

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062618\
 Data File : BF106860.D
 Acq On : 27 Jun 2018 8:37
 Operator : JU/SJ
 Sample : J3693-01
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W-15

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.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	5.190	481	485	493	rBV	33979	40030	3.07%	0.474%
2	5.596	551	554	568	rBB	537241	513830	39.44%	6.090%
3	6.595	721	724	736	rBV	358635	326325	25.05%	3.868%
4	6.737	743	748	751	rBV	936319	862656	66.21%	10.224%
5	6.954	782	785	789	rBV	280649	239260	18.36%	2.836%
6	7.113	808	812	815	rBV	881013	752191	57.73%	8.915%
7	7.525	877	882	885	rBV	667453	638020	48.97%	7.562%
8	8.237	1000	1003	1009	rBV	276105	262596	20.15%	3.112%
9	8.972	1121	1128	1134	rBV	44056	59110	4.54%	0.701%
10	9.072	1142	1145	1152	rVB	24797	25218	1.94%	0.299%
11	9.213	1165	1169	1176	rVB4	30356	38108	2.92%	0.452%
12	9.313	1182	1186	1189	rBV	1183056	1034539	79.40%	12.261%
13	9.378	1195	1197	1205	rVB5	13128	21709	1.67%	0.257%
14	9.701	1249	1252	1261	rVB	17542	18489	1.42%	0.219%
15	9.995	1299	1302	1308	rBV2	285744	263784	20.25%	3.126%
16	10.189	1332	1335	1338	rBV	21225	20198	1.55%	0.239%
17	10.283	1348	1351	1354	rVB	32877	28367	2.18%	0.336%
18	10.613	1402	1407	1411	rVB	39335	34565	2.65%	0.410%
19	10.795	1434	1438	1441	rBV	694880	677397	51.99%	8.028%
20	10.907	1454	1457	1460	rVB	34575	29258	2.25%	0.347%
21	10.960	1463	1466	1468	rBV	67439	52605	4.04%	0.623%
22	10.989	1468	1471	1475	rVV2	72430	87211	6.69%	1.034%
23	11.025	1475	1477	1480	rVV	70059	61379	4.71%	0.727%
24	11.054	1480	1482	1484	rVB	32085	24938	1.91%	0.296%
25	11.142	1493	1497	1500	rVB3	49574	55503	4.26%	0.658%
26	11.178	1500	1503	1505	rVV	56018	48029	3.69%	0.569%
27	11.219	1508	1510	1513	rVB	33307	26967	2.07%	0.320%
28	11.489	1552	1556	1562	rBV	327923	275882	21.17%	3.270%
29	12.872	1789	1791	1796	rBV3	13231	13665	1.05%	0.162%
30	13.077	1821	1826	1829	rBV	1344658	1302921	100.00%	15.442%
31	13.619	1915	1918	1921	rBV	41568	34671	2.66%	0.411%
32	13.954	1972	1975	1982	rBV	38630	41676	3.20%	0.494%
33	14.142	2003	2007	2011	rBV	354401	325663	24.99%	3.860%
34	15.683	2263	2269	2275	rVB	150467	200664	15.40%	2.378%

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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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

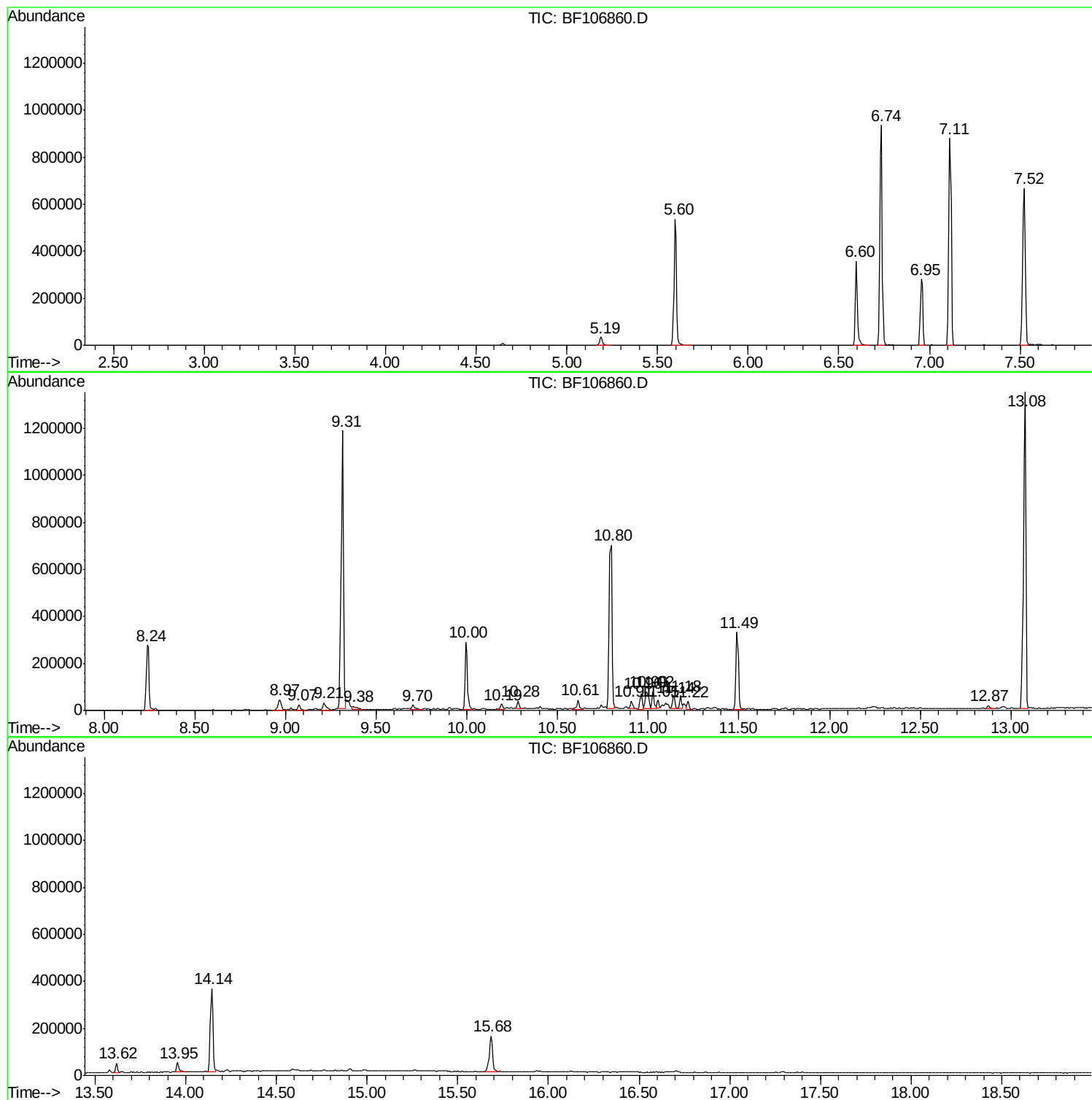
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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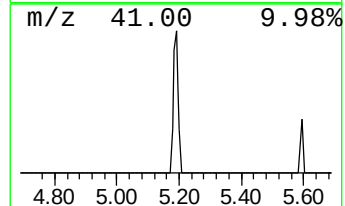
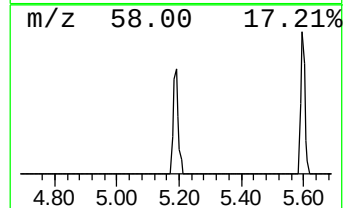
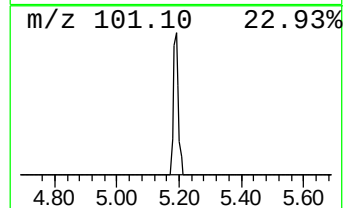
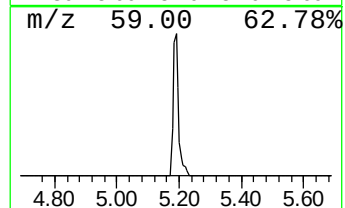
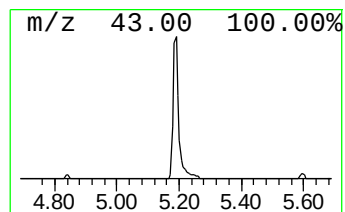
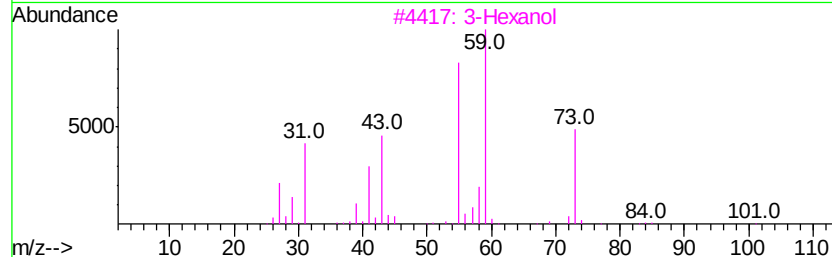
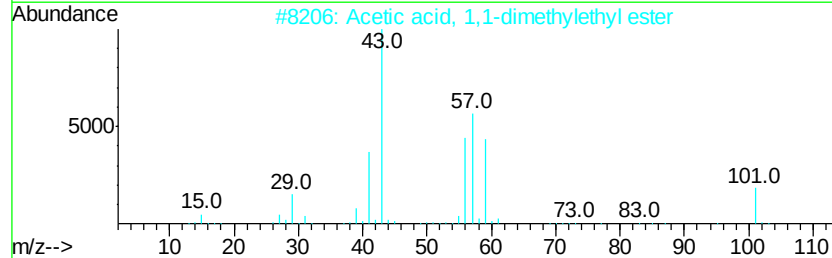
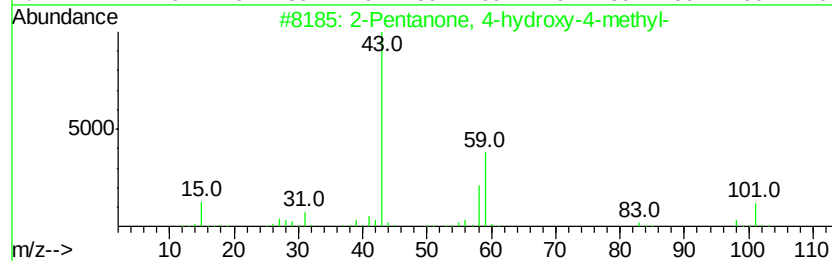
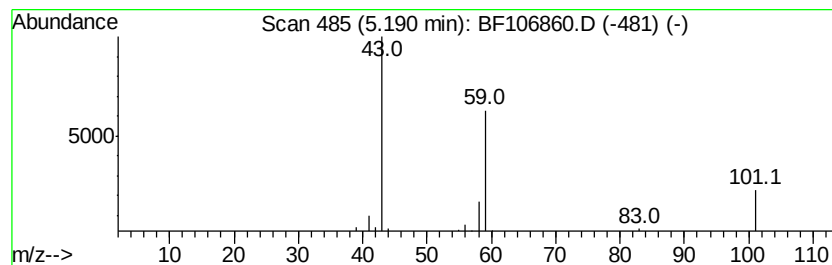
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	3.35 ng	40030	1,4-Dichlorobenzene-d4	6.95

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	40
3		3-Hexanol	102	C6H14O	000623-37-0	33
4		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	32
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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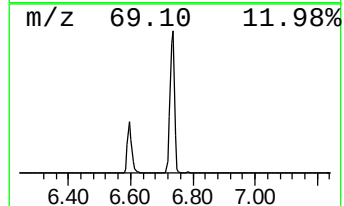
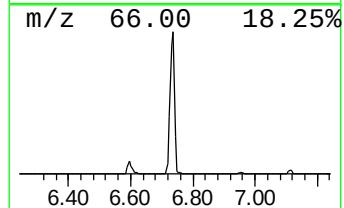
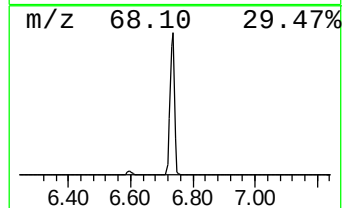
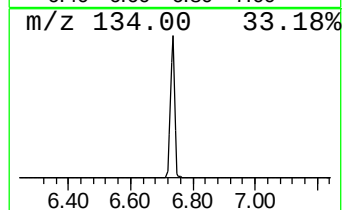
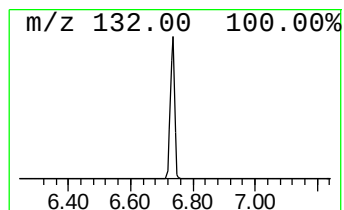
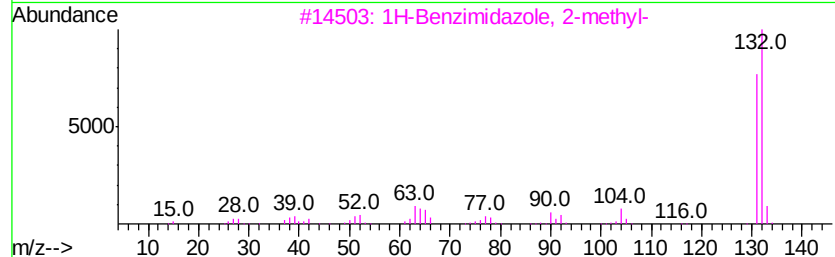
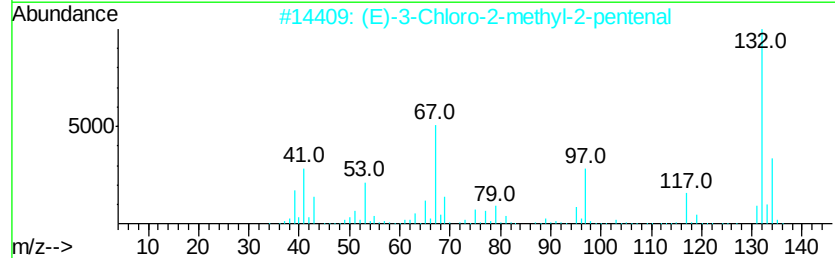
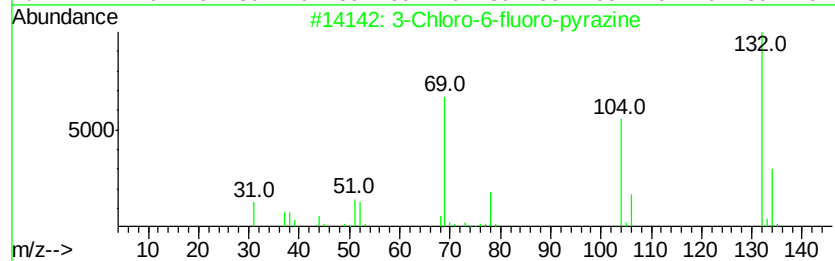
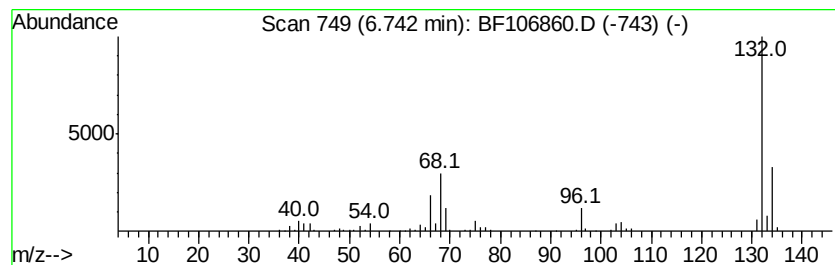
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	72.11 ng	862656	1,4-Dichlorobenzene-d4	6.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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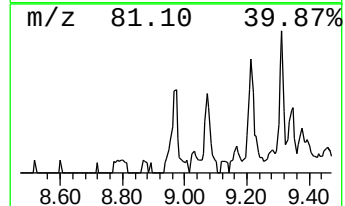
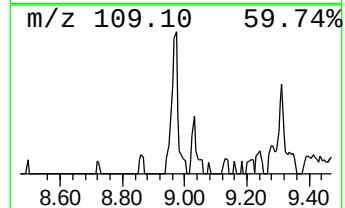
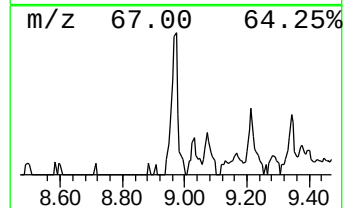
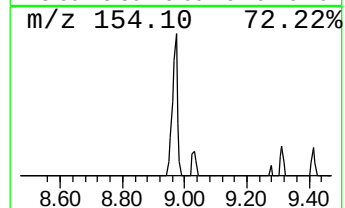
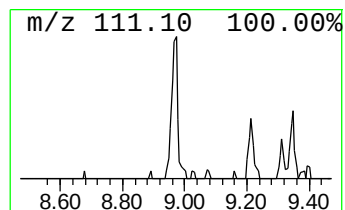
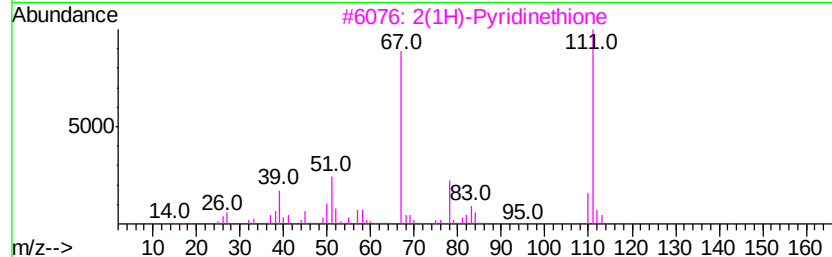
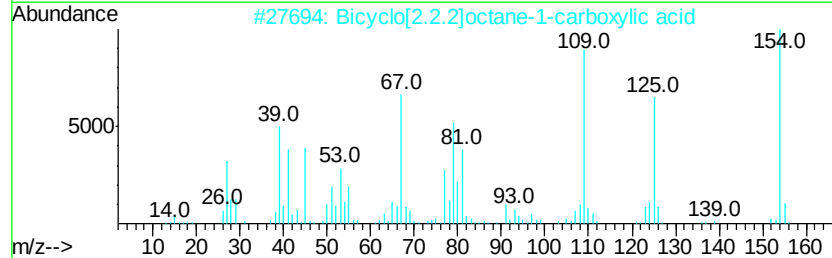
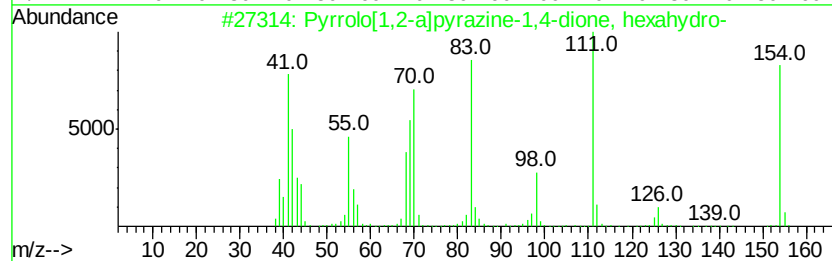
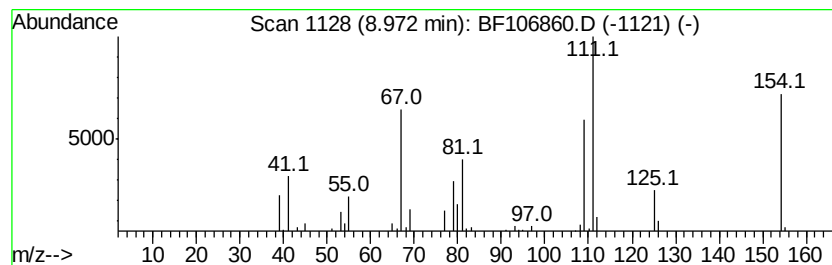
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown8.97 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.97	4.50 ng	59110	Naphthalene-d8	8.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrrolo[1,2-a]pyrazine-1,4-dione...	154	C7H10N2O2	019179-12-5	43
2		Bicyclo[2.2.2]octane-1-carboxyli...	154	C9H14O2	000699-55-8	43
3		2(1H)-Pyridinethione	111	C5H5NS	002637-34-5	35
4		Octahydronaphthalen-4a-ol	154	C10H18O	055693-34-0	32
5		2(1H)-Pyridinone, 3-hydroxy-	111	C5H5NO2	016867-04-2	18



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ALS Vial : 46 Sample Multiplier: 1

Instrument :
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ClientSampled :
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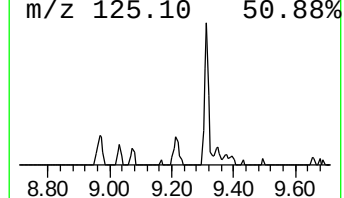
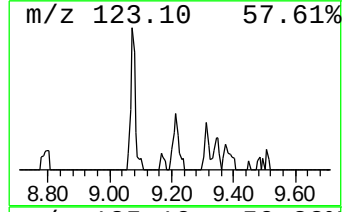
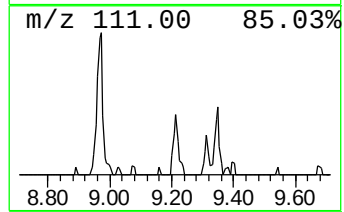
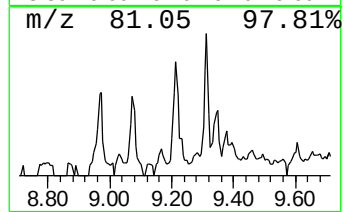
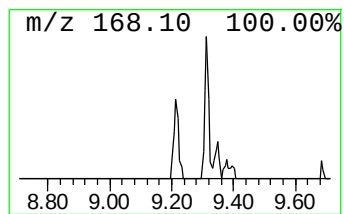
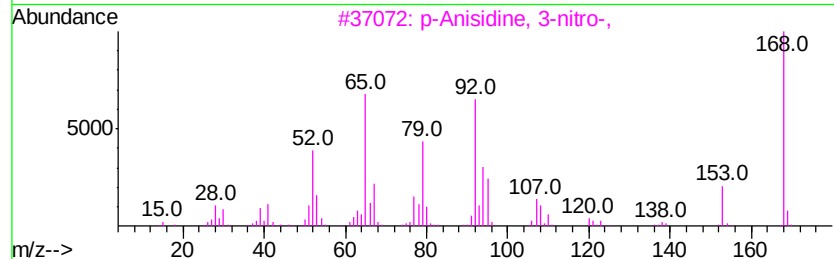
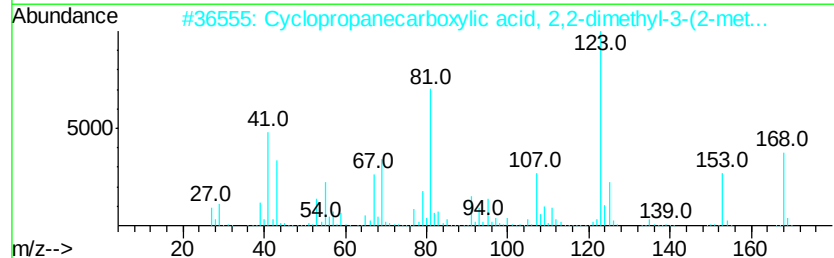
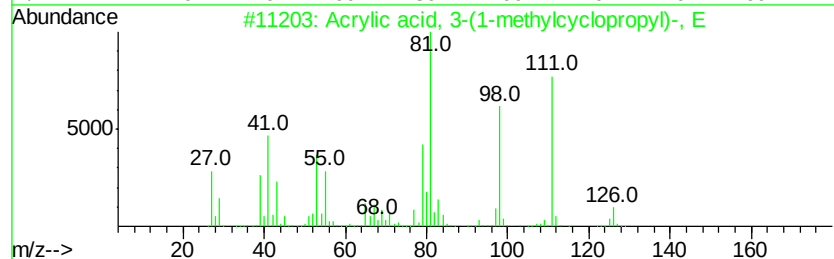
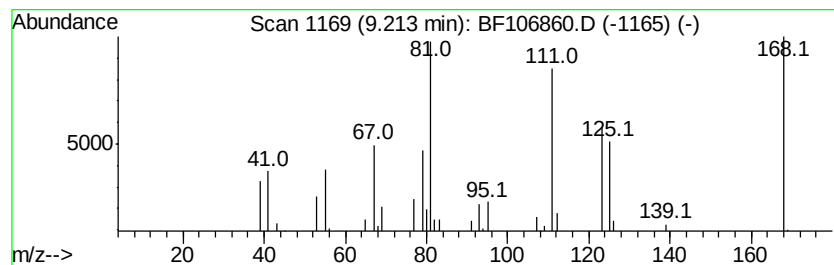
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown9.21 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.21	2.89 ng	38108	Acenaphthene-d10	10.00

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Acrylic acid, 3-(1-methylcyclopr...	126	C7H10O2	1000149-48-7	43
2		Cyclopropanecarboxylic acid, 2,2...	168	C10H16O2	015259-78-6	27
3		p-Anisidine, 3-nitro-,	168	C7H8N2O3	000577-72-0	25
4		2-Cyclohexen-1-one, 2-hydroxy-4,...	168	C10H16O2	064637-01-0	22
5		Bicyclo[3.3.1]nonane-1-carboxyli...	168	C10H16O2	017530-63-1	22



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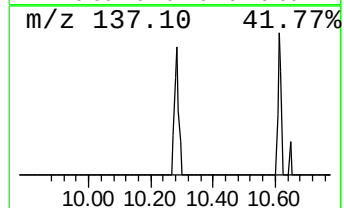
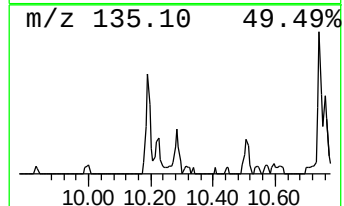
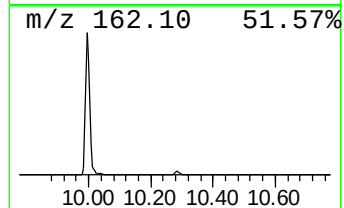
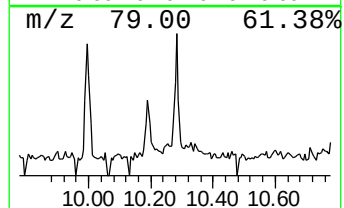
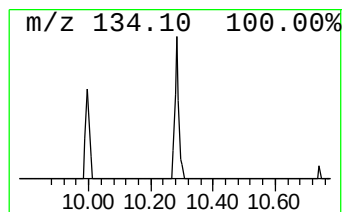
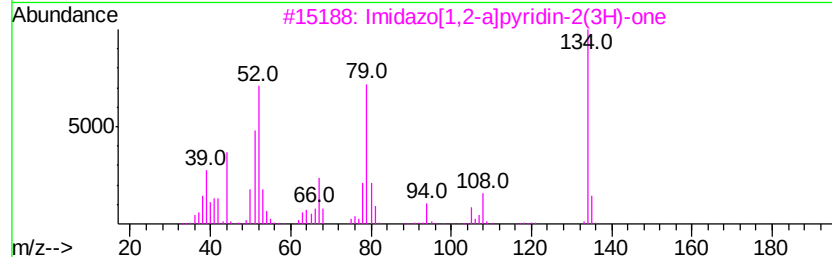
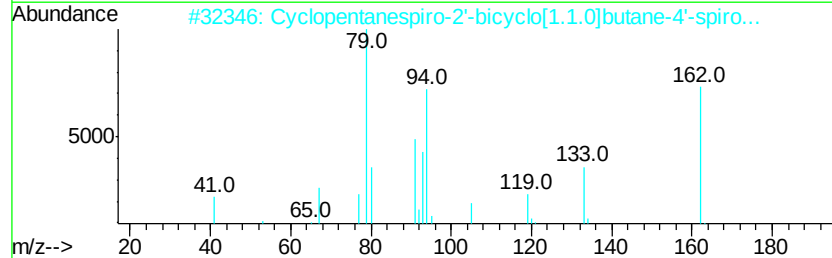
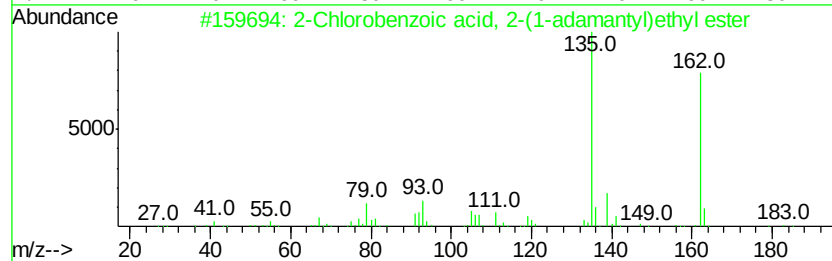
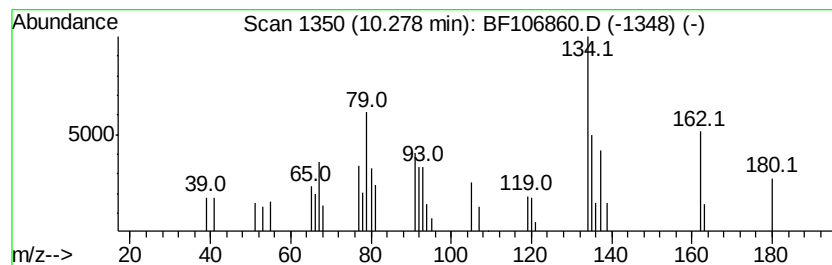
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 unknown10.28 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.28	2.15 ng	28367	Acenaphthene-d10	10.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Chlorobenzoic acid, 2-(1-adama...	318	C19H23ClO2	1000292-27-9	43
2		Cyclopentanespiro-2'-bicyclo[1.1...	162	C12H18	114310-66-6	38
3		Imidazo[1,2-a]pyridin-2(3H)-one	134	C7H6N2O	003999-06-2	38
4		Benzoic acid, 2-hydroxy-6-methyl...	180	C10H12O3	006555-40-4	30
5		Benzenamine, N,N,3-trimethyl-	135	C9H13N	000121-72-2	30



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Sample : J3693-01
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
W-15

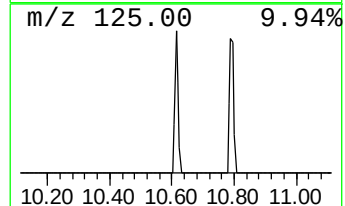
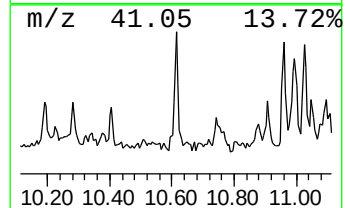
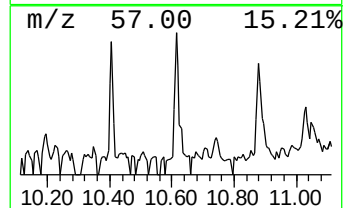
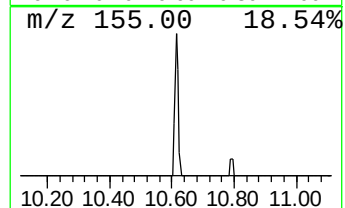
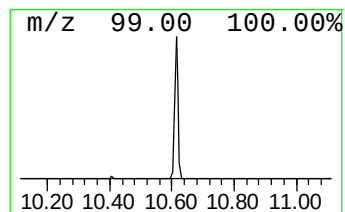
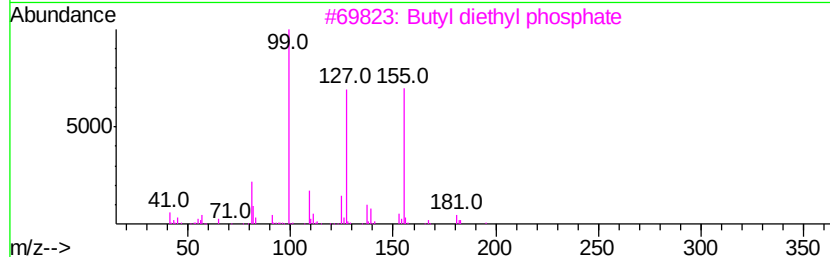
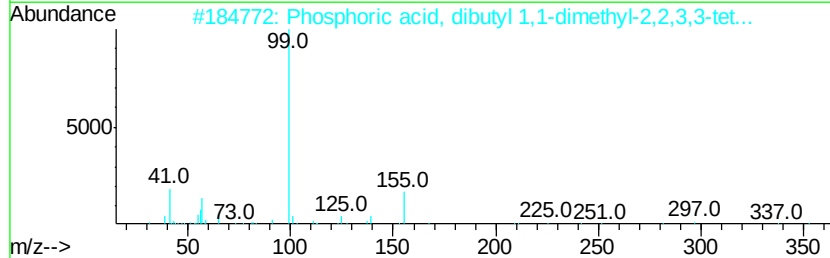
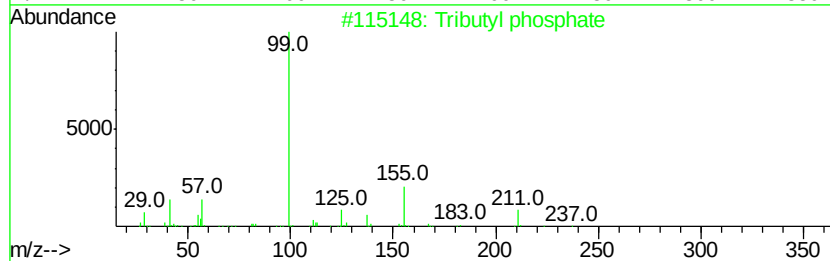
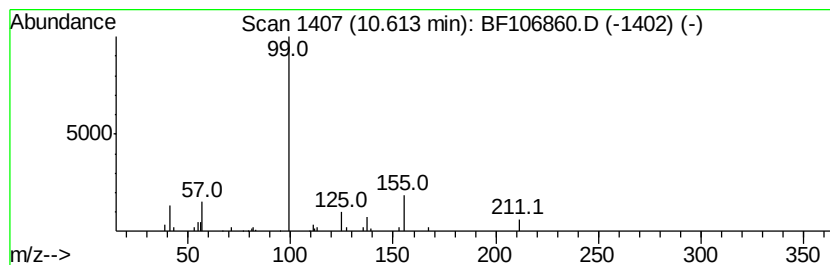
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tributyl phosphate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.61	2.62 ng	34565	Acenaphthene-d10	10.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tributyl phosphate	266	C12H27O4P	000126-73-8	90
2		Phosphoric acid, dibutyl 1,1-dim...	352	C13H25F4O4P	1000298-93-9	43
3		Butyl diethyl phosphate	210	C8H19O4P	1000298-58-0	38
4		Phosphoric acid, dipropyl nonyl ...	308	C15H33O4P	1000309-01-0	36
5		Tripropyl phosphate	224	C9H21O4P	000513-08-6	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062618\
Data File : BF106860.D
Acq On : 27 Jun 2018 8:37
Operator : JU/SJ
Sample : J3693-01
Misc :
ALS Vial : 46 Sample Multiplier: 1

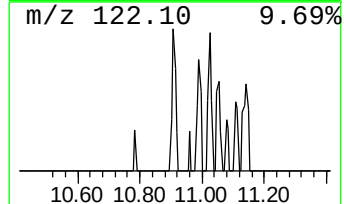
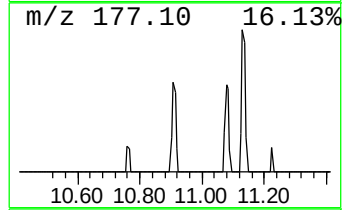
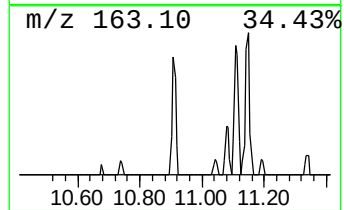
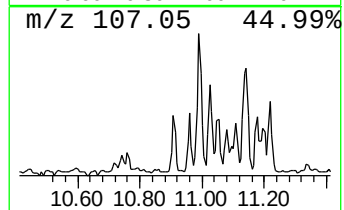
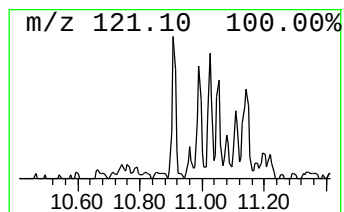
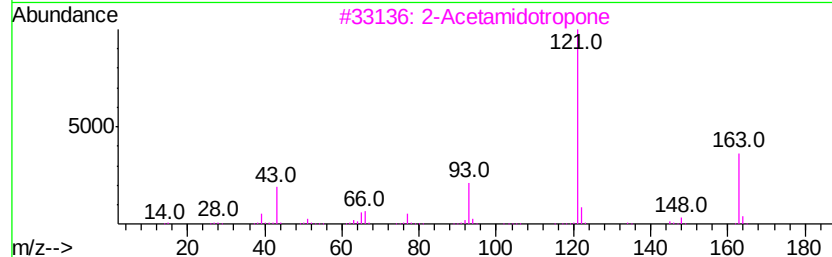
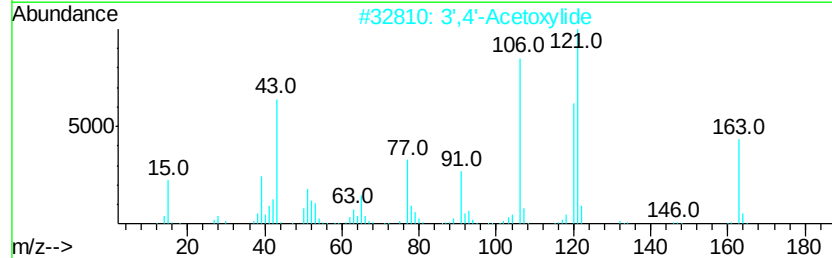
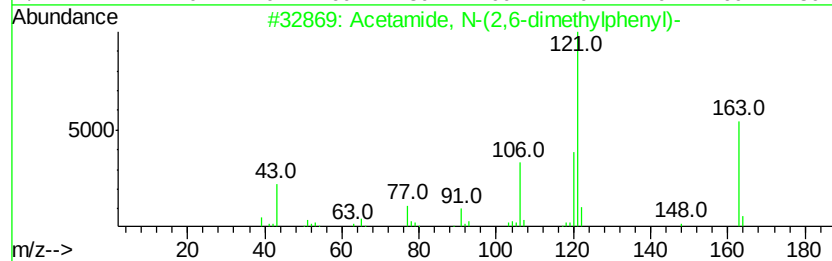
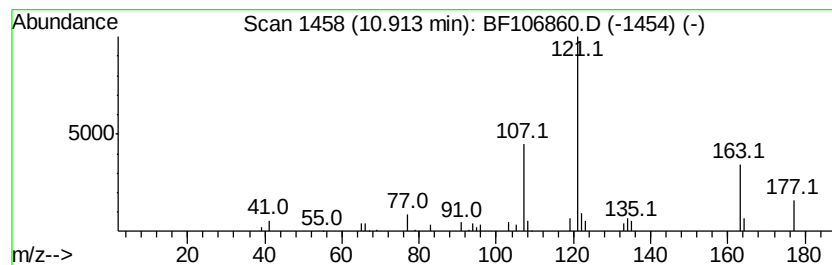
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ClientSampled :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Acetamide, N-(2,6-dimethylp... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.91	2.12 ng	29258	Phenanthrene-d10		11.49	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Acetamide, N-(2,6-dimethylphenyl)-	163	C10H13NO	002198-53-0	52
2		3',4'-Acetoxylide	163	C10H13NO	002198-54-1	50
3		2-Acetamidotropone	163	C9H9NO2	006422-12-4	49
4		Acetamide, N-methyl-N-(4-methylp...	163	C10H13NO	000612-03-3	47
5		Oxalic acid, isobutyl 2-isopropy...	264	C15H20O4	1000309-56-2	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062618\
Data File : BF106860.D
Acq On : 27 Jun 2018 8:37
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Sample : J3693-01
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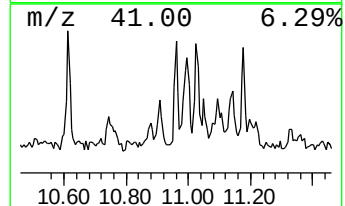
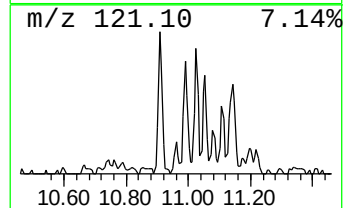
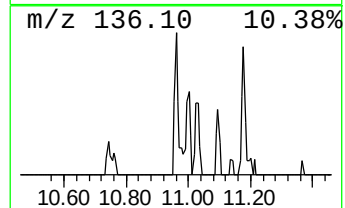
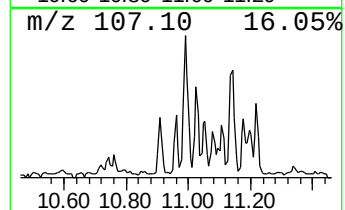
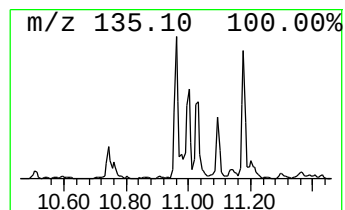
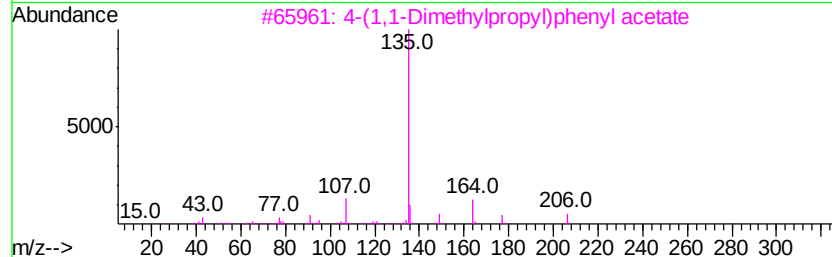
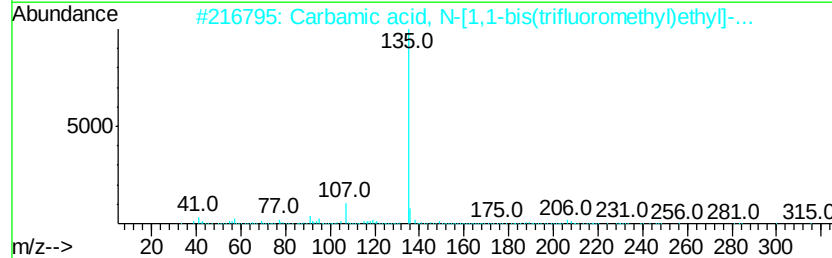
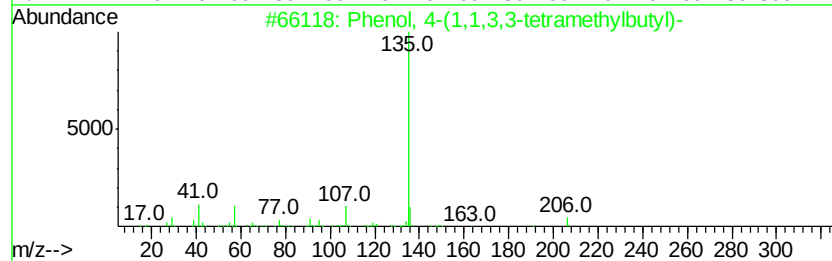
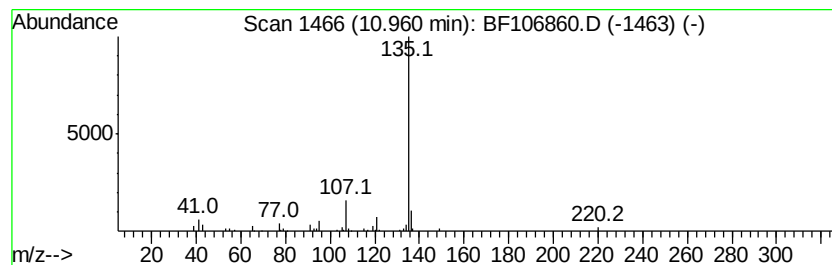
Instrument :
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.96	3.81 ng	52605	Phenanthrene-d10	11.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
2	Carbamic acid, N-[1,1-bis(triflu...	413	C19H25F6N02	296242-69-8	72
3	4-(1,1-Dimethylpropyl)phenyl ace...	206	C13H18O2	1000373-45-3	64
4	1-Adamantanecarboxylic acid chlo...	198	C11H15ClO	002094-72-6	59
5	4-tert-Butylphenyl acetate	192	C12H16O2	003056-64-2	56



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062618\
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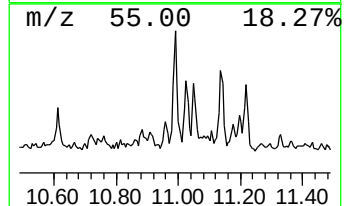
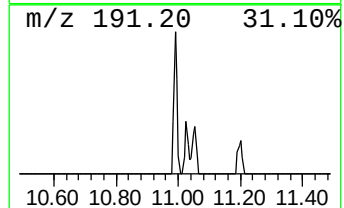
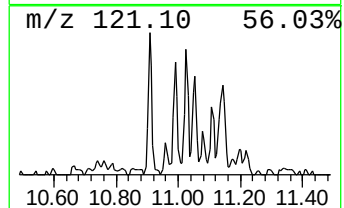
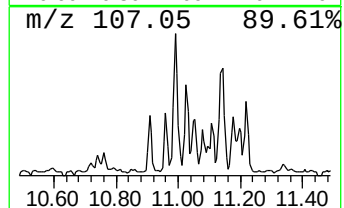
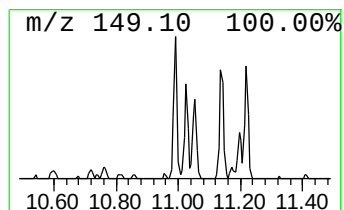
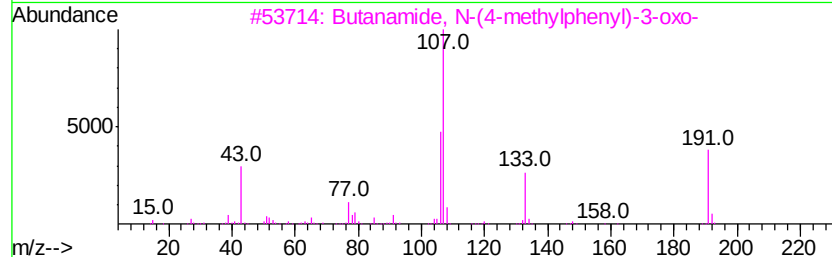
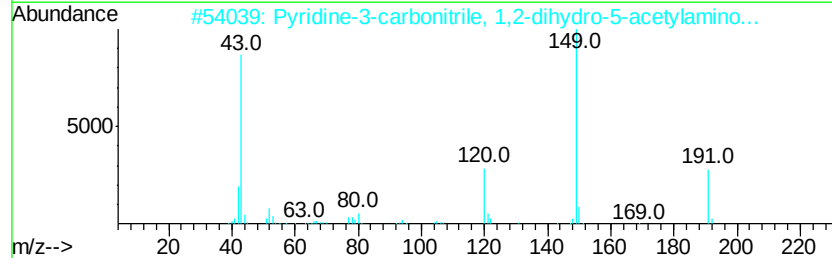
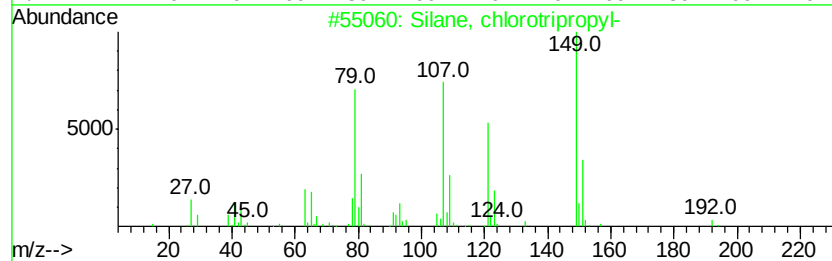
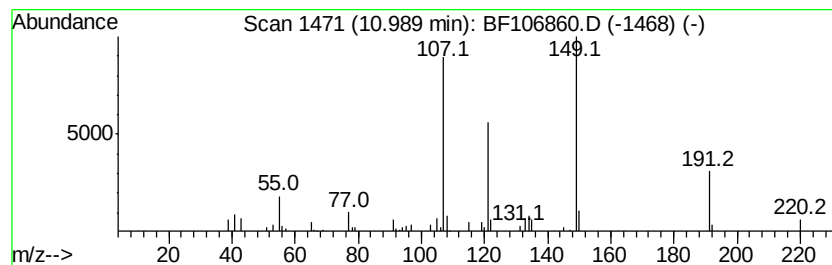
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Silane, chlorotripropyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.99	6.32 ng	87211	Phenanthrene-d10		11.49
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Silane, chlorotripropyl-	192	C9H21ClSi	000995-25-5	50
2	Pyridine-3-carbonitrile, 1,2-dih...	191	C9H9N3O2	220098-92-0	38
3	Butanamide, N-(4-methylphenyl)-3...	191	C11H13NO2	002415-85-2	38
4	Phenol, 2-(1,1-dimethylethyl)-5-...	164	C11H16O	000088-60-8	35
5	Butanamide, N-(2-methylphenyl)-3...	191	C11H13NO2	000093-68-5	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062618\
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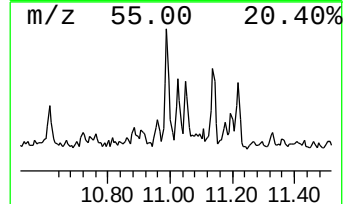
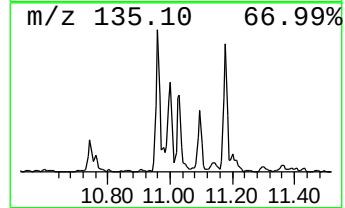
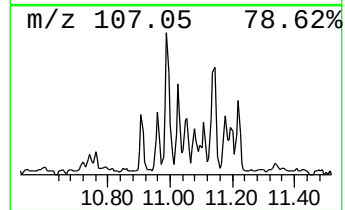
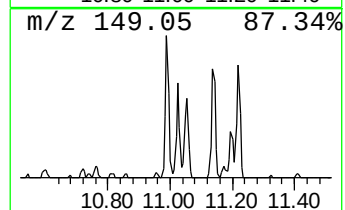
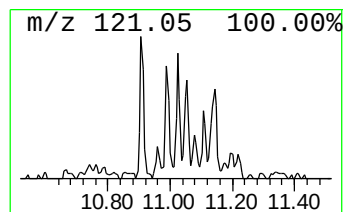
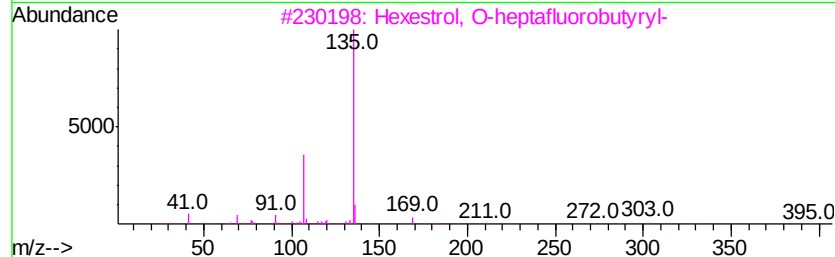
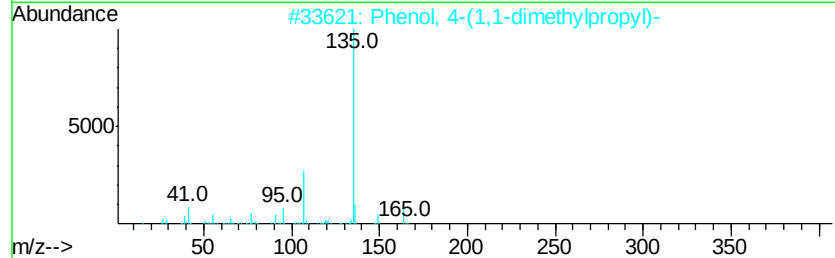
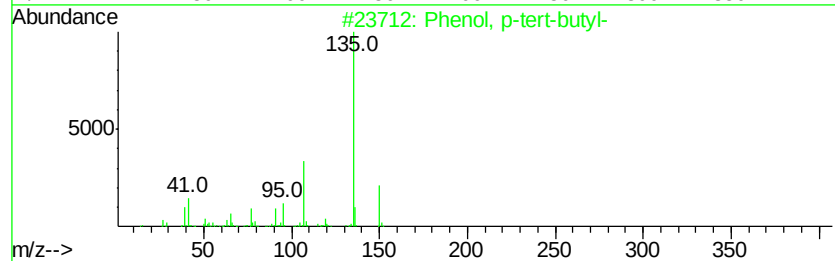
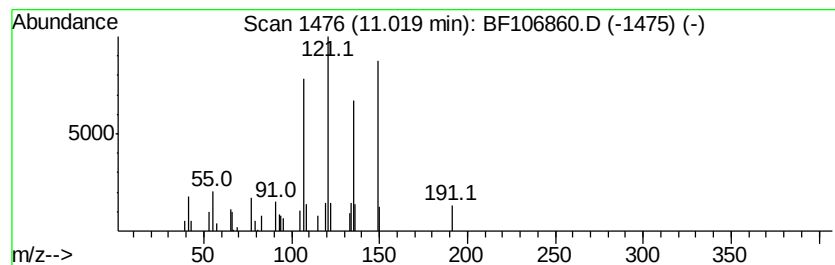
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Phenol, p-tert-butyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.02	4.45 ng	61379	Phenanthrene-d10	11.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	58
2	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	50
3	Hexestrol, O-heptafluorobutyl-	466	C22H21F7O3	1000365-31-9	50
4	Hexestrol, O-acetyl-	312	C20H24O3	1000374-87-8	50
5	Hexestrol	270	C18H22O2	005635-50-7	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062618\
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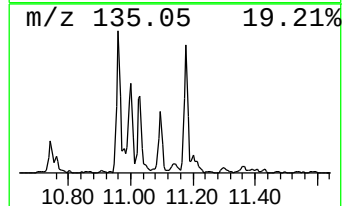
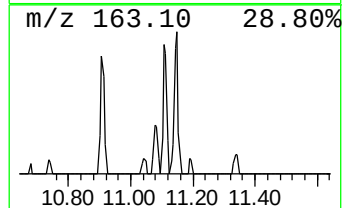
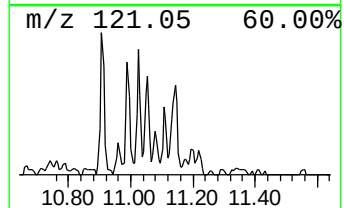
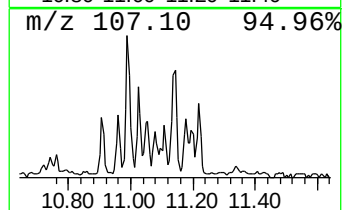
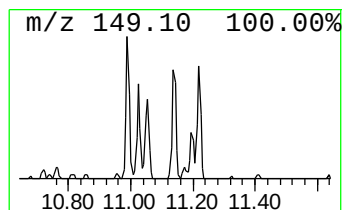
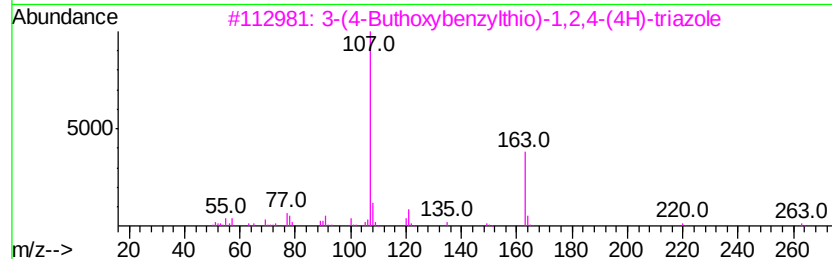
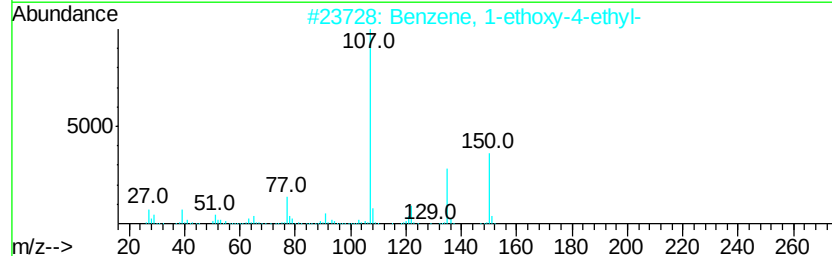
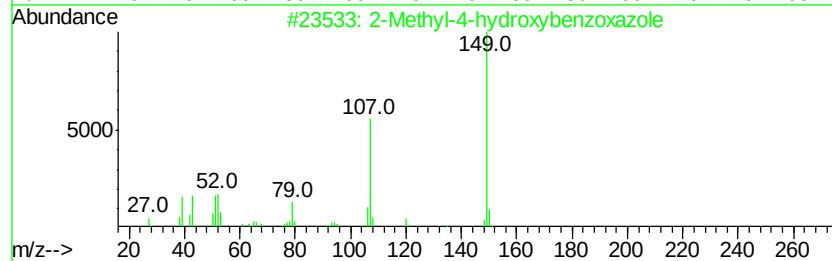
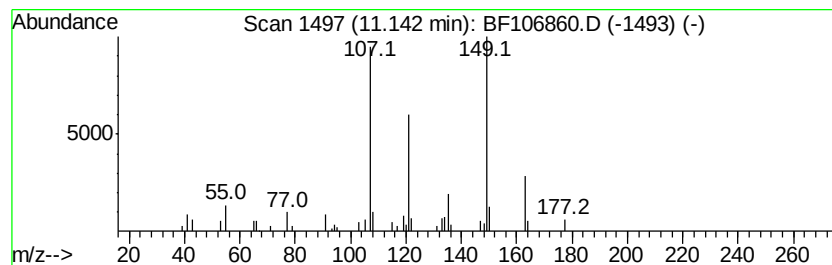
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown11.14 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.14	4.02 ng	55503	Phenanthrene-d10	11.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	47
2			Benzene, 1-ethoxy-4-ethyl-	150	C10H14O	001585-06-4	43
3			3-(4-Buthoxybenzylthio)-1,2,4-(4...	263	C13H17N3OS	057736-49-9	38
4			Phenol, 2-(1,1-dimethylethyl)-5-...	164	C11H16O	000088-60-8	38
5			5H-Pyrrolo(3,2-d)pyrimidine-2,4-...	149	C6H7N5	1000244-21-4	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062618\
Data File : BF106860.D
Acq On : 27 Jun 2018 8:37
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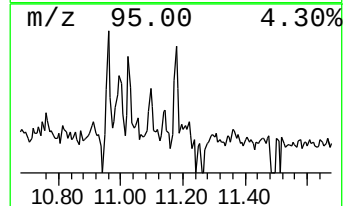
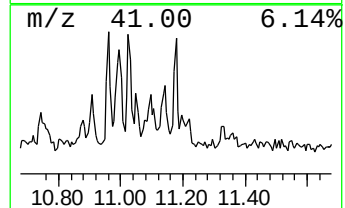
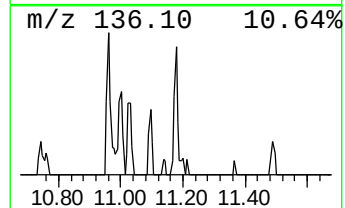
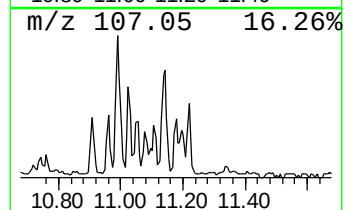
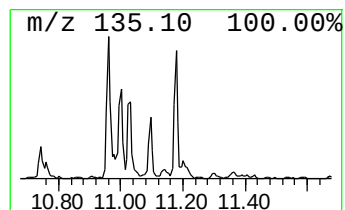
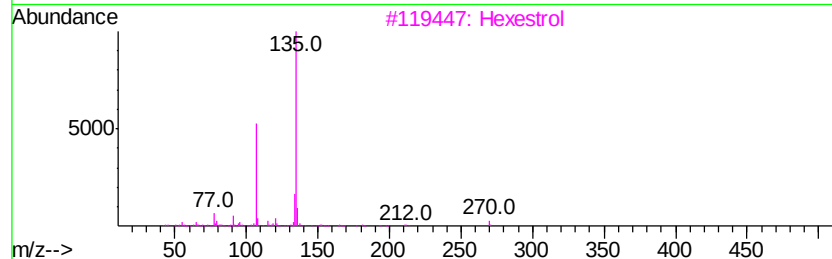
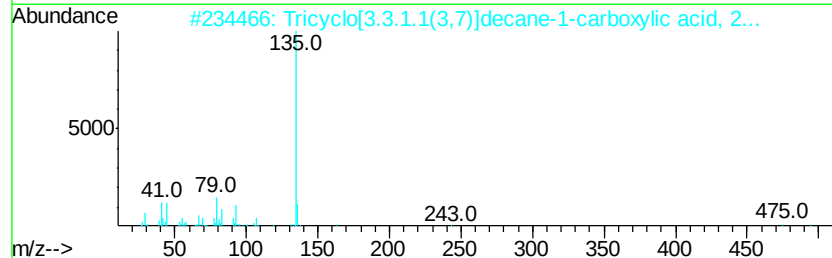
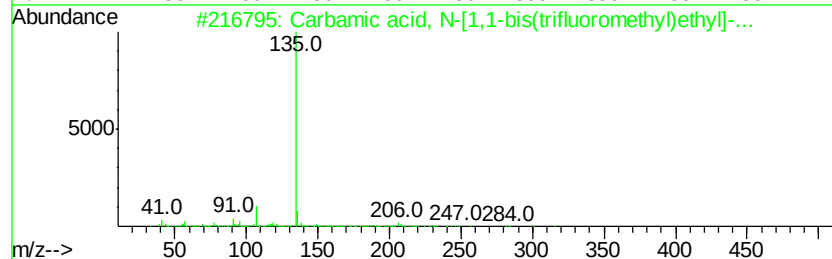
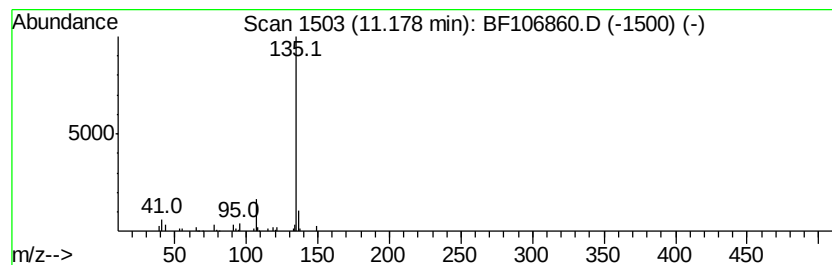
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Carbamic acid, N-[1,1-bis(t... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.18	3.48 ng	48029	Phenanthrene-d10	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Carbamic acid, N-[1,1-bis(triflu...	413	C19H25F6N02	296242-69-8	72
2		Tricyclo[3.3.1.1(3,7)]decane-1-c...	494	C18H18F1202	1000141-65-1	50
3		Hexestrol	270	C18H2202	000084-16-2	50
4		Hexestrol, 0-trifluoroacetyl-	366	C20H21F303	1000365-31-7	45
5		o-Anisic acid, tridec-2-ynyl ester	330	C21H3003	1000292-57-9	45



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Data File : BF106860.D
Acq On : 27 Jun 2018 8:37
Operator : JU/SJ
Sample : J3693-01
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
W-15

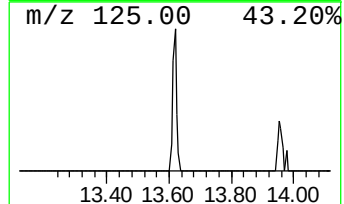
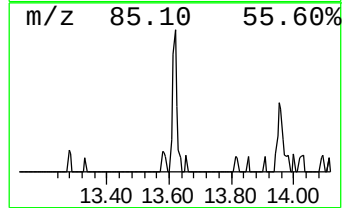
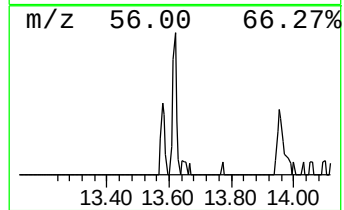
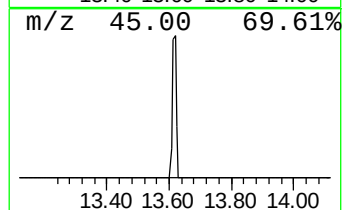
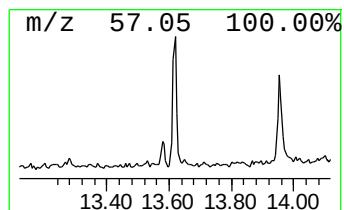
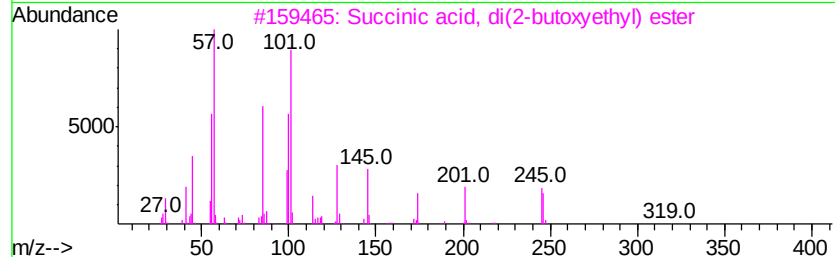
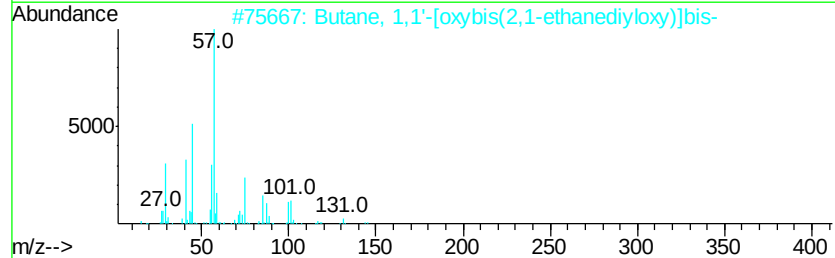
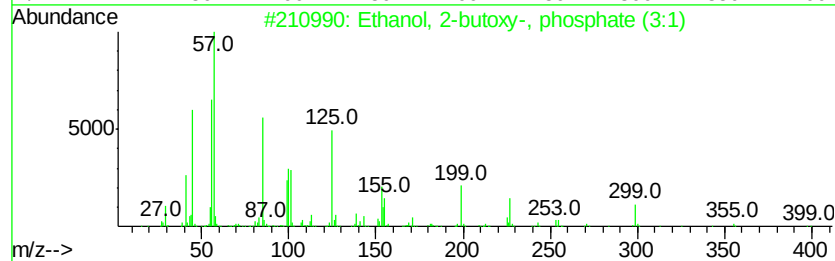
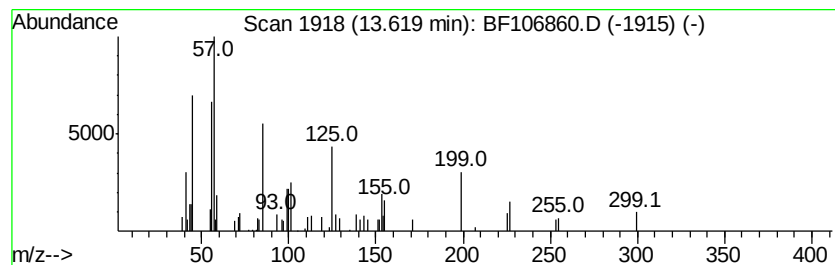
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Ethanol, 2-butoxy-, phospho... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.62	2.13 ng	34671	Chrysene-d12	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-butoxy-, phosphate (3:1)	398	C18H39O7P	000078-51-3	91
2		Butane, 1,1'-[oxybis(2,1-ethaned...	218	C12H26O3	000112-73-2	27
3		Succinic acid, di(2-butoxyethyl)...	318	C16H30O6	1000325-84-7	25
4		3-Cyclopentylpropionic acid, 4-m...	242	C14H26O3	1000354-33-2	25
5		trans-2,3-Epoxyoctane	128	C8H16O	028180-70-3	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062618\
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Sample : J3693-01
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
W-15

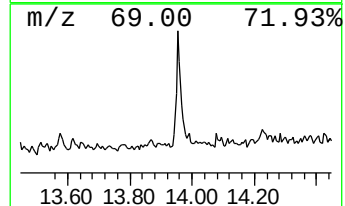
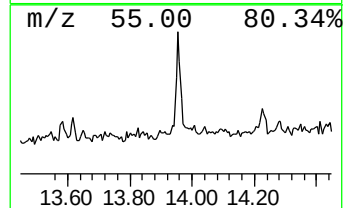
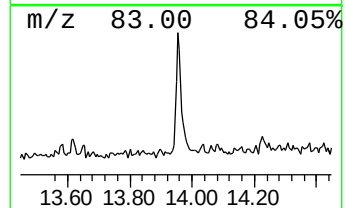
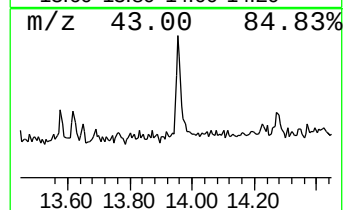
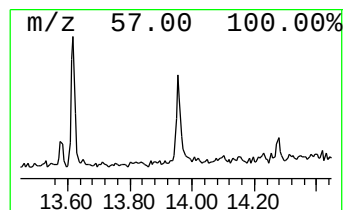
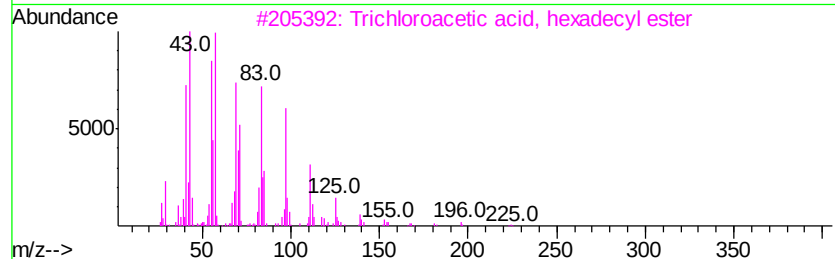
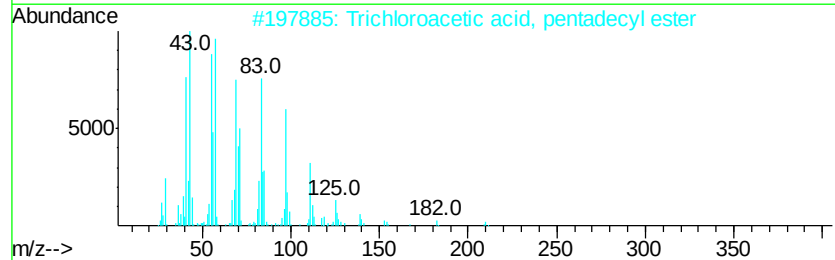
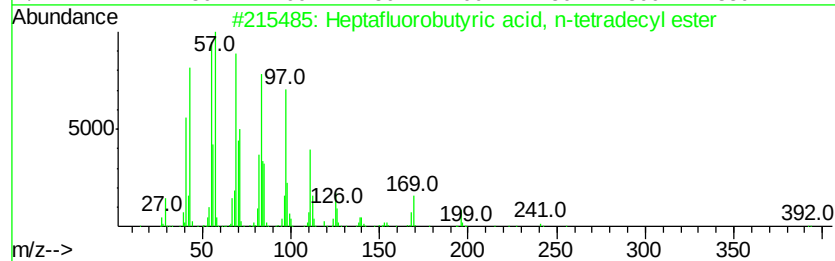
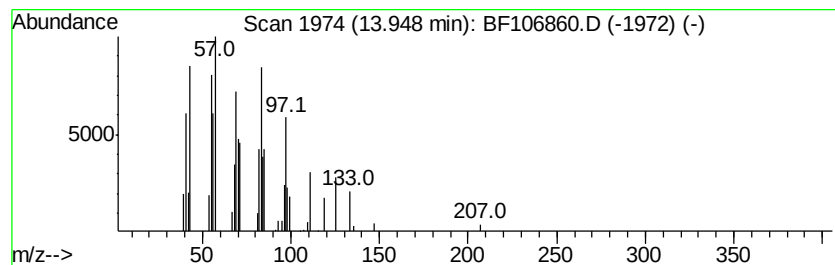
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Heptafluorobutyric acid, n-... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	2.56 ng	41676	Chrysene-d12	14.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	91
2			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
3			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
4			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	90
5			n-Tetracosanol-1	354	C24H50O	000506-51-4	87



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Instrument :
 BNA_F
 ClientSampleId :
 W-15

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	3.4	ng	40030	1	6.95	239260	20.0
unknown6.74	6.74	72.1	ng	862656	1	6.95	239260	20.0
unknown8.97	8.97	4.5	ng	59110	2	8.24	262596	20.0
unknown9.21	9.21	2.9	ng	38108	3	10.00	263784	20.0
unknown10.28	10.28	2.1	ng	28367	3	10.00	263784	20.0
Tributyl phosphate	10.61	2.6	ng	34565	3	10.00	263784	20.0
Acetamide, N-(2,6...	10.91	2.1	ng	29258	4	11.49	275882	20.0
Phenol, 4-(1,1,3,...	10.96	3.8	ng	52605	4	11.49	275882	20.0
Silane, chlorotri...	10.99	6.3	ng	87211	4	11.49	275882	20.0
Phenol, p-tert-bu...	11.02	4.5	ng	61379	4	11.49	275882	20.0
unknown11.14	11.14	4.0	ng	55503	4	11.49	275882	20.0
Carbamic acid, N-...	11.18	3.5	ng	48029	4	11.49	275882	20.0
Ethanol, 2-butoxy...	13.62	2.1	ng	34671	5	14.14	325663	20.0
Heptafluorobutyri...	13.95	2.6	ng	41676	5	14.14	325663	20.0