

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050515\
 Data File : BF078946.D
 Acq On : 5 May 2015 17:56
 Operator : TP/IZ
 Sample : G2032-06
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B3(5-7)

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-BF043015.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	1.518	26	29	31	rVB	4891470	6021795	100.00%	20.487%
2	1.655	38	41	44	rVB	186060	330627	5.49%	1.125%
3	1.770	50	51	55	rBV	277728	382189	6.35%	1.300%
4	1.838	55	57	61	rVV	226306	295090	4.90%	1.004%
5	1.907	61	63	84	rVV	1315885	2963192	49.21%	10.081%
6	2.158	84	85	93	rVB	35425	69729	1.16%	0.237%
7	5.164	346	348	355	rVB	184782	250253	4.16%	0.851%
8	5.656	388	391	393	rBV	1434260	1654931	27.48%	5.630%
9	6.902	498	500	503	rBV	1749467	1624773	26.98%	5.528%
10	7.062	512	514	517	rBV	1751063	1747588	29.02%	5.946%
11	7.359	537	540	542	rBV	444733	431944	7.17%	1.470%
12	7.542	553	556	559	rVB	1462032	1294872	21.50%	4.405%
13	8.045	597	600	602	rBV	1262945	1087470	18.06%	3.700%
14	8.582	643	647	649	rBV	60720	104021	1.73%	0.354%
15	8.936	675	678	680	rBV	491541	583545	9.69%	1.985%
16	10.273	793	795	798	rBV	1575507	1690629	28.08%	5.752%
17	10.776	837	839	842	rBV	73558	75617	1.26%	0.257%
18	11.108	866	868	871	rBV	724995	675043	11.21%	2.297%
19	11.451	895	898	901	rBV	65365	81895	1.36%	0.279%
20	12.091	951	954	956	rBV	1115781	1282313	21.29%	4.363%
21	12.959	1027	1030	1031	rBV	500701	697858	11.59%	2.374%
22	13.668	1090	1092	1102	rBV	175237	246447	4.09%	0.838%
23	13.851	1106	1108	1115	rVB	55781	79147	1.31%	0.269%
24	14.537	1166	1168	1172	rVB3	227857	347740	5.77%	1.183%
25	14.651	1176	1178	1184	rBV5	30567	68146	1.13%	0.232%
26	14.811	1190	1192	1195	rBV	632279	725686	12.05%	2.469%
27	14.868	1195	1197	1199	rVV	51423	66756	1.11%	0.227%
28	14.948	1201	1204	1207	rVV	2090788	2084982	34.62%	7.094%
29	16.125	1304	1307	1315	rBV	131087	260152	4.32%	0.885%
30	16.251	1315	1318	1320	rBV	690859	847865	14.08%	2.885%
31	16.937	1376	1378	1383	rBV2	34663	65028	1.08%	0.221%
32	17.726	1445	1447	1451	rBV2	29957	61493	1.02%	0.209%
33	18.000	1468	1471	1474	rBV	656985	840146	13.95%	2.858%
34	18.754	1535	1537	1545	rVB3	47341	114828	1.91%	0.391%

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ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
B3(5-7)

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-BF043015.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	19.200	1573	1576	1580	rVB2	57018	103001	1.71%	0.350%
36	20.160	1656	1660	1666	rVB5	48632	135789	2.25%	0.462%

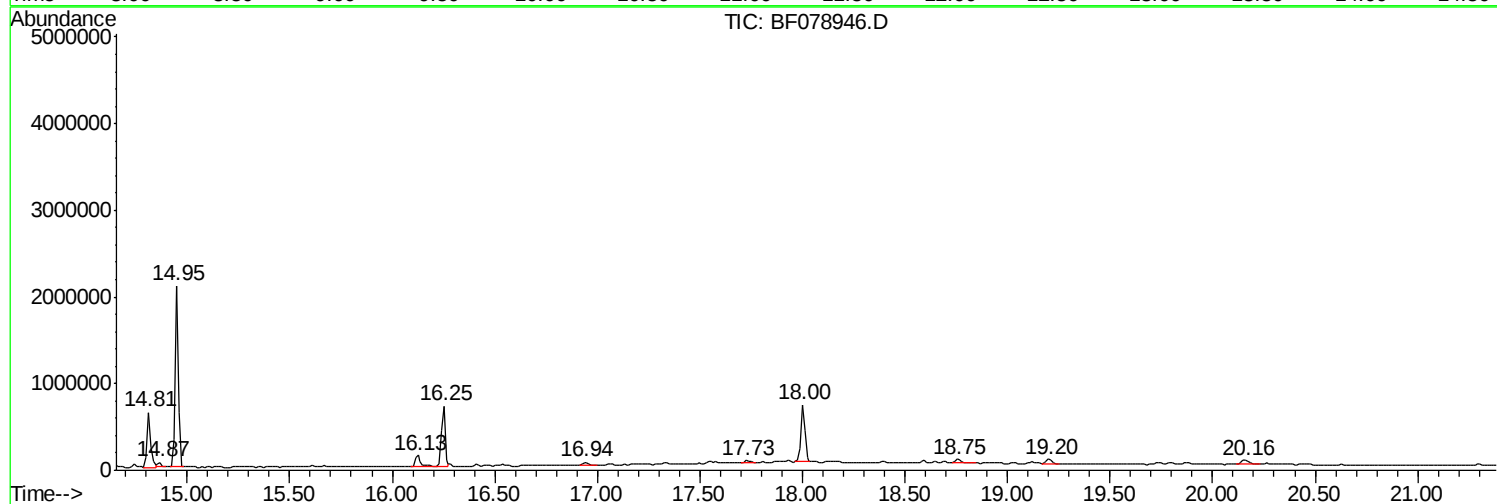
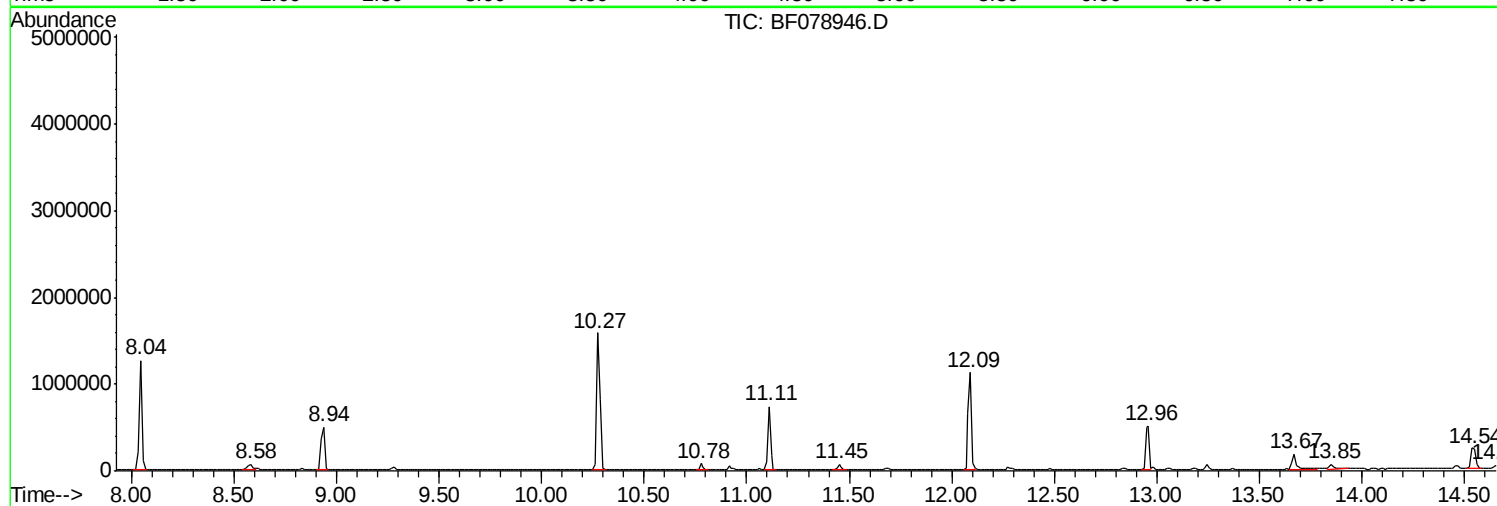
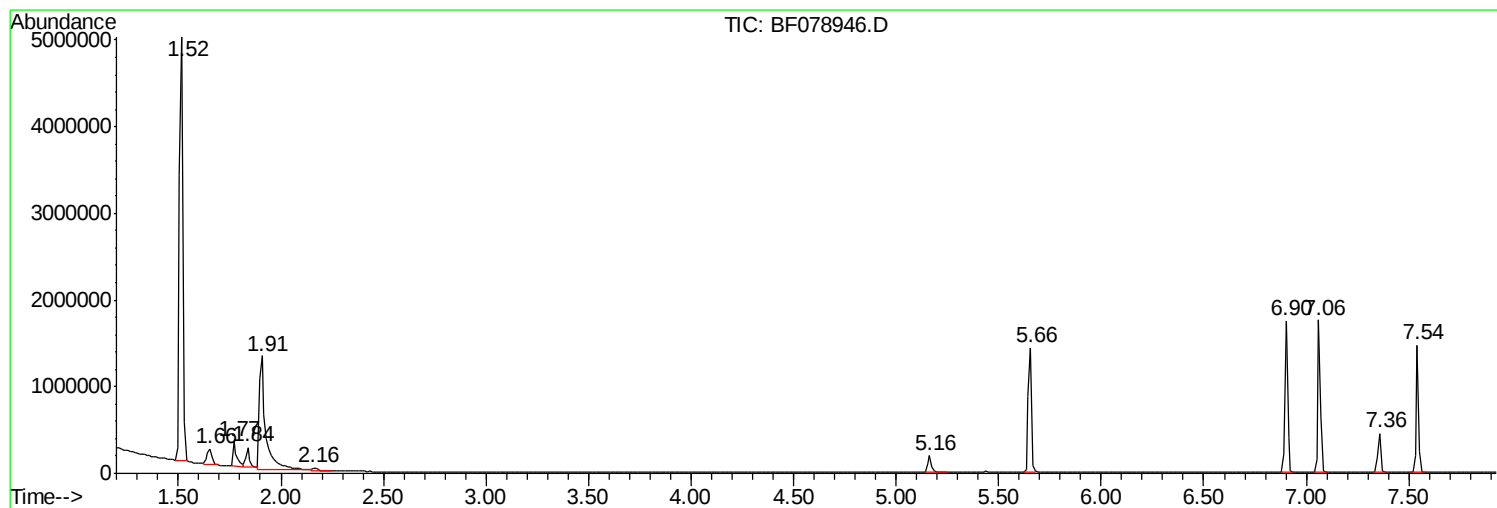
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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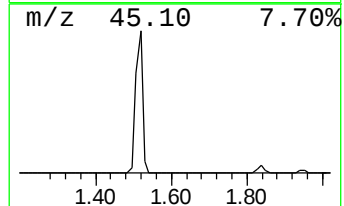
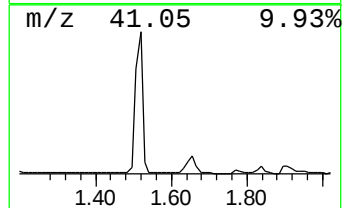
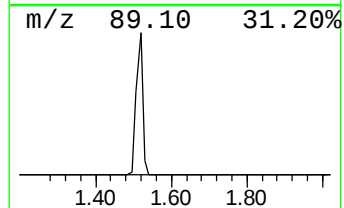
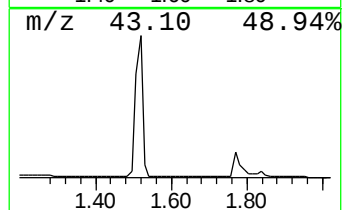
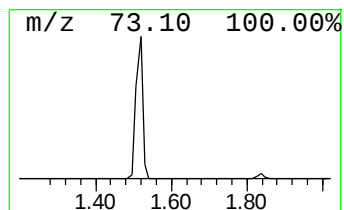
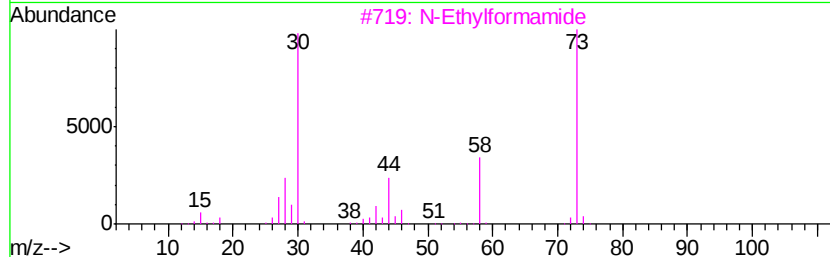
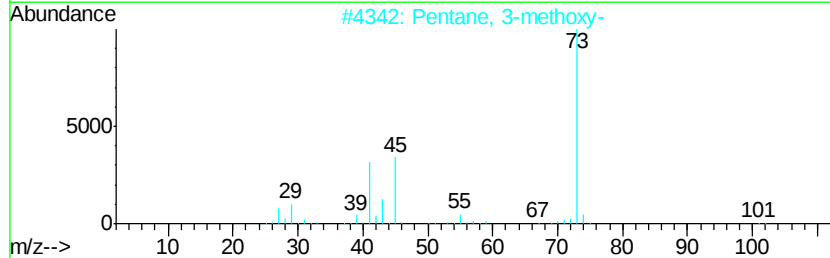
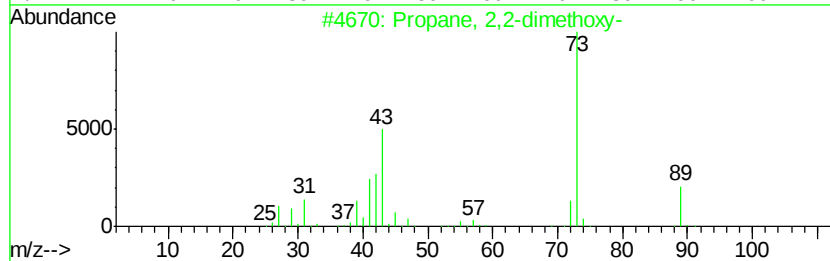
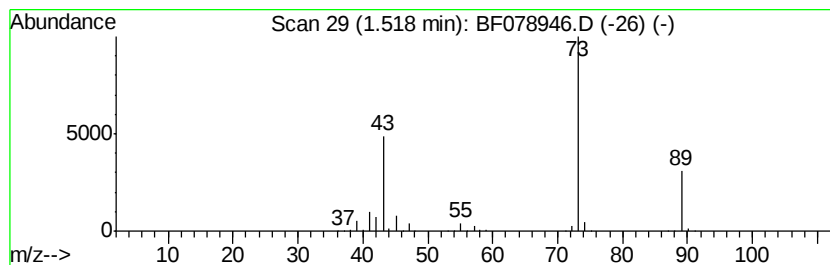
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown1.52 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.52	278.82 ng	6021800	1,4-Dichlorobenzene-d4	7.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	38
2		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
3		N-Ethylformamide	73	C3H7NO	000627-45-2	9
4		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
5		2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9



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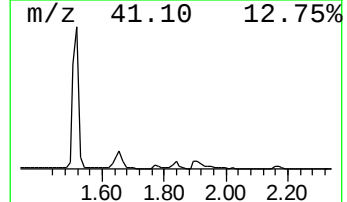
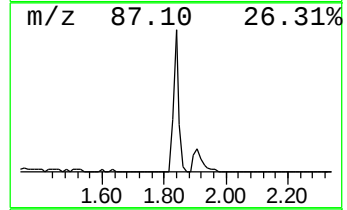
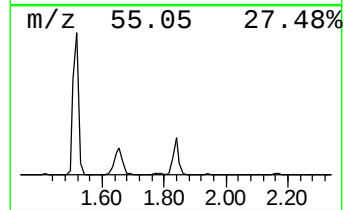
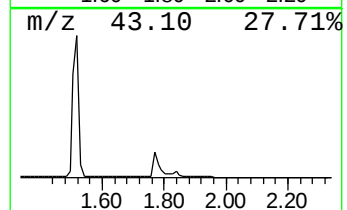
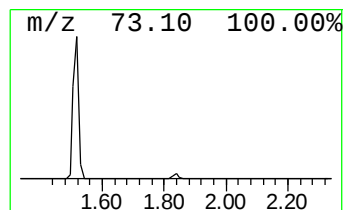
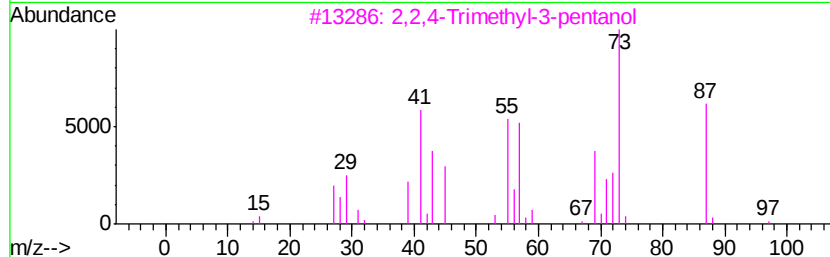
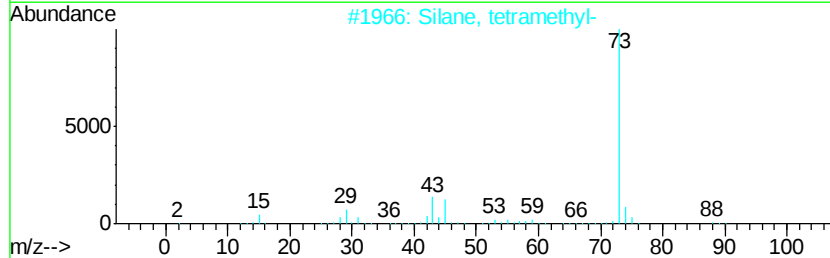
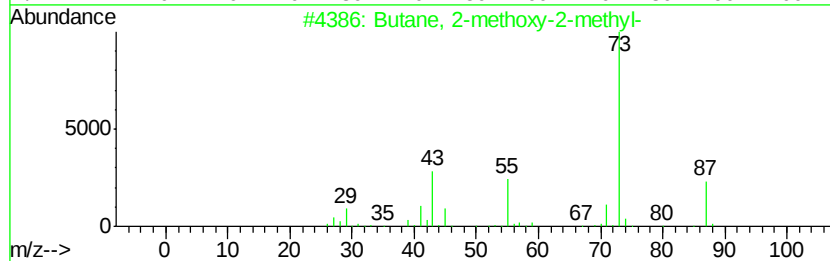
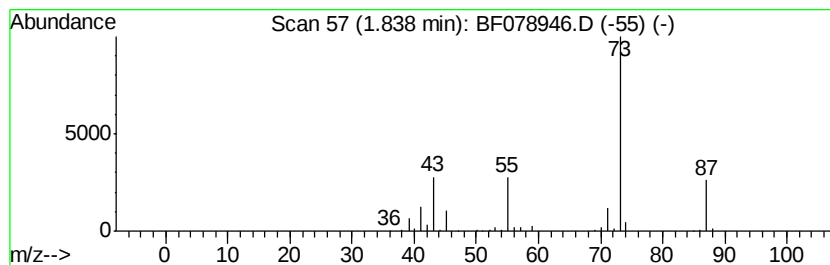
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.84	13.66 ng	295090	1,4-Dichlorobenzene-d4	7.36

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	86
2		Silane, tetramethyl-	88	C4H12Si	000075-76-3	43
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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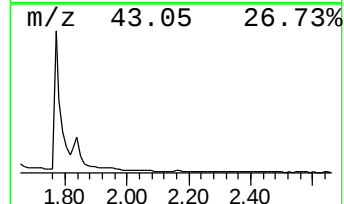
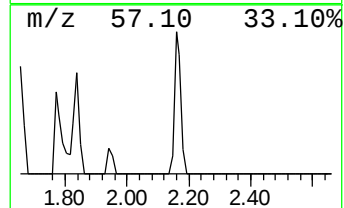
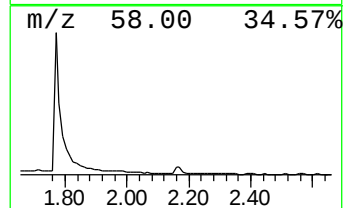
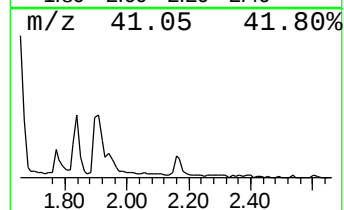
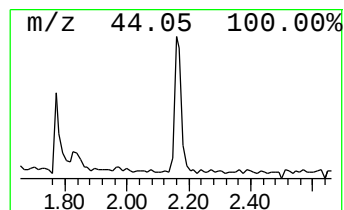
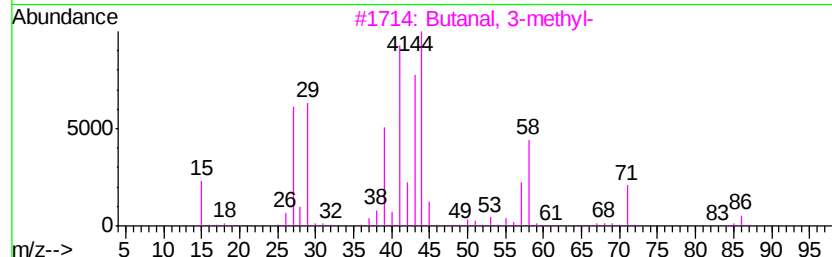
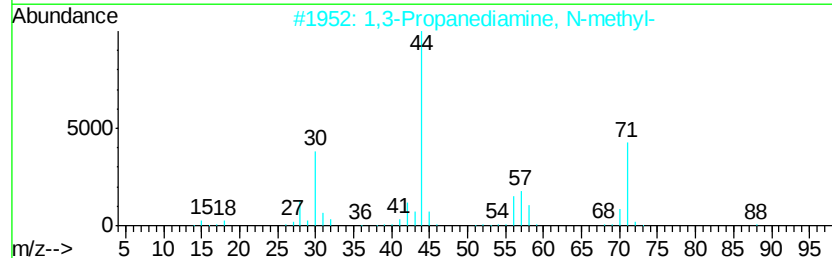
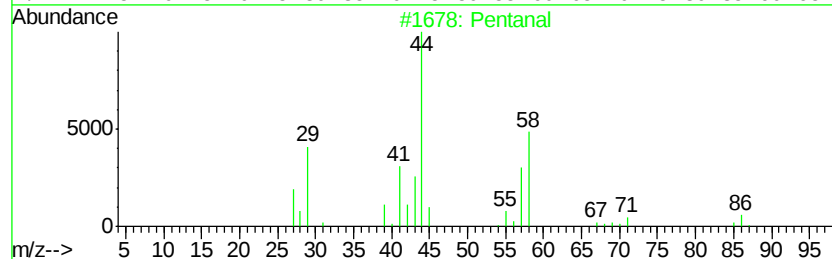
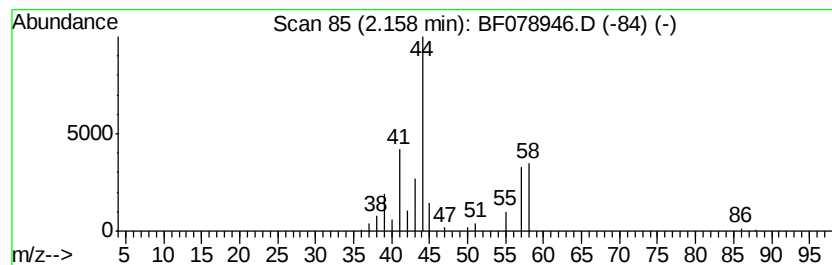
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown2.16 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.16	3.23 ng	69729	1,4-Dichlorobenzene-d4	7.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentanal	86	C5H10O	000110-62-3	39
2		1,3-Propanediamine, N-methyl-	88	C4H12N2	006291-84-5	9
3		Butanal, 3-methyl-	86	C5H10O	000590-86-3	7
4		Ethene, methoxy-	58	C3H6O	000107-25-5	5
5		Butanal, 3-hydroxy-	88	C4H8O2	000107-89-1	5



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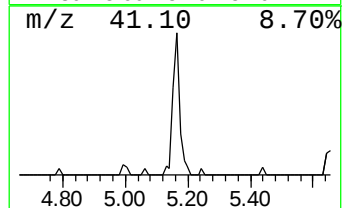
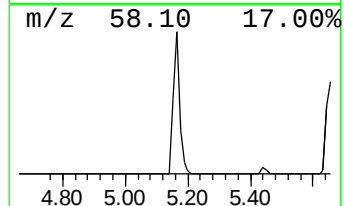
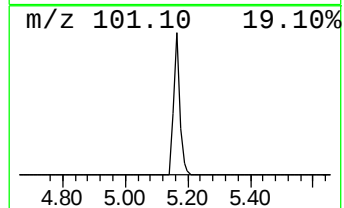
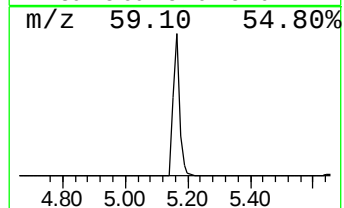
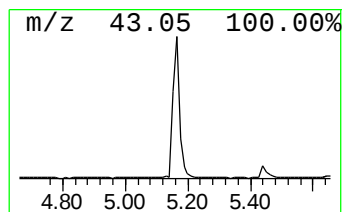
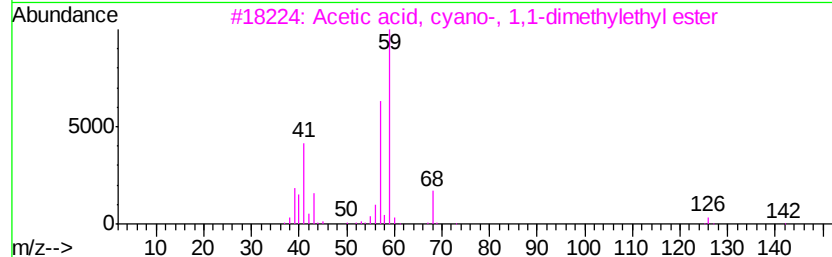
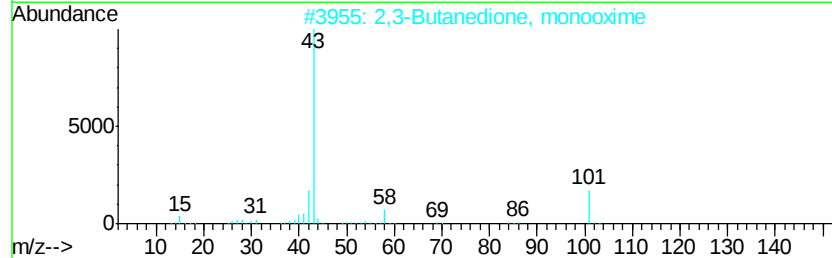
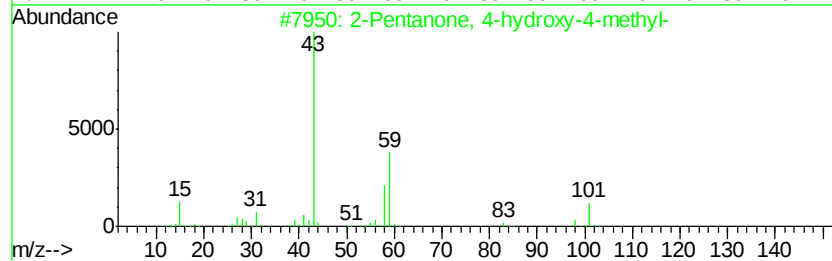
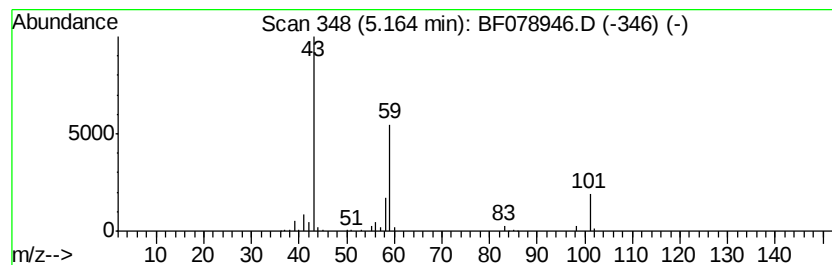
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	11.59 ng	250253	1,4-Dichlorobenzene-d4	7.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25		
3	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17		
4	5-Hexen-2-one	98	C6H10O	000109-49-9	9		
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9		



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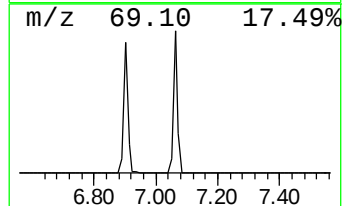
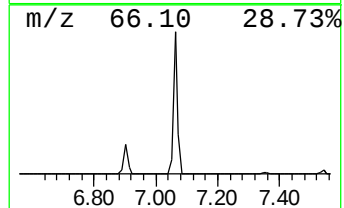
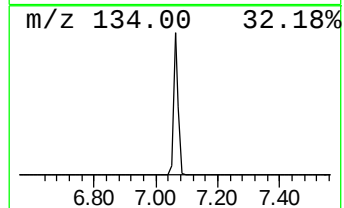
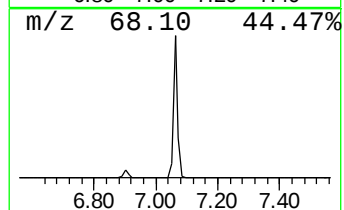
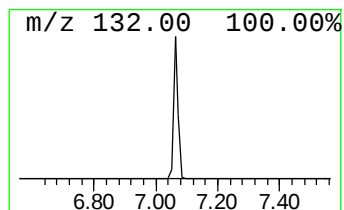
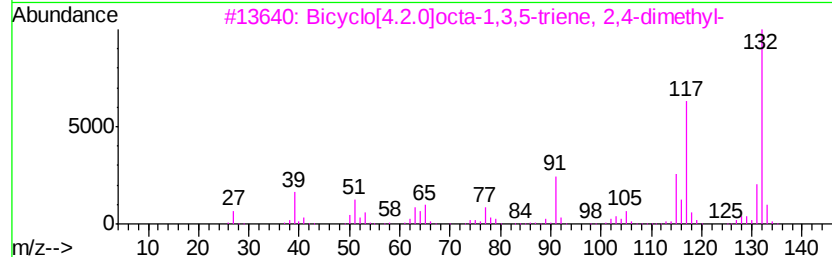
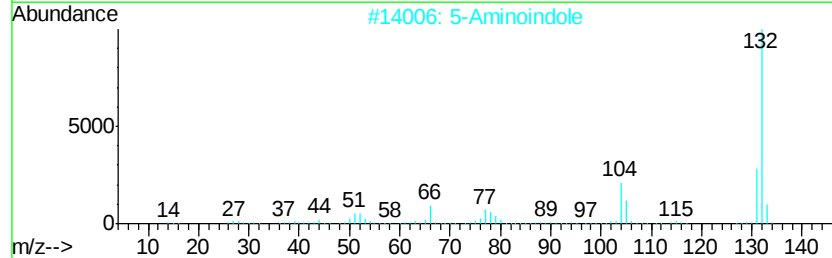
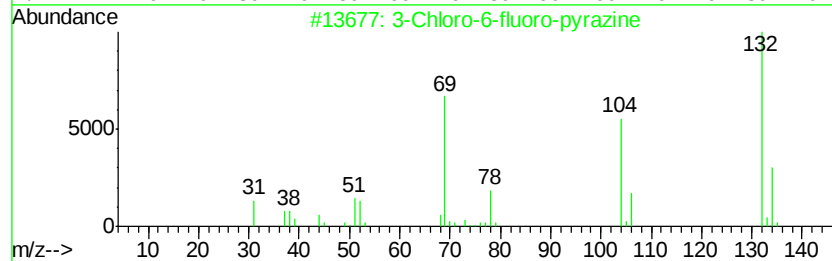
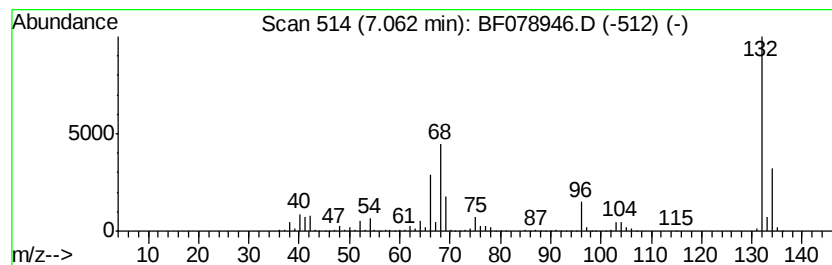
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 8 unknown7.06 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.06	80.92 ng	1747590	1,4-Dichlorobenzene-d4	7.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Aminoindole	132	C8H8N2	005192-03-0	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050515\
Data File : BF078946.D
Acq On : 5 May 2015 17:56
Operator : TP/IZ
Sample : G2032-06
Misc :
ALS Vial : 18 Sample Multiplier: 1

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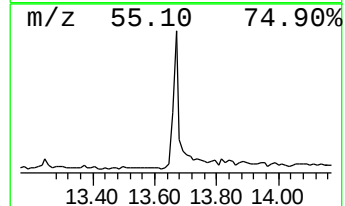
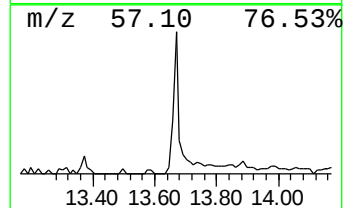
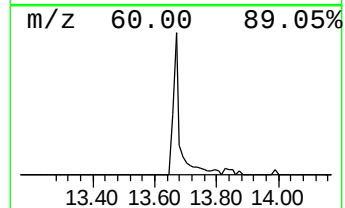
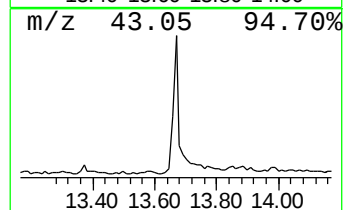
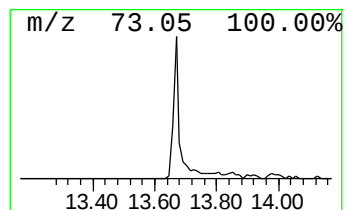
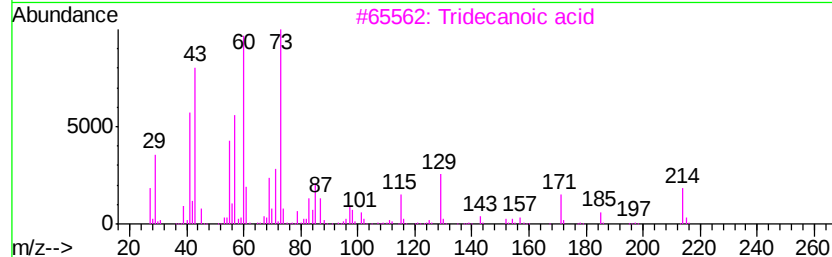
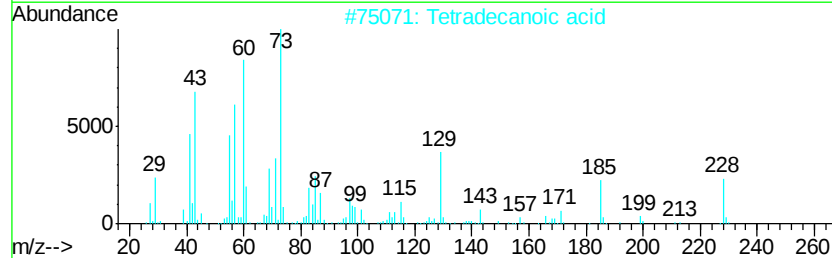
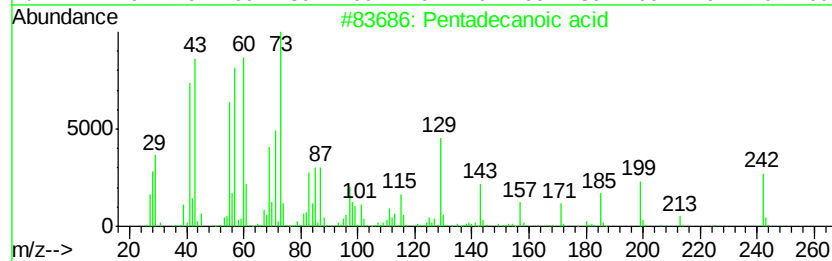
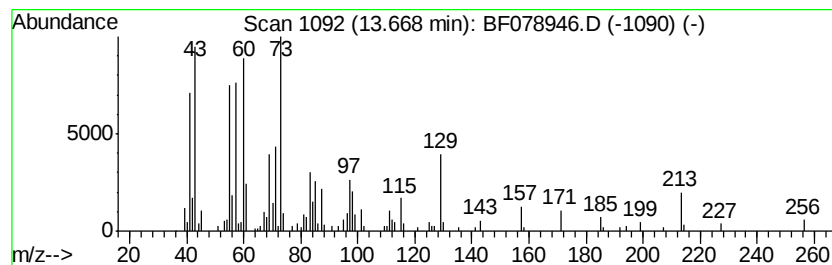
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Pentadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.67	7.06 ng	246447	Phenanthrene-d10	12.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecanoic acid	242	C15H30O2	001002-84-2	74
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	72
3			Tridecanoic acid	214	C13H26O2	000638-53-9	68
4			5-Acetoxypentadecane	270	C17H34O2	1000245-62-3	14
5			Tetradecane	198	C14H30	000629-59-4	11



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050515\
Data File : BF078946.D
Acq On : 5 May 2015 17:56
Operator : TP/IZ
Sample : G2032-06
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B3(5-7)

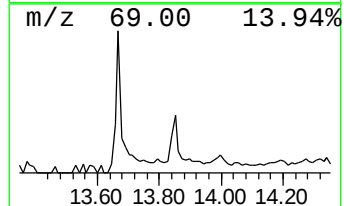
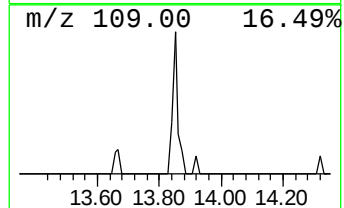
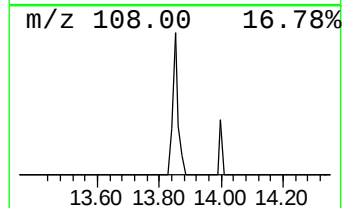
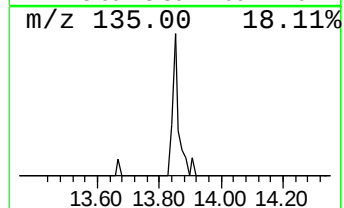
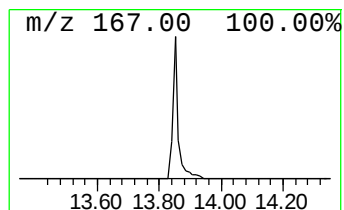
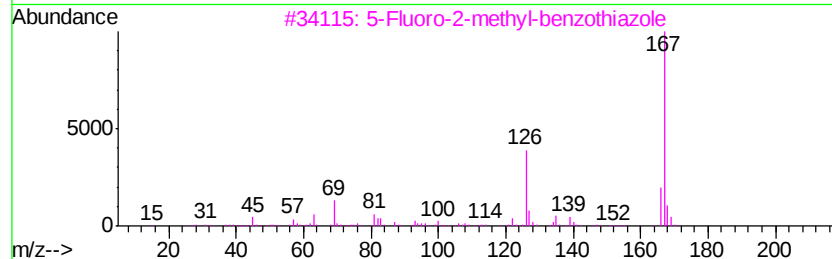
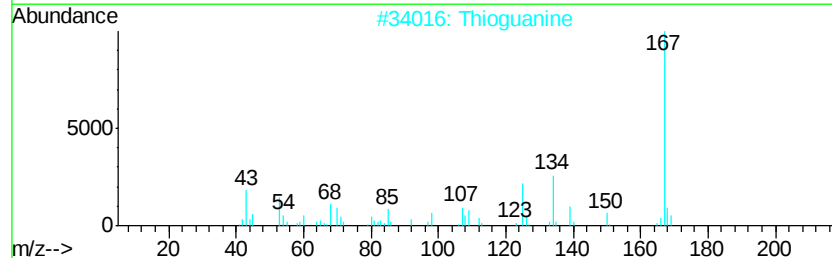
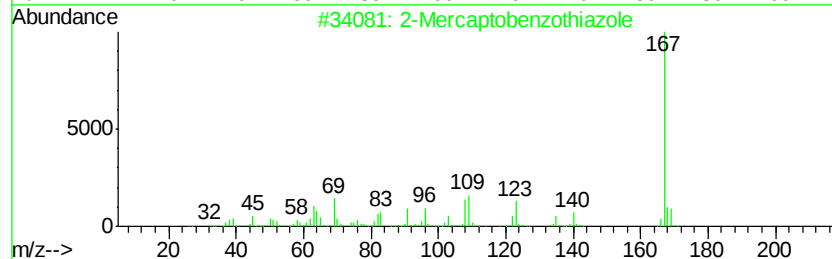
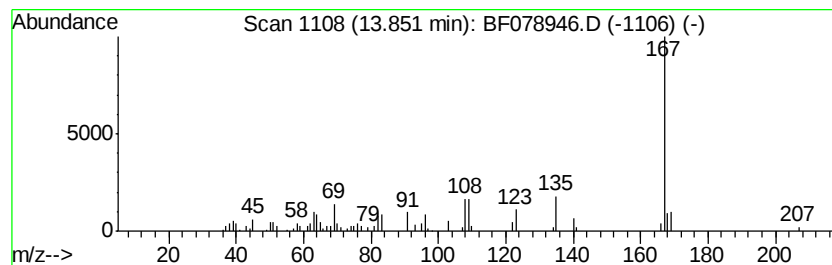
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2-Mercaptobenzothiazole Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.85	2.27 ng	79147	Phenanthrene-d10	12.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Mercaptobenzothiazole	167	C7H5NS2	000149-30-4	96
2			Thioquanine	167	C5H5N5S	000154-42-7	59
3			5-Fluoro-2-methyl-benzothiazole	167	C8H6FNS	000399-75-7	53
4			3-Methoxythiobenzamide	167	C8H9NOS	064559-06-4	53
5			5-Fluoro-2-methylphenyl isothioc...	167	C8H6FNS	175205-39-7	40



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050515\
Data File : BF078946.D
Acq On : 5 May 2015 17:56
Operator : TP/IZ
Sample : G2032-06
Misc :
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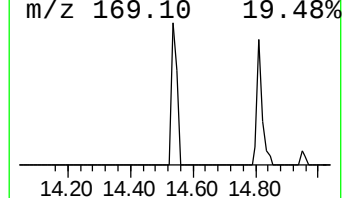
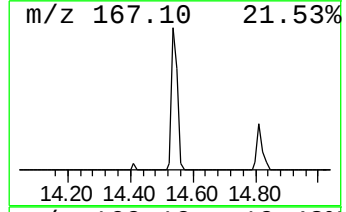
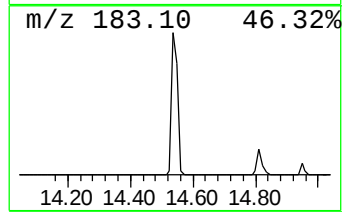
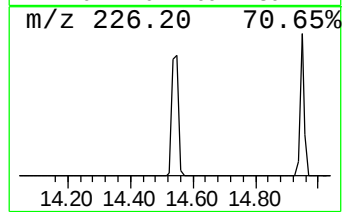
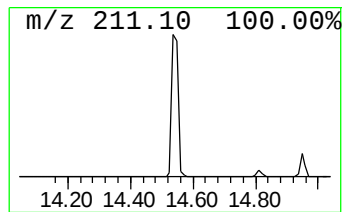
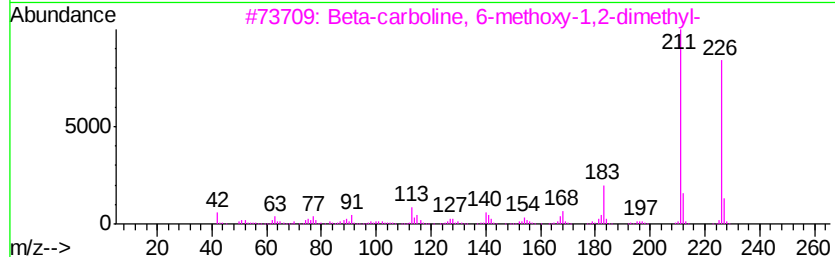
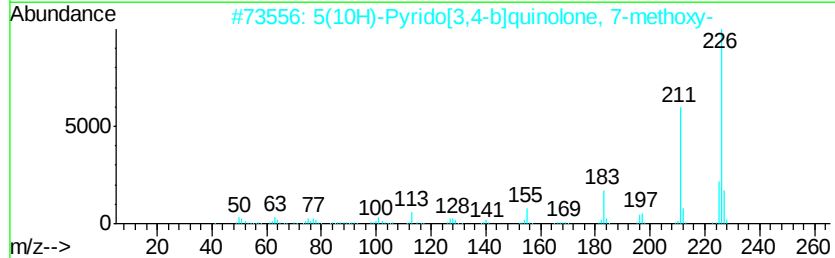
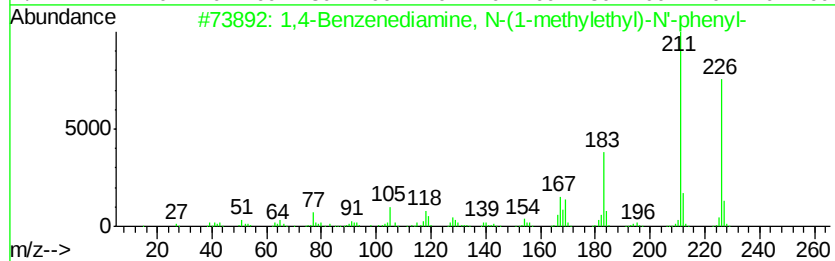
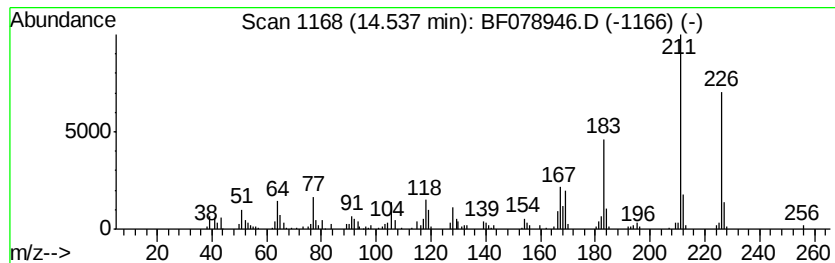
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1,4-Benzenediamine, N-(1-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.54	9.97 ng	347740	Phenanthrene-d10	12.96		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Benzenediamine, N-(1-methyle...	226	C15H18N2	000101-72-4	99	
2	5(10H)-Pyrido[3,4-b]quinolone, 7...	226	C13H10N2O2	1000256-99-5	81	
3	Beta-carboline, 6-methoxy-1,2-di...	226	C14H14N2O	154505-42-7	58	
4	1,1'-Biphenyl, 3-(1,1-dimethylet...	226	C16H18O	055059-21-7	49	
5	9H-Purine, 6-chloro-9-(trimethyl...	226	C8H11ClN4Si	032865-86-4	49	



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050515\
Data File : BF078946.D
Acq On : 5 May 2015 17:56
Operator : TP/IZ
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Misc :
ALS Vial : 18 Sample Multiplier: 1

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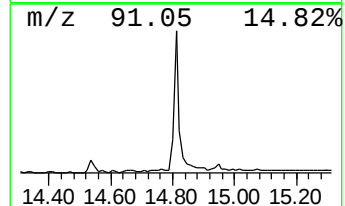
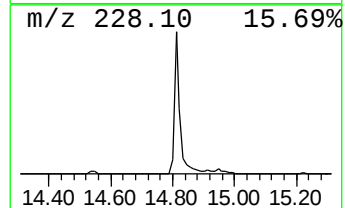
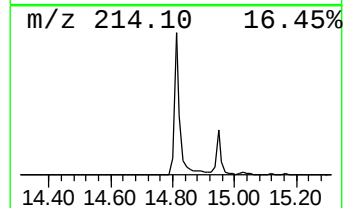
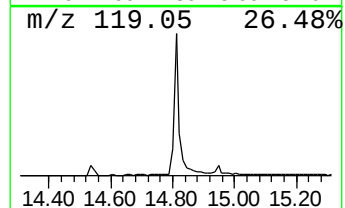
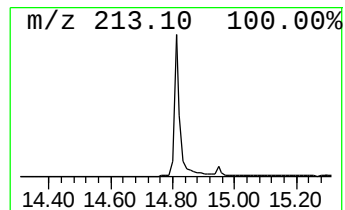
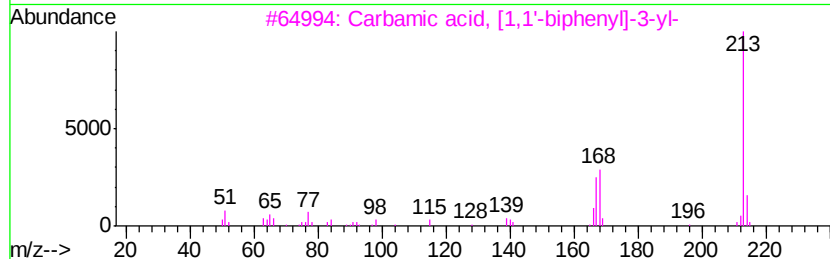
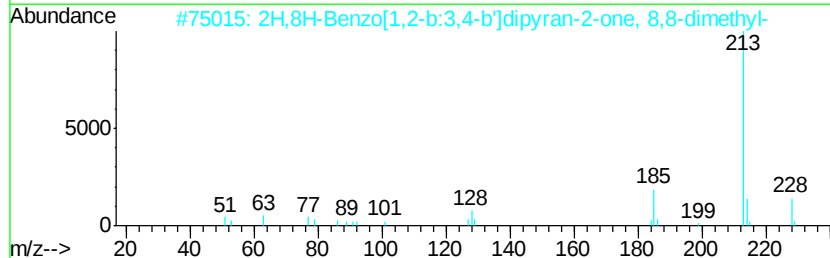
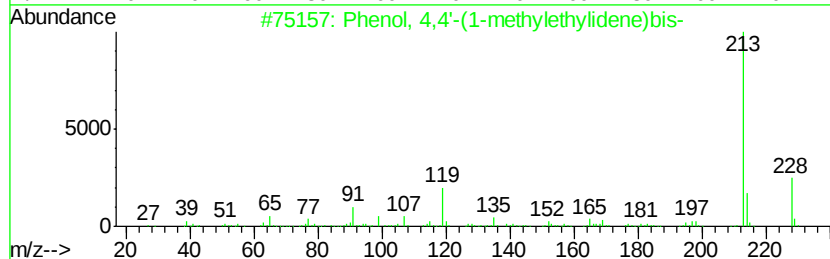
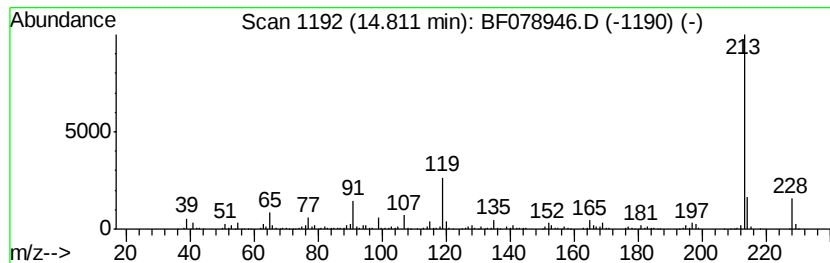
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Phenol, 4,4'-(1-methylethyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.81	17.12 ng	725686	Chrysene-d12	16.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	97
2		2H,8H-Benzo[1,2-b:3,4-b']dipyrans...	228	C14H12O3	000523-59-1	53
3		Carbamic acid, [1,1'-biphenyl]-3...	213	C13H11NO2	055030-29-0	49
4		Carbanilic acid, p-phenyl-	213	C13H11NO2	004474-53-7	49
5		1,4-Puran, 5,6-dihydro-3-[3-indo...	213	C14H15NO	1000128-72-9	40



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050515\
Data File : BF078946.D
Acq On : 5 May 2015 17:56
Operator : TP/IZ
Sample : G2032-06
Misc :
ALS Vial : 18 Sample Multiplier: 1

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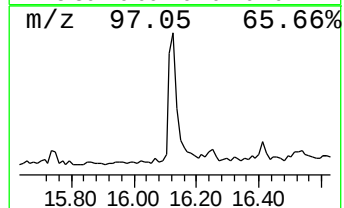
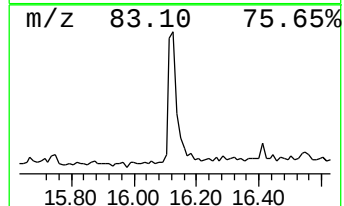
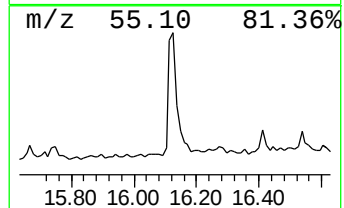
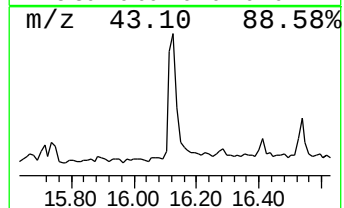
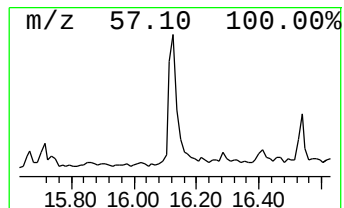
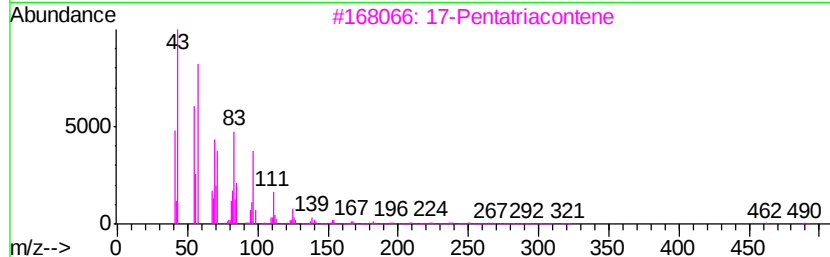
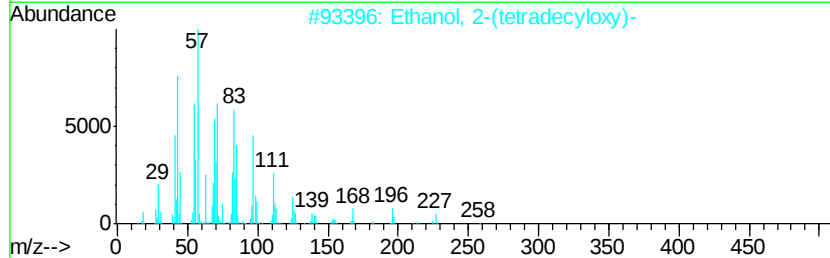
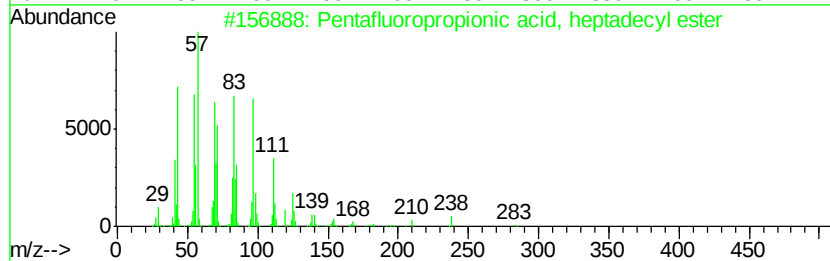
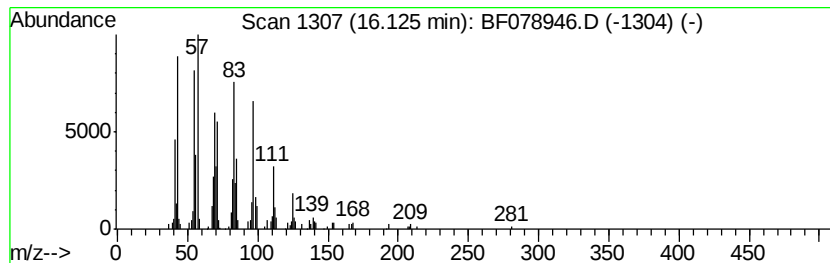
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Pentafluoropropionic acid, ... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.13	6.14 ng	260152	Chrysene-d12	16.25

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	91
2			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
3			17-Pentatriacontene	491	C35H70	006971-40-0	91
4			2- Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	90
5			2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050515\
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Acq On : 5 May 2015 17:56
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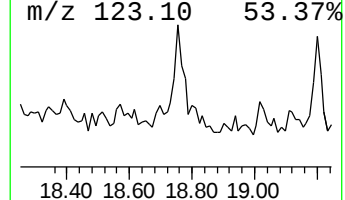
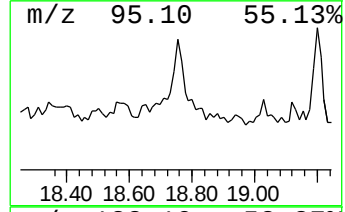
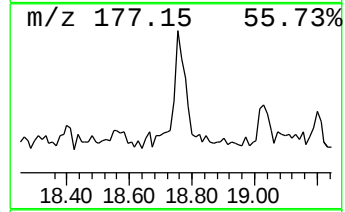
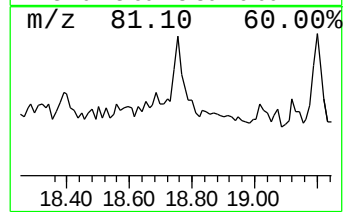
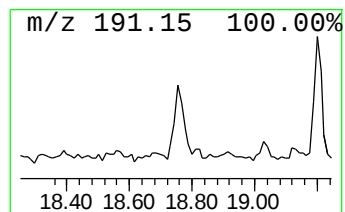
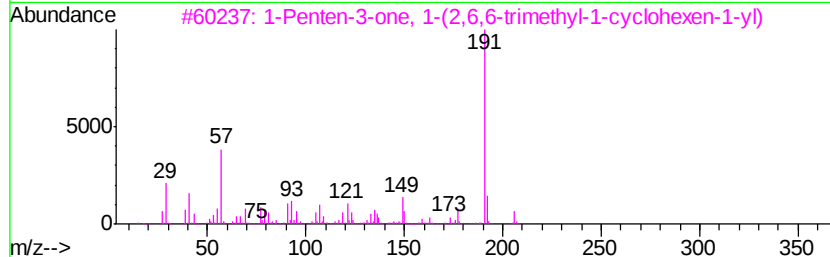
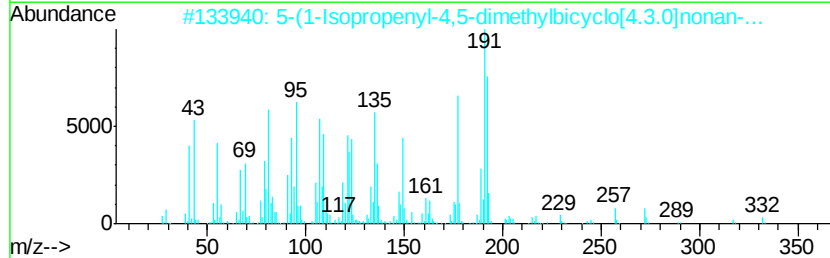
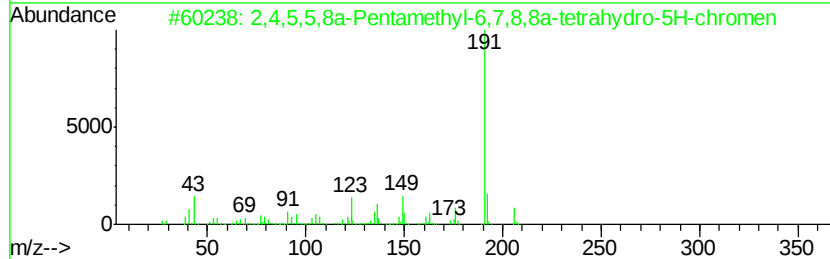
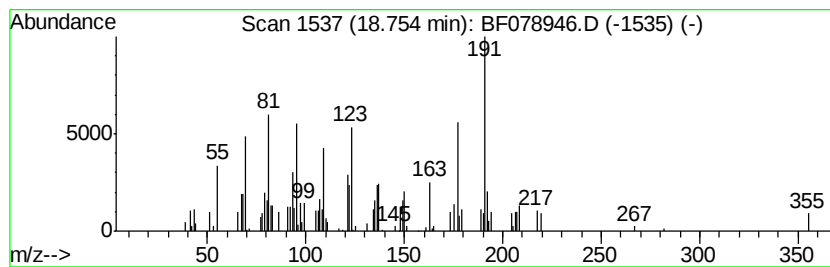
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown18.75 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.75	2.73 ng	114828	Perylene-d12	18.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	43
2		5-(1-Isopropenyl-4,5-dimethylbic...	332	C22H36O2	1000195-69-8	38
3		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	27
4		3,7-Diisopropyl-5-methyl-1,2(4H)...	192	C12H20N2	088829-71-4	25
5		2-Propenamide, N-(4-methoxyphenyl)-	177	C10H11NO2	007766-37-2	18



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050515\
Data File : BF078946.D
Acq On : 5 May 2015 17:56
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ClientSampleId :
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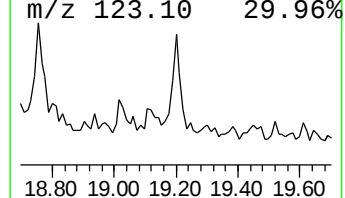
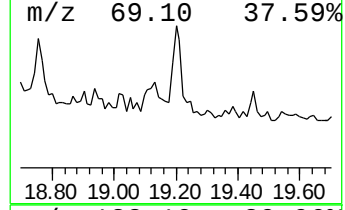
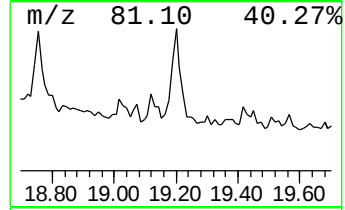
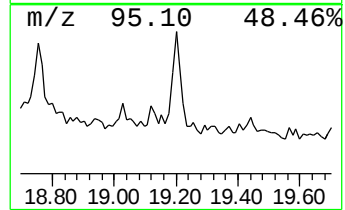
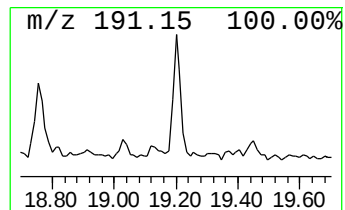
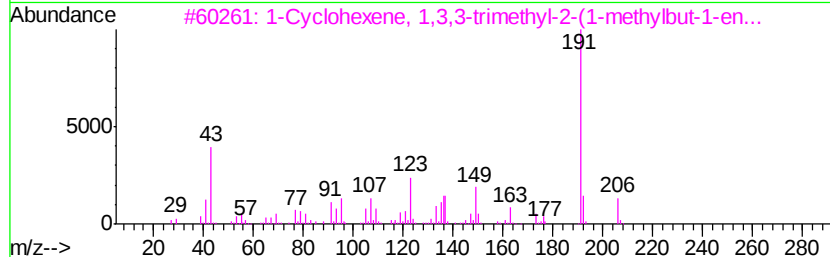
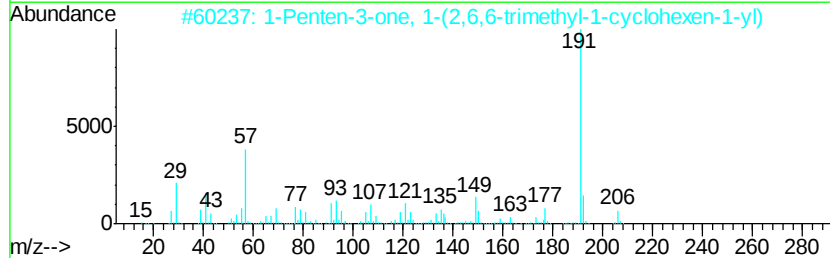
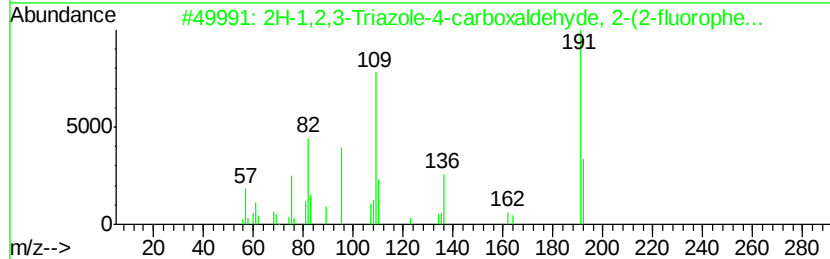
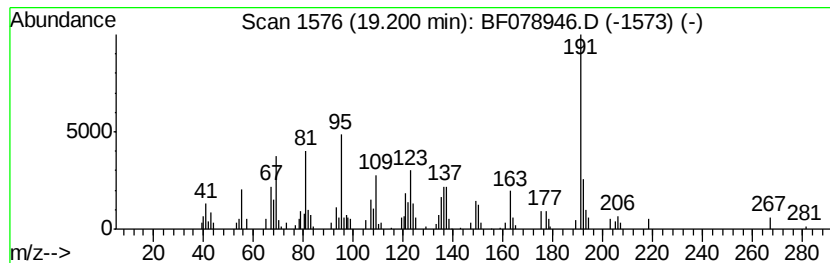
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown19.20 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.20	2.45 ng	103001	Perylene-d12	18.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H-1,2,3-Triazole-4-carboxaldehv...	191	C9H6FN3O	051306-43-5	46
2		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	38
3		1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	38
4		Anthracene, 9-dodecyltetradecahv...	360	C26H48	055401-75-7	38
5		2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050515\
Data File : BF078946.D
Acq On : 5 May 2015 17:56
Operator : TP/IZ
Sample : G2032-06
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B3(5-7)

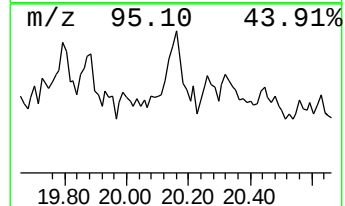
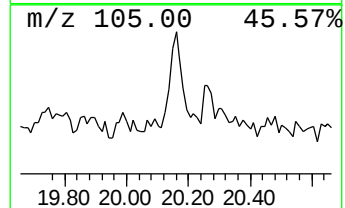
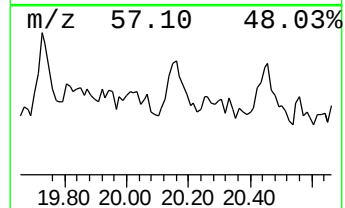
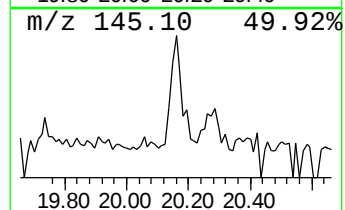
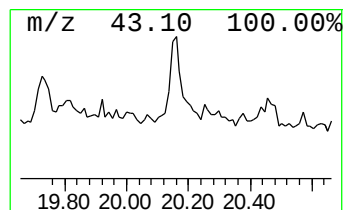
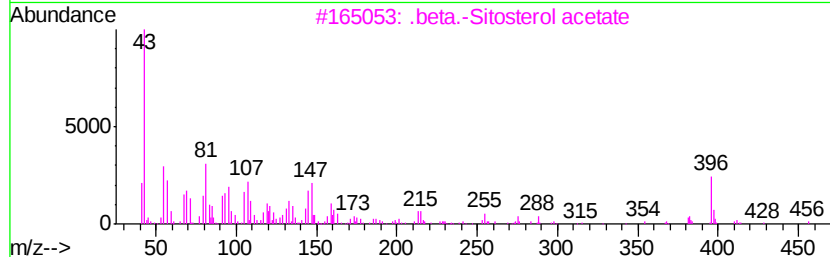
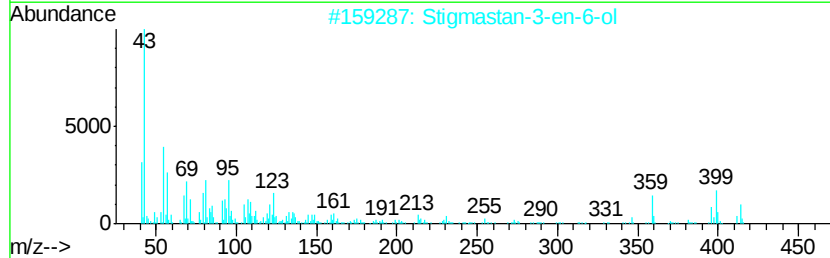
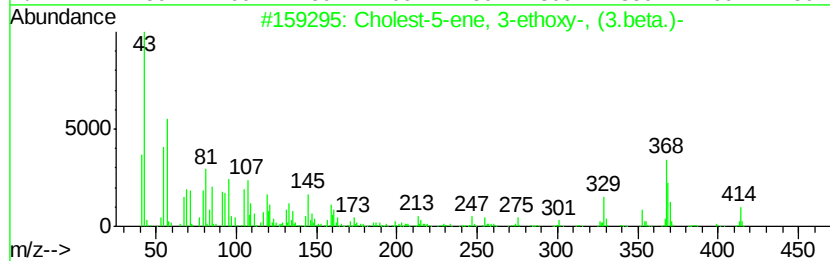
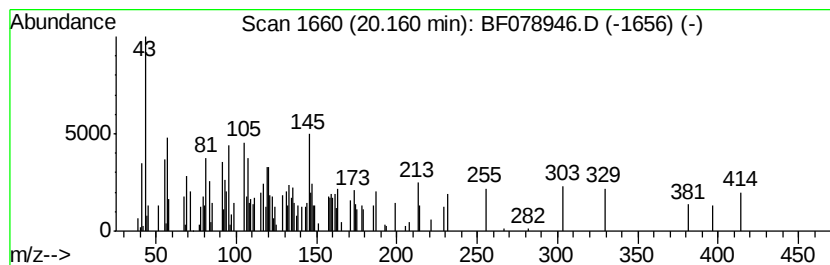
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 unknown20.16 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.16	3.23 ng	135789	Perylene-d12	18.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cholest-5-ene, 3-ethoxy-, (3.bet...	414	C29H50O	000986-19-6	46
2		Stigmastan-3-en-6-ol	414	C29H50O	1000214-19-7	12
3		.beta.-Sitosterol acetate	456	C31H52O2	000915-05-9	10
4		3-Hydroxypyridine monoacetate	137	C7H7NO2	017747-43-2	9
5		2-Hepten-6-one, 5-phenyl-, (Z)-	188	C13H16O	040931-42-8	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050515\
 Data File : BF078946.D
 Acq On : 5 May 2015 17:56
 Operator : TP/IZ
 Sample : G2032-06
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B3(5-7)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.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
unknown1.52	1.52	278.8	ng	6021800	1	7.36	431944	20.0
Butane, 2-methoxy...	1.84	13.7	ng	295090	1	7.36	431944	20.0
unknown2.16	2.16	3.2	ng	69729	1	7.36	431944	20.0
2-Pentanone, 4-hy...	5.16	11.6	ng	250253	1	7.36	431944	20.0
unknown7.06	7.06	80.9	ng	1747590	1	7.36	431944	20.0
Pentadecanoic acid	13.67	7.1	ng	246447	4	12.96	697858	20.0
2-Mercaptobenzoth...	13.85	2.3	ng	79147	4	12.96	697858	20.0
1,4-Benzenediamin...	14.54	10.0	ng	347740	4	12.96	697858	20.0
Phenol, 4,4'-(1-m...	14.81	17.1	ng	725686	5	16.25	847865	20.0
Pentafluoropropio...	16.13	6.1	ng	260152	5	16.25	847865	20.0
unknown18.75	18.75	2.7	ng	114828	6	18.00	840146	20.0
unknown19.20	19.20	2.5	ng	103001	6	18.00	840146	20.0
unknown20.16	20.16	3.2	ng	135789	6	18.00	840146	20.0