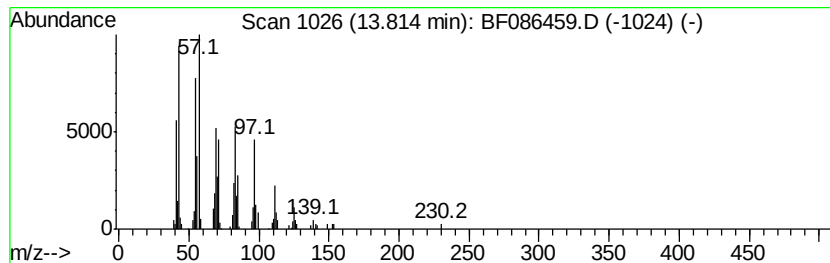
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 17-Pentatriacontene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.81	5.11 ng	228546	Chrysene-d12	13.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			17-Pentatriacontene	491	C35H70	006971-40-0	91
2			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
3			1-Tetracosanol	354	C24H50O	000506-51-4	91
4			Bromoacetic acid, pentadecyl ester	348	C17H33BrO2	131143-01-6	87
5			3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042716\
Data File : BF086459.D
Acq On : 28 Apr 2016 8:05
Operator : UM/SJ
Sample : H2736-02
Misc :
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Instrument :
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ClientSampleId :
RICHMOND-AMBOY-SB02(0.5-8.0)

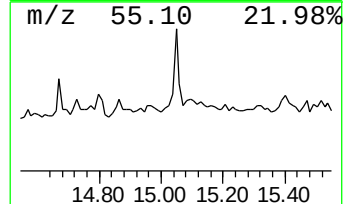
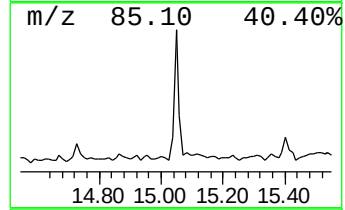
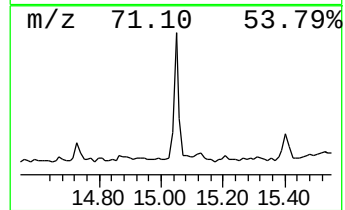
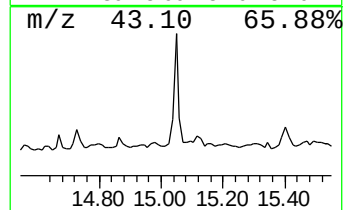
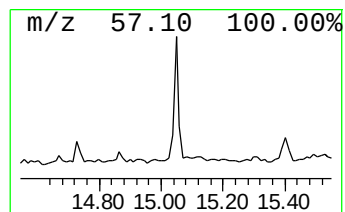
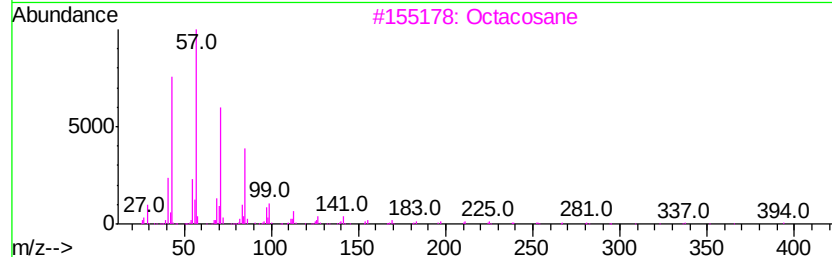
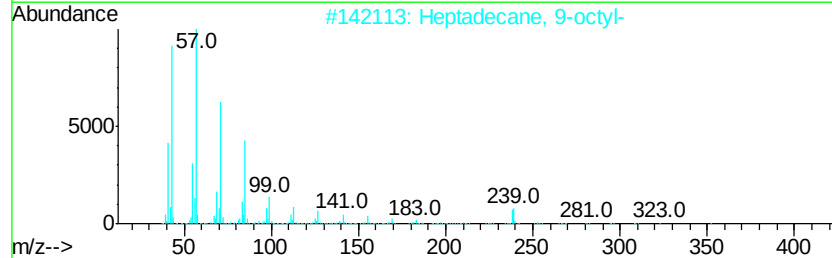
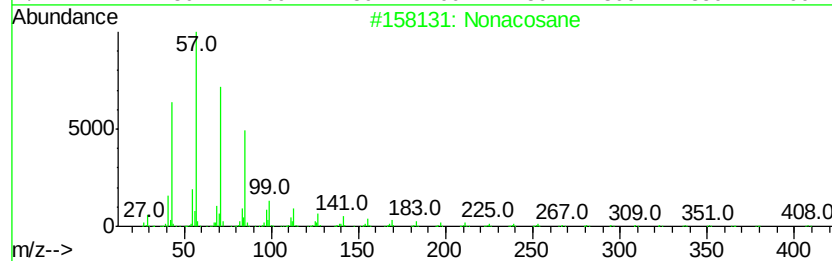
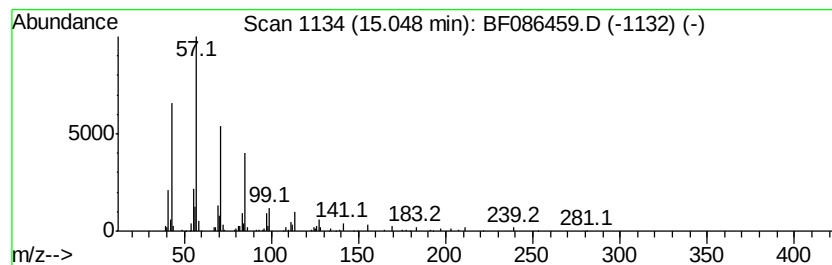
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Nonacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.05	3.76 ng	110552	Perylene-d12	15.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonacosane	408	C29H60	000630-03-5	91
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Hexacosane	366	C26H54	000630-01-3	90
5		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042716\
Data File : BF086459.D
Acq On : 28 Apr 2016 8:05
Operator : UM/SJ
Sample : H2736-02
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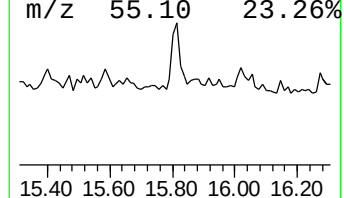
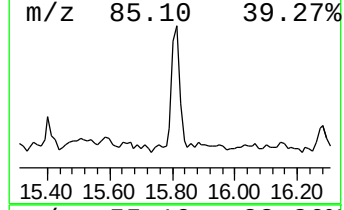
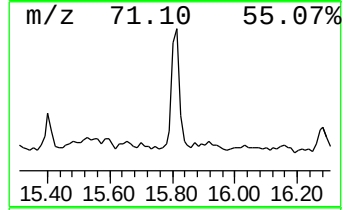
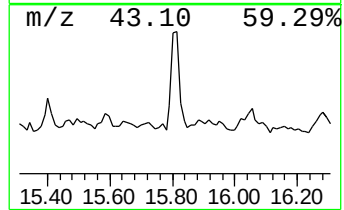
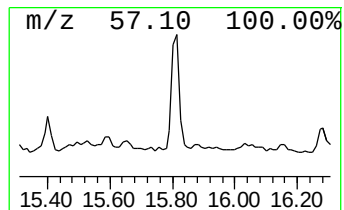
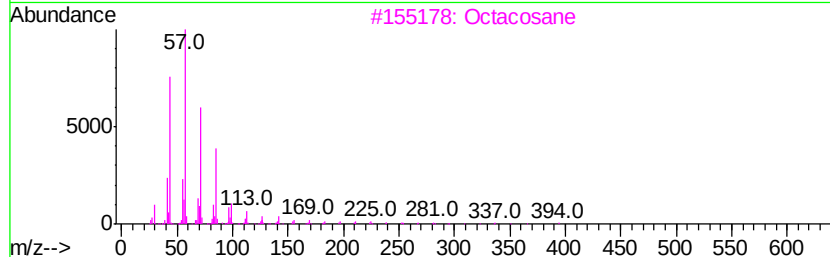
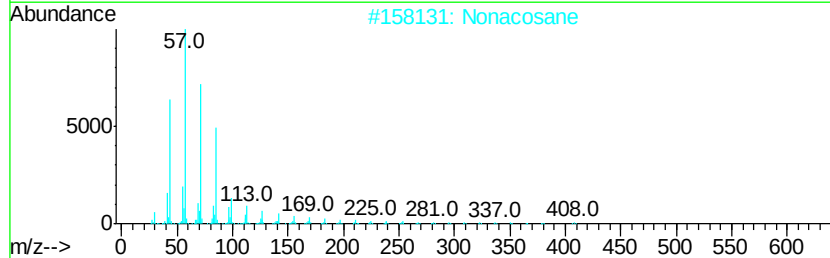
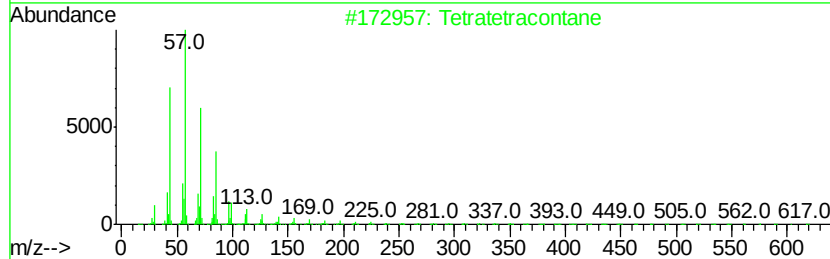
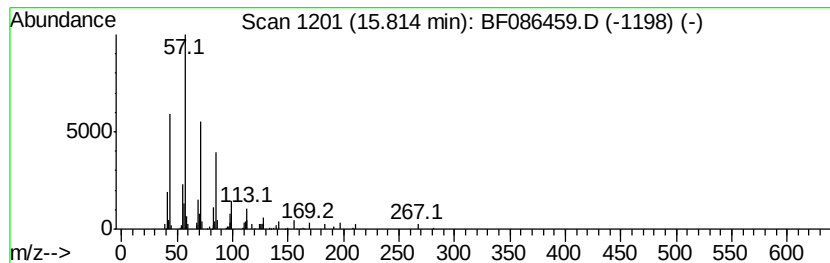
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tetratetracontane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.81	3.61 ng	105916	Perylene-d12	15.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	91
2			Nonacosane	408	C29H60	000630-03-5	91
3			Octacosane	394	C28H58	000630-02-4	90
4			Heneicosane	296	C21H44	000629-94-7	90
5			Eicosane, 3-methyl-	296	C21H44	006418-46-8	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042716\
Data File : BF086459.D
Acq On : 28 Apr 2016 8:05
Operator : UM/SJ
Sample : H2736-02
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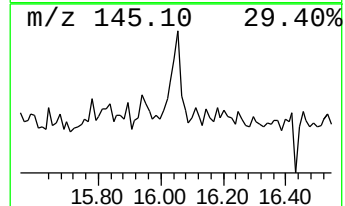
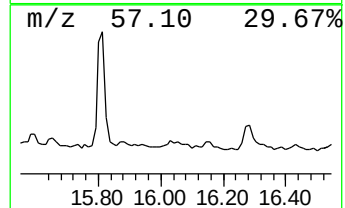
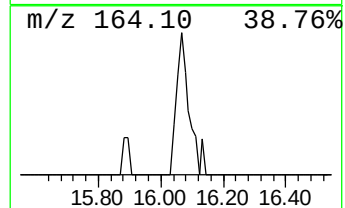
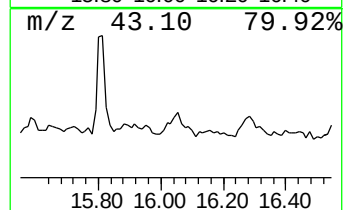
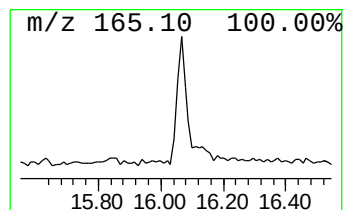
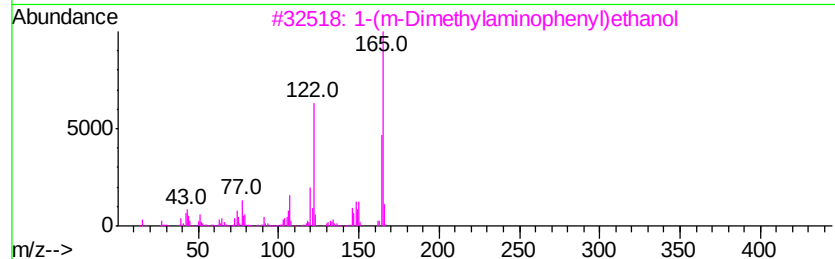
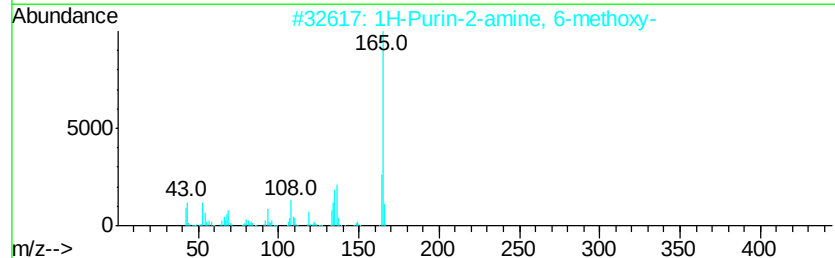
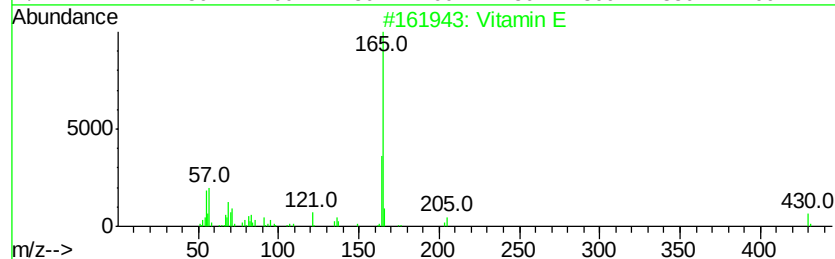
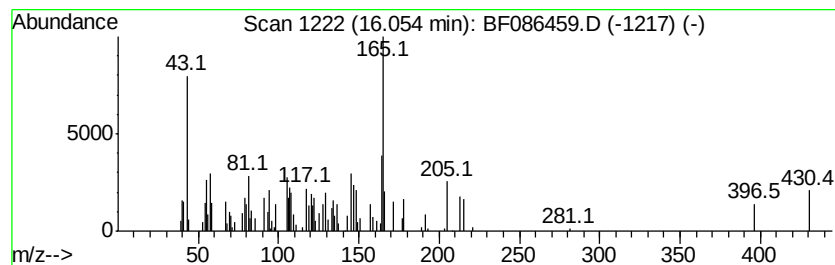
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown16.05 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.05	2.63 ng	77325	Perylene-d12	15.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Vitamin E	430	C29H50O2	000059-02-9	43
2		1H-Purin-2-amine, 6-methoxy-	165	C6H7N5O	020535-83-5	38
3		1-(m-Dimethylaminophenyl)ethanol	165	C10H15NO	005339-01-5	38
4		4-Amino-3-methoxypyrazolo[3,4-d]...	165	C6H7N5O	111375-22-5	35
5		5,6-Dimethoxy-3-pyridazinecarbon...	165	C7H7N3O2	020552-46-9	35



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042716\
Data File : BF086459.D
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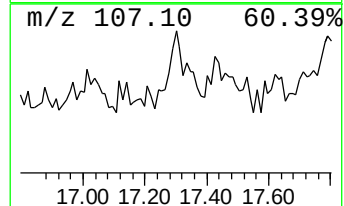
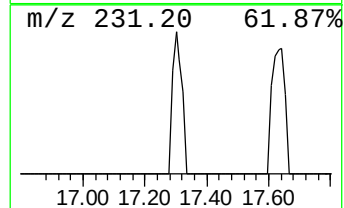
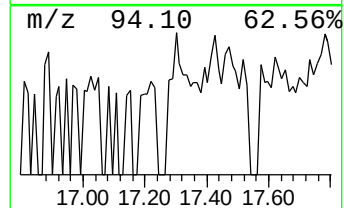
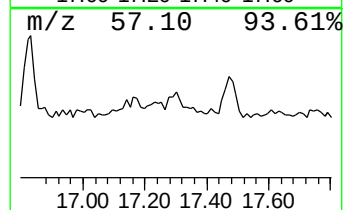
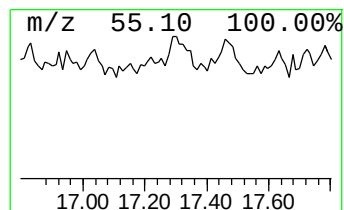
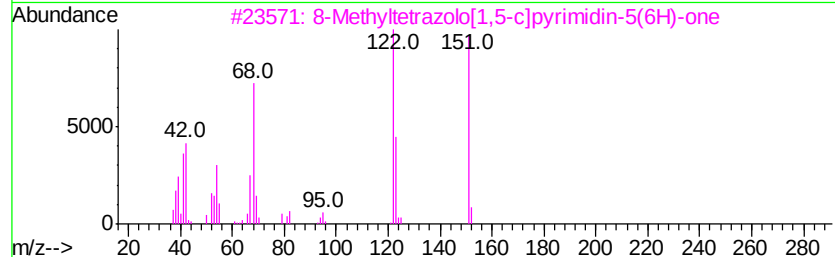
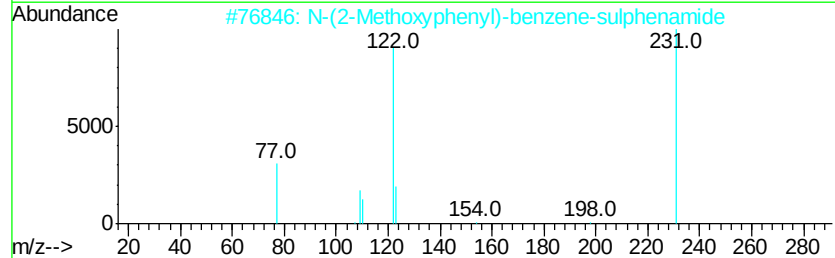
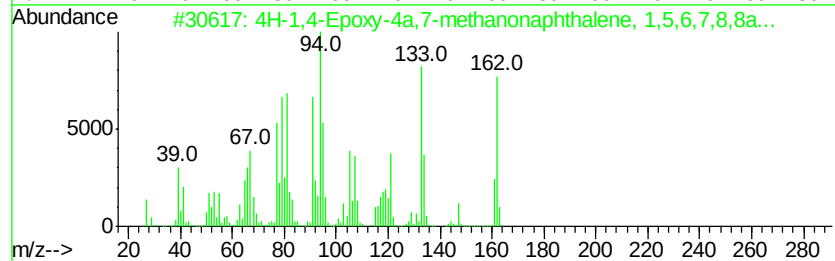
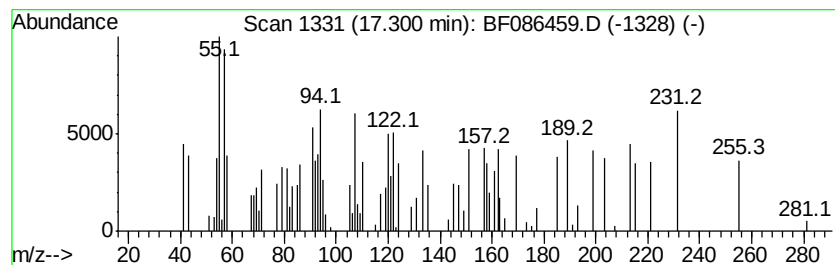
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown17.30 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.30	2.24 ng	65868	Perylene-d12	15.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-1,4-Epoxy-4a,7-methanonaphtha...	162	C11H14O	032763-64-7	15
2		N-(2-Methoxyphenyl)-benzene-sulp...	231	C13H13NOS	111685-76-8	10
3		8-Methyltetrazolo[1,5-c]pyrimidi...	151	C5H5N5O	040595-05-9	7
4		3,4-Pyridinedimethanol, 5-hydrox...	169	C8H11NO3	000065-23-6	7
5		Tetracyclo[5.2.1.0(2,6).0(3,5)]d...	162	C12H18	074646-38-1	7



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042716\
Data File : BF086459.D
Acq On : 28 Apr 2016 8:05
Operator : UM/SJ
Sample : H2736-02
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RICHMOND-AMBOY-SB02(0.5-8.0)

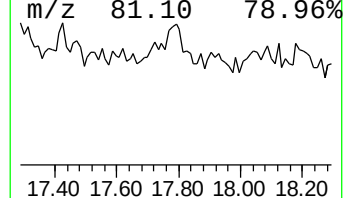
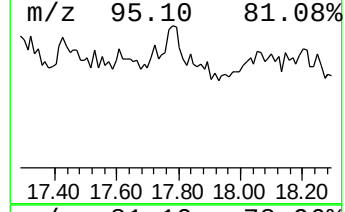
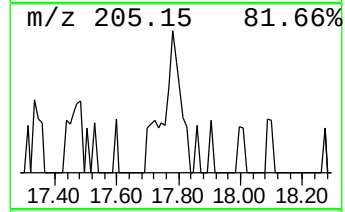
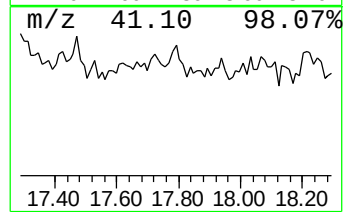
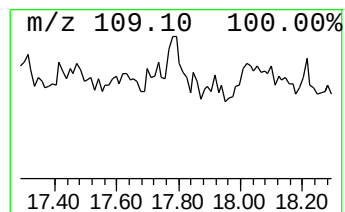
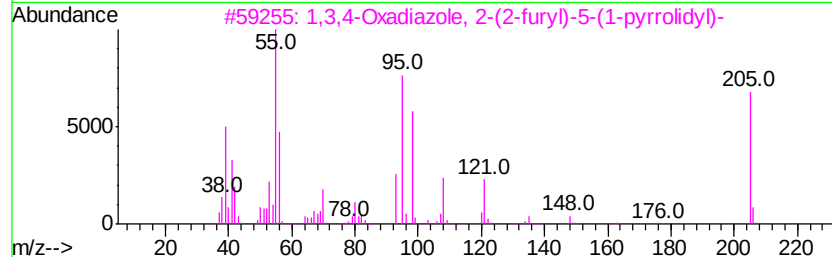
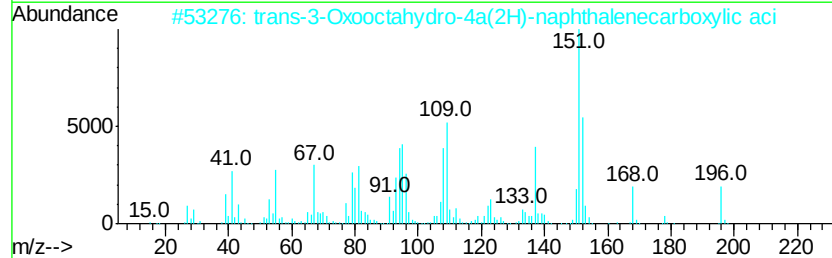
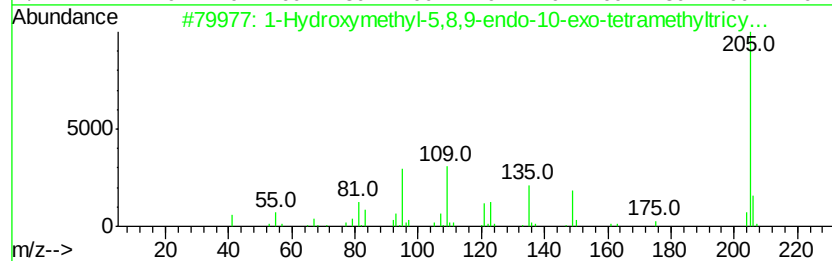
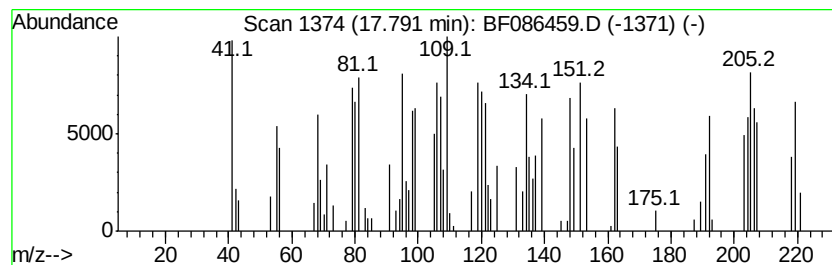
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown17.79 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.79	2.26 ng	66521	Perylene-d12	15.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hydroxymethyl-5,8,9-endo-10-ex...	236	C16H28O	1000140-32-9	27
2		trans-3-Oxoctahydro-4a(2H)-naph...	196	C11H16O3	071198-51-1	12
3		1,3,4-Oxadiazole, 2-(2-furyl)-5-...	205	C10H11N3O2	284674-14-2	11
4		Methoxy(methyl)chlorosilane	110	C2H7ClOSi	018151-27-4	11
5		2,4,4-Trimethyl-3-(3-methylbuta-...	206	C14H22O	097452-05-6	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042716\
Data File : BF086459.D
Acq On : 28 Apr 2016 8:05
Operator : UM/SJ
Sample : H2736-02
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RICHMOND-AMBOY-SB02(0.5-8.0)

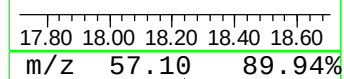
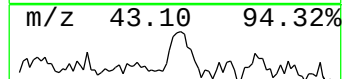
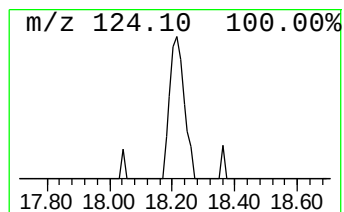
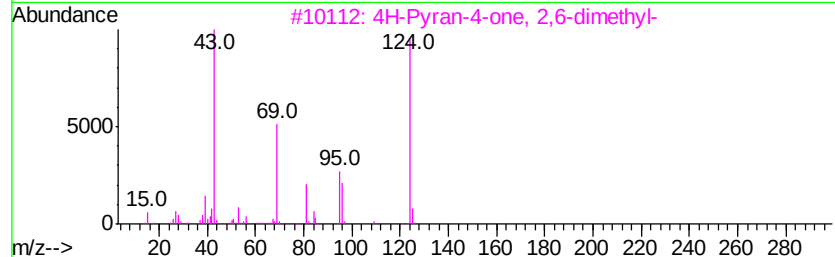
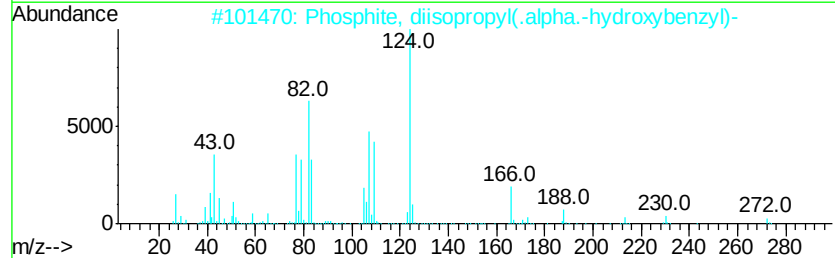
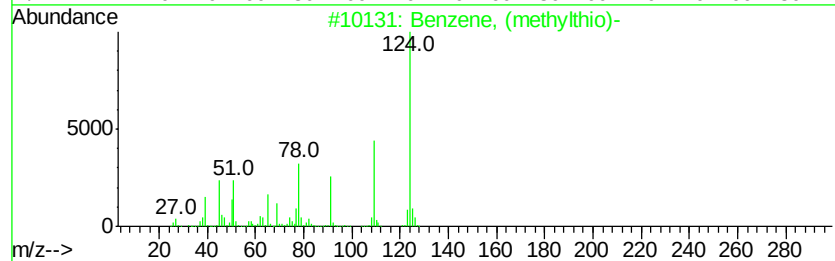
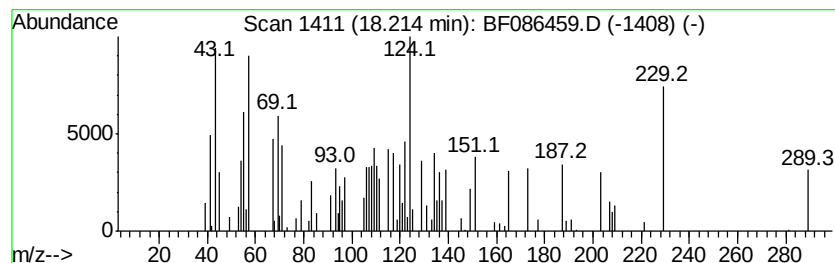
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown18.21 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.21	2.40 ng	70448	Perylene-d12	15.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (methylthio)-	124	C7H8S	000100-68-5	14
2			Phosphite, diisopropyl(.alpha.-h...	272	C13H21O4P	020386-43-0	14
3			4H-Pyran-4-one, 2,6-dimethyl-	124	C7H8O2	001004-36-0	11
4			4,5-Pyrimidinediamine, 6-methyl-	124	C5H8N4	022715-28-2	10
5			1H-1,2,3-Triazol-1-amine, N-[(4-...	234	C11H11ClN4	113261-29-3	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042716\
 Data File : BF086459.D
 Acq On : 28 Apr 2016 8:05
 Operator : UM/SJ
 Sample : H2736-02
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RICHMOND-AMBOY-SB02(0.5-8.0)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.99	2.5	ng	94564	1	6.81	768653	20.0
unknown6.57	6.57	84.0	ng	3230320	1	6.81	768653	20.0
Hexanoic acid, 2-...	7.47	2.0	ng	90196	2	8.09	901010	20.0
Hexadecane	10.28	2.5	ng	96439	3	9.85	766912	20.0
Eicosane	11.20	2.1	ng	71689	4	11.32	671024	20.0
1-Hexadecanol	11.56	5.3	ng	178361	4	11.32	671024	20.0
1-Docosene	12.39	6.2	ng	208679	4	11.32	671024	20.0
17-Pentatriacontene	13.81	5.1	ng	228546	5	13.95	894742	20.0
Nonacosane	15.05	3.8	ng	110552	6	15.37	587392	20.0
Tetratetracontane	15.81	3.6	ng	105916	6	15.37	587392	20.0
unknown16.05	16.05	2.6	ng	77325	6	15.37	587392	20.0
unknown17.30	17.30	2.2	ng	65868	6	15.37	587392	20.0
unknown17.79	17.79	2.3	ng	66521	6	15.37	587392	20.0
unknown18.21	18.21	2.4	ng	70448	6	15.37	587392	20.0