

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102716\
 Data File : BF090893.D
 Acq On : 28 Oct 2016 3:24
 Operator : UM/SJ
 Sample : H5430-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EME-BLEND-007

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-BF102616.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	4.933	246	249	256	rBV	398836	551937	17.09%	1.754%
2	5.367	284	287	290	rBV	2727207	2935087	90.86%	9.327%
3	6.419	376	379	381	rBV	2687762	3130880	96.92%	9.949%
4	6.544	387	390	392	rBV	2854438	3082933	95.44%	9.796%
5	6.773	408	410	419	rVB	986803	924452	28.62%	2.938%
6	6.933	421	424	426	rBV	2427154	2626314	81.30%	8.345%
7	7.344	457	460	462	rBV	1440383	1761390	54.53%	5.597%
8	8.064	520	523	525	rBV	1097848	1170600	36.24%	3.720%
9	9.139	615	617	619	rBV	3722168	3230287	100.00%	10.265%
10	9.539	649	652	654	rBV	573850	566928	17.55%	1.801%
11	9.813	674	676	678	rBV	1471718	1217723	37.70%	3.869%
12	10.259	711	715	716	rBV2	48077	73285	2.27%	0.233%
13	10.476	730	734	735	rBV	26604	38108	1.18%	0.121%
14	10.602	742	745	747	rBV	1925336	1793872	55.53%	5.700%
15	10.728	754	756	760	rVB	83248	91690	2.84%	0.291%
16	11.173	793	795	797	rBV	44067	40924	1.27%	0.130%
17	11.299	803	806	808	rBV	908962	965848	29.90%	3.069%
18	12.888	942	945	947	rBV	2007943	2336004	72.32%	7.423%
19	13.791	1021	1024	1029	rBV	73523	164507	5.09%	0.523%
20	13.928	1034	1036	1039	rBV	790978	956971	29.62%	3.041%
21	14.111	1048	1052	1053	rBV3	27568	71315	2.21%	0.227%
22	14.419	1077	1079	1085	rBV3	45399	112599	3.49%	0.358%
23	14.785	1109	1111	1113	rVB	31733	40659	1.26%	0.129%
24	15.025	1130	1132	1136	rVB	39107	55930	1.73%	0.178%
25	15.345	1157	1160	1162	rBV	597049	747819	23.15%	2.376%
26	15.562	1177	1179	1181	rBV	40864	70588	2.19%	0.224%
27	16.031	1218	1220	1225	rVB2	48955	115433	3.57%	0.367%
28	16.522	1261	1263	1267	rVB4	30552	57100	1.77%	0.181%
29	16.865	1291	1293	1297	rBV5	29592	51899	1.61%	0.165%
30	17.243	1322	1326	1329	rBV2	89278	229840	7.12%	0.730%
31	17.437	1337	1343	1348	rBV3	67085	211000	6.53%	0.670%
32	17.540	1349	1352	1355	rBV2	74998	165283	5.12%	0.525%
33	17.757	1368	1371	1380	rVB5	88746	290748	9.00%	0.924%
34	17.974	1387	1390	1391	rBV3	31551	63901	1.98%	0.203%

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ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EME-BLEND-007

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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35	18.111	1399	1402	1405	rBV3	36283	98292	3.04%	0.312%
36	18.180	1405	1408	1415	rVB3	59821	181576	5.62%	0.577%
37	18.980	1473	1478	1486	rBV2	130546	468541	14.50%	1.489%
38	19.140	1487	1492	1503	rVB2	231426	777896	24.08%	2.472%

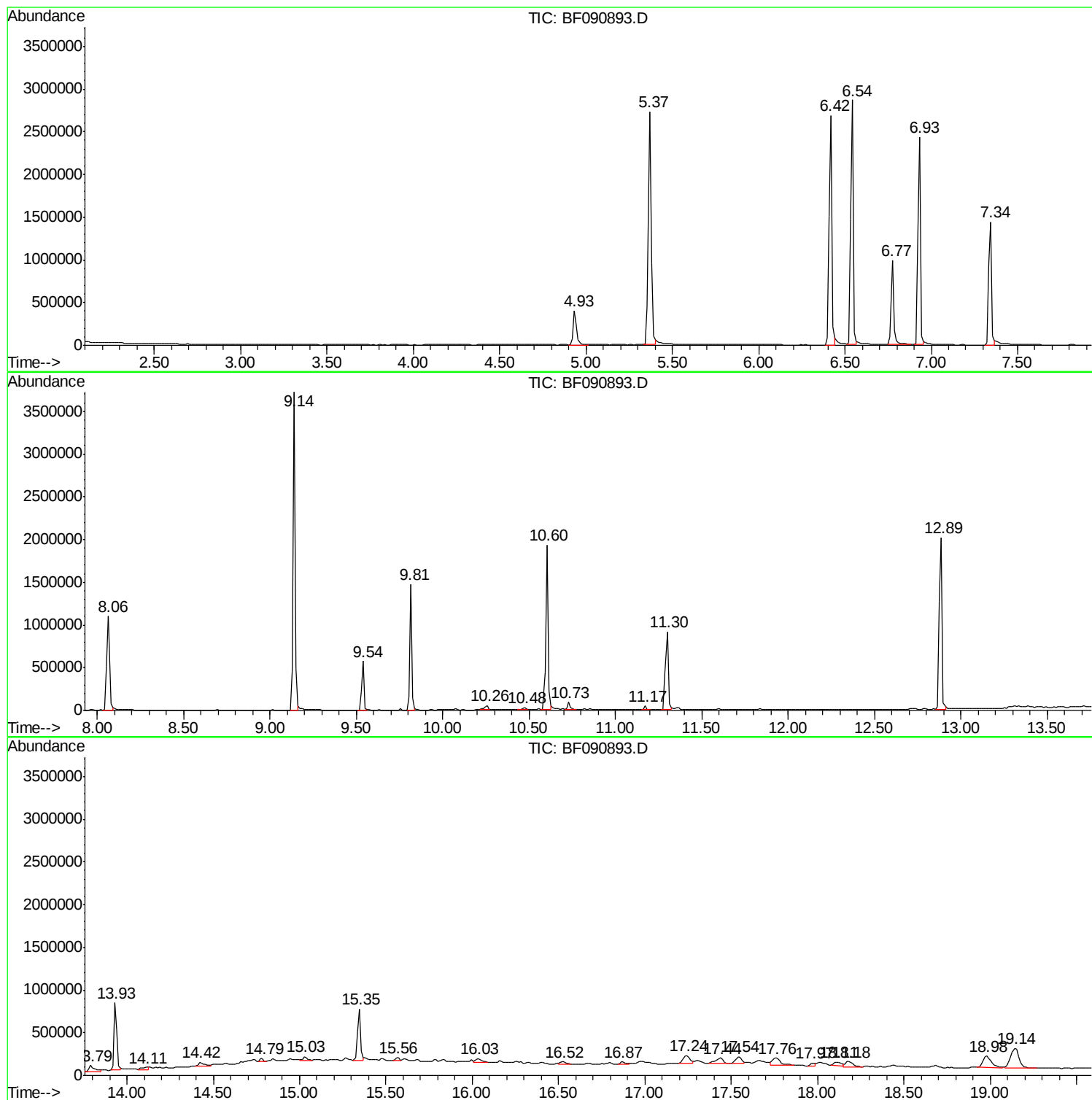
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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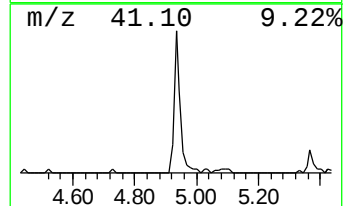
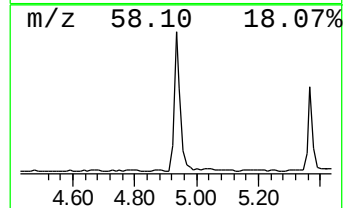
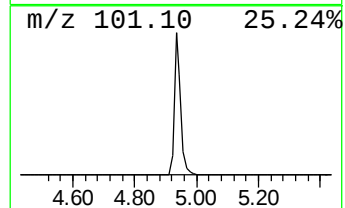
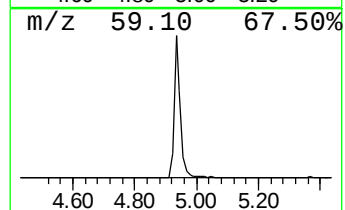
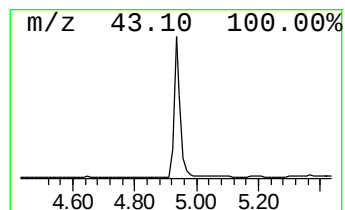
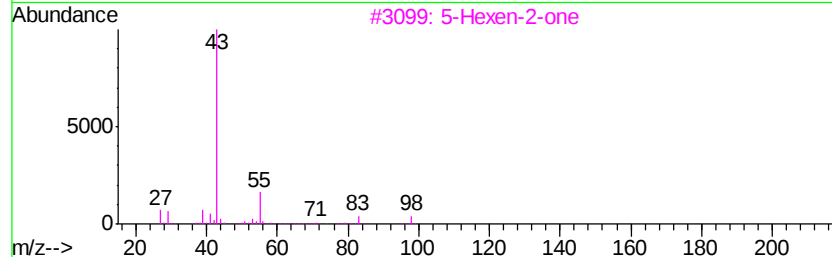
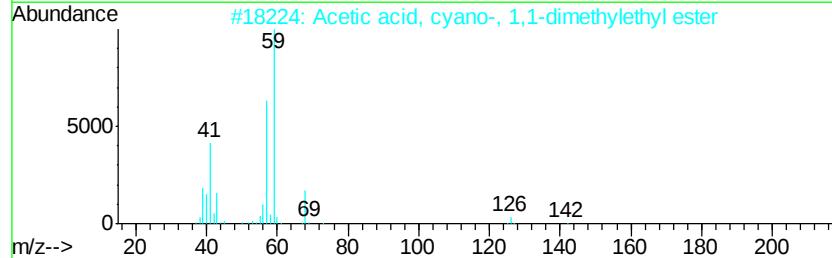
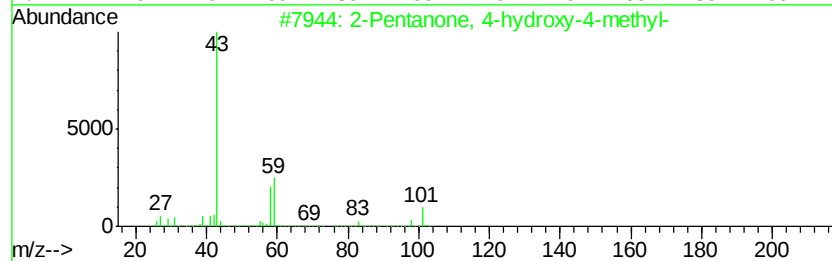
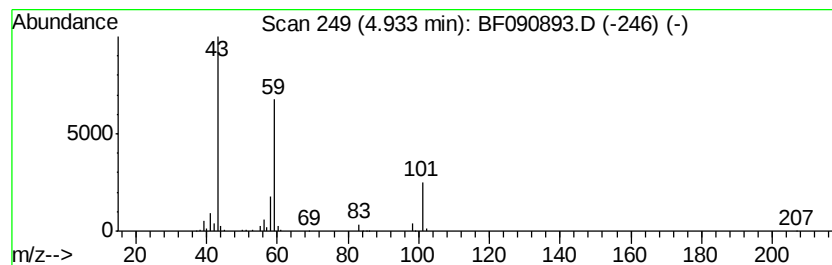
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.93	11.94 ng	551937	1,4-Dichlorobenzene-d4	6.77

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59	
2	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17	
3	5-Hexen-2-one	98	C6H10O	000109-49-9	9	
4	Acetone	58	C3H6O	000067-64-1	9	
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9	



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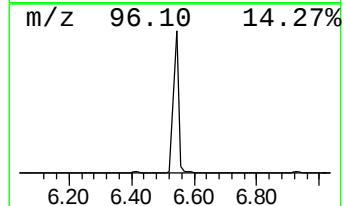
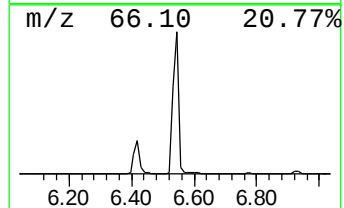
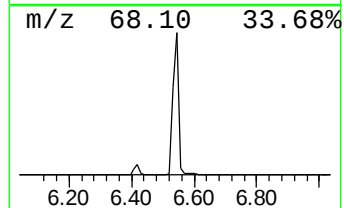
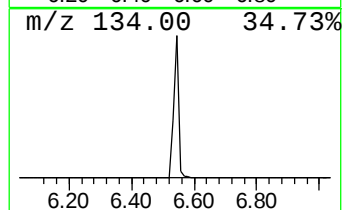
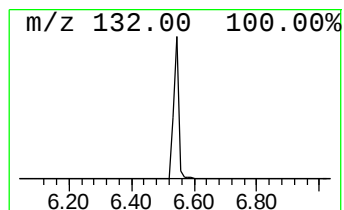
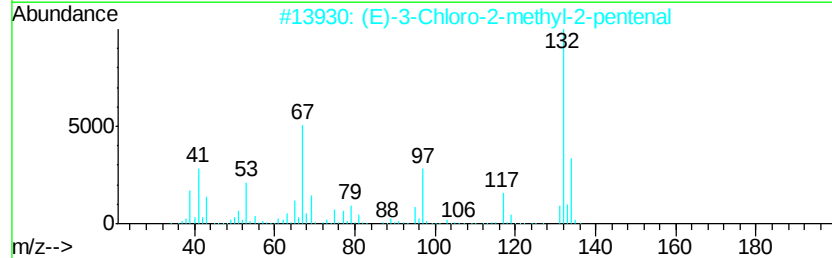
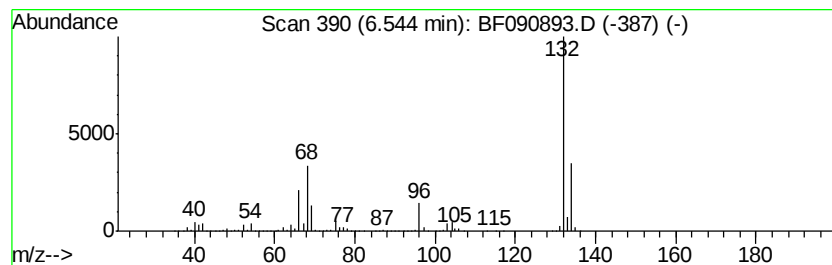
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.54 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.54	66.70 ng	3082930	1,4-Dichlorobenzene-d4	6.77

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	12
3		Tranvlcvpromine-propionyl	189	C12H15NO	1000123-86-3	12
4		Xenon	132	Xe	007440-63-3	9
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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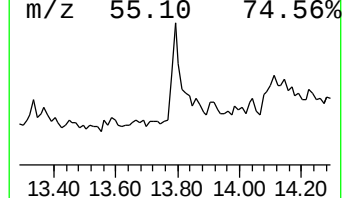
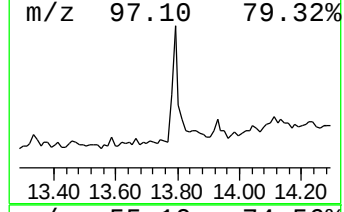
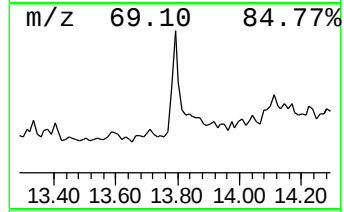
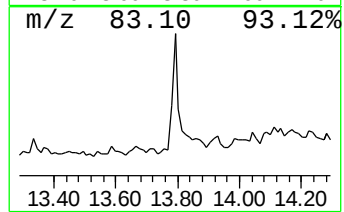
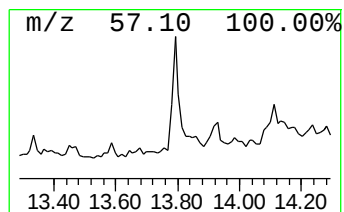
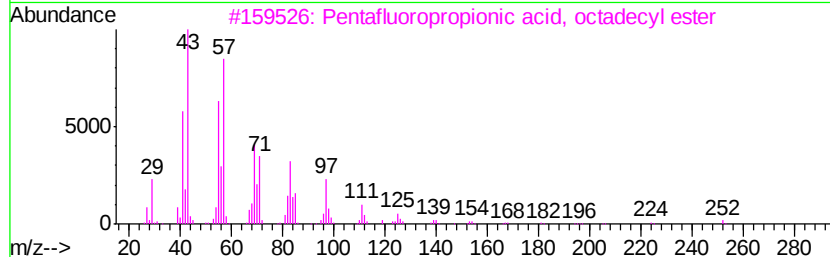
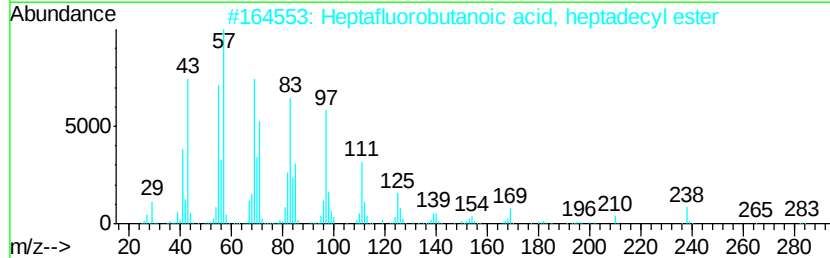
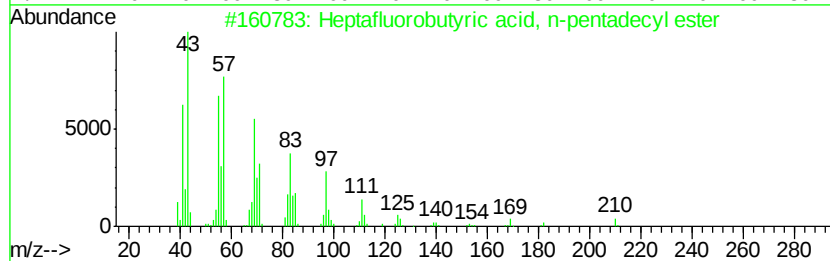
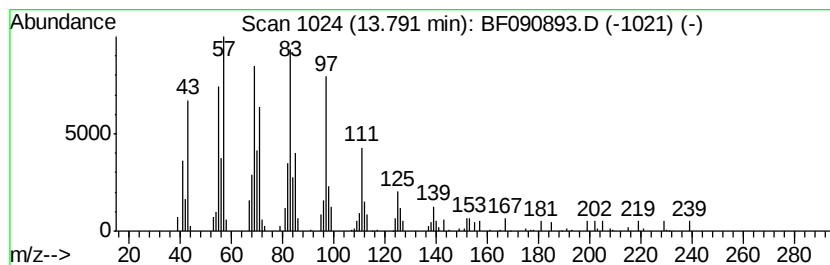
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Heptafluorobutyric acid, n-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	3.44 ng	164507	Chrysene-d12	13.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptafluorobutyric acid, n-penta...	424	C19H31F7O2	1000216-79-3	94
2		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	93
3		Pentafluoropropionic acid, octad...	416	C21H37F5O2	1000280-07-7	91
4		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
5		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	91



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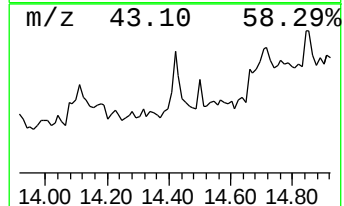
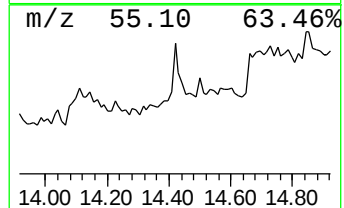
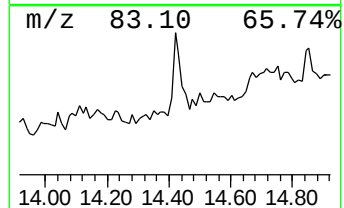
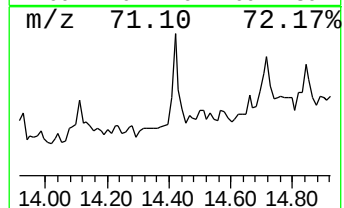
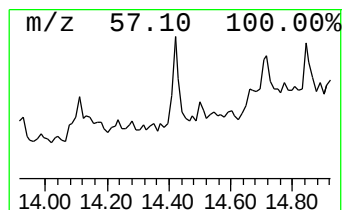
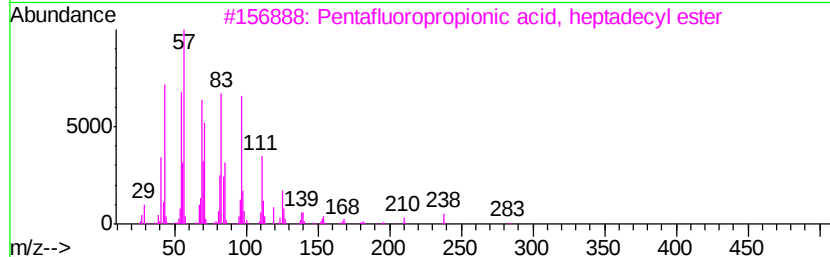
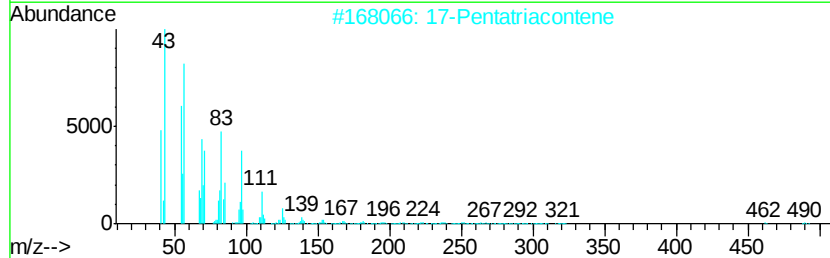
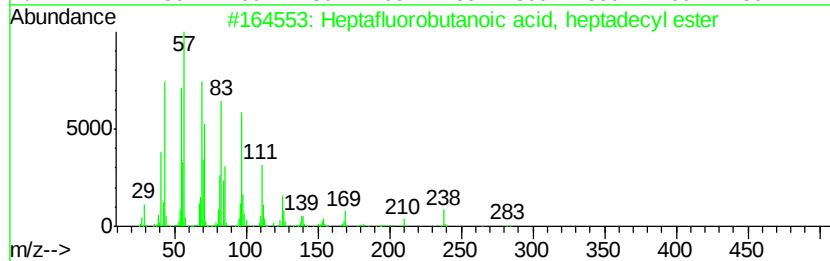
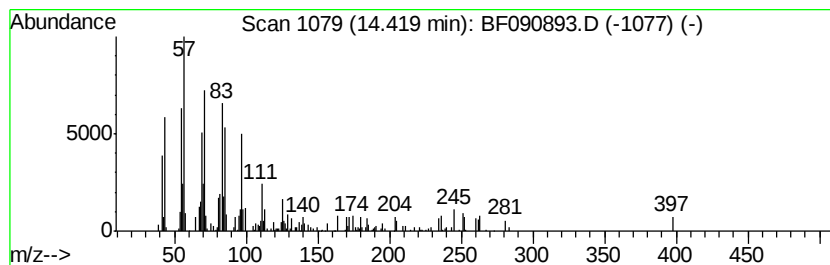
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Heptafluorobutanoic acid, h... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.42	2.35 ng	112599	Chrysene-d12	13.93

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	58
2		17-Pentatriacontene	491	C35H70	006971-40-0	53
3		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	52
4		Tridecane	184	C13H28	000629-50-5	51
5		E-7-Octadecene	252	C18H36	1000130-92-0	46



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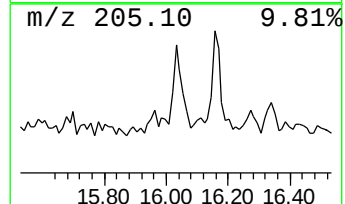
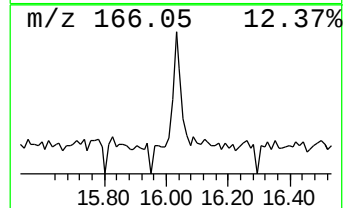
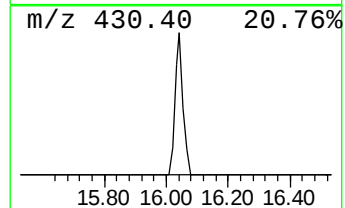
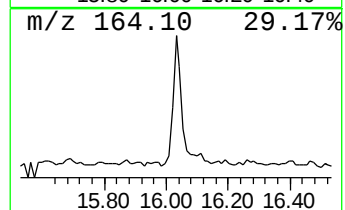
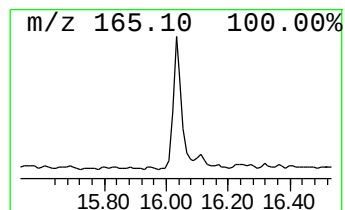
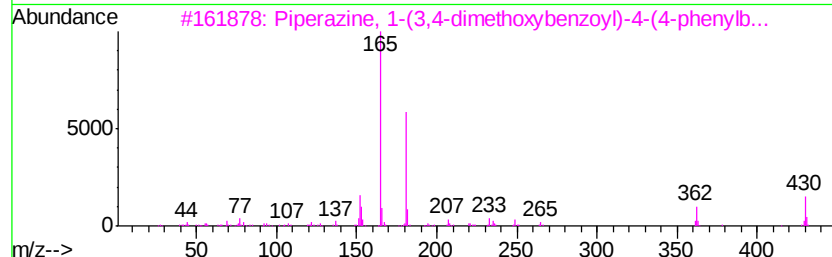
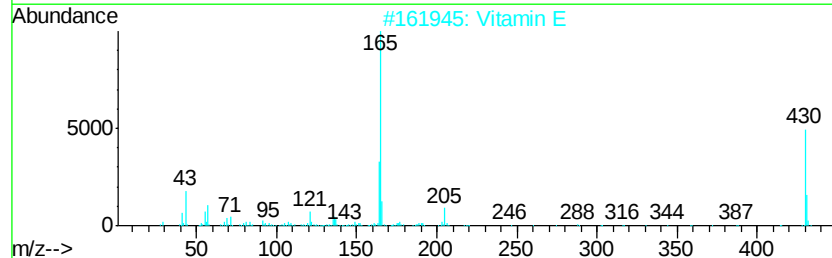
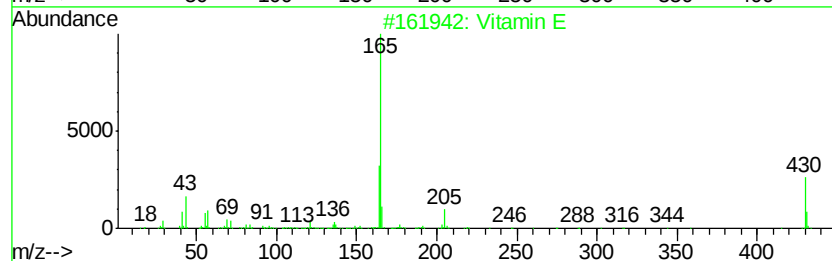
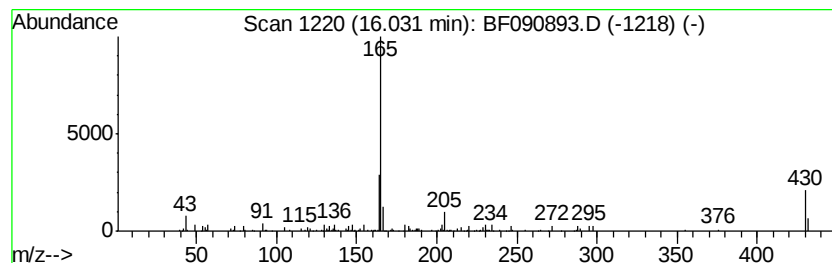
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Vitamin E Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.03	3.09 ng	115433	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Vitamin E	430	C29H50O2	010191-41-0	90
2		Vitamin E	430	C29H50O2	000059-02-9	90
3		Piperazine, 1-(3,4-dimethoxybenz...	430	C26H26N2O4	333756-87-9	45
4		2-Oxo-2-phenylethyl 2,6-dimethox...	300	C17H16O5	283170-58-1	28
5		4,5,6,7-Tetrahydro-benzo[c]thiop...	221	C12H15NOS	1000296-67-7	28



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ClientSampleId :
EME-BLEND-007

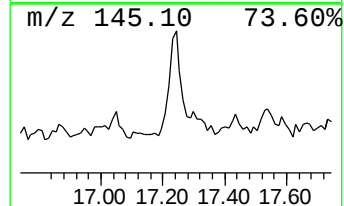
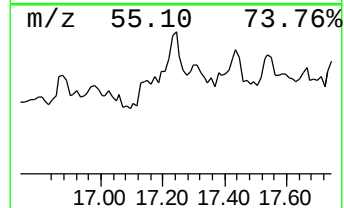
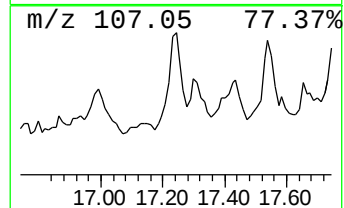
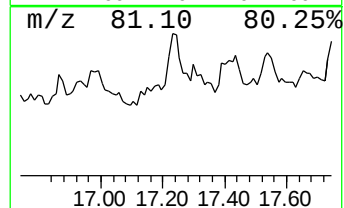
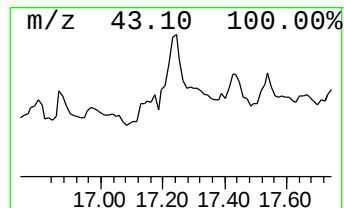
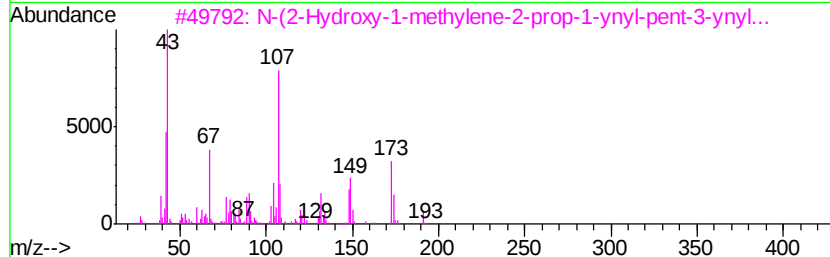
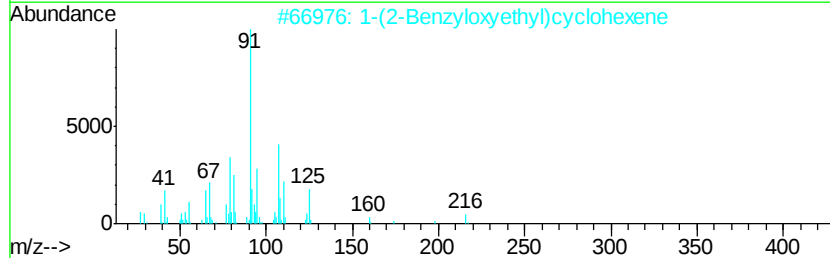
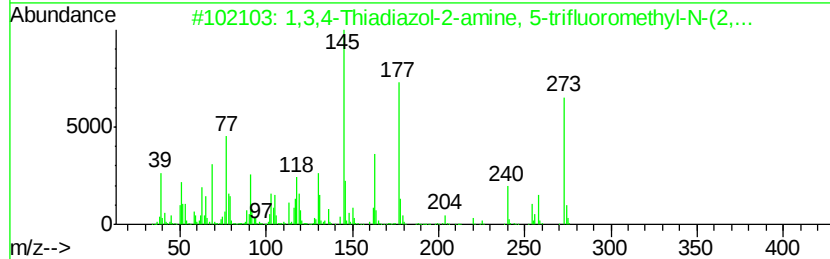
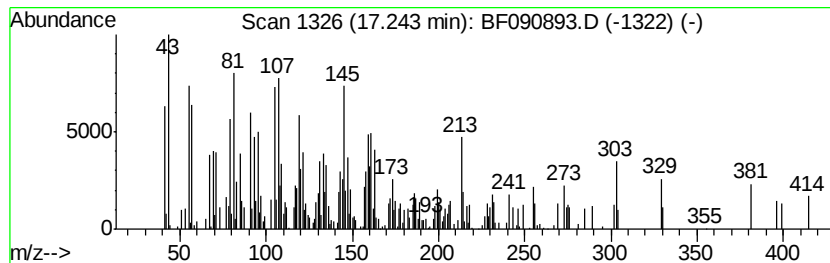
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown17.24 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.24	6.15 ng	229840	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3,4-Thiadiazol-2-amine, 5-trif...	273	C11H10F3N3S	1000267-89-8	38
2		1-(2-Benzoyloxyethyl)cyclohexene	216	C15H20O	082725-62-0	25
3		N-(2-Hydroxy-1-methylene-2-prop-...	191	C11H13NO2	1000194-35-2	25
4		Benzeneacetamide, 4-(oxiranylmeth...	207	C11H13NO3	029122-69-8	25
5		1H-Cycloprop[e]azulene, 1a,2,3,5...	204	C15H24	021747-46-6	15



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102716\
Data File : BF090893.D
Acq On : 28 Oct 2016 3:24
Operator : UM/SJ
Sample : H5430-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EME-BLEND-007

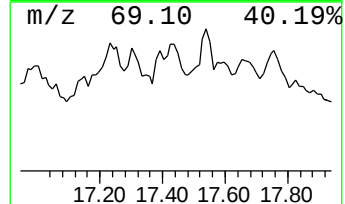
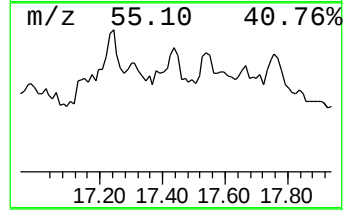
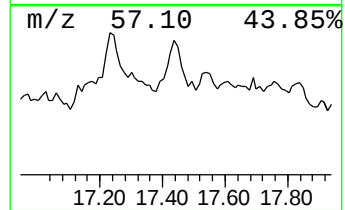
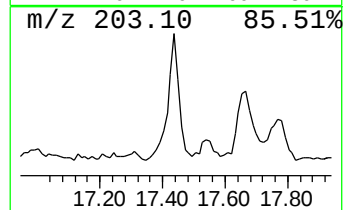
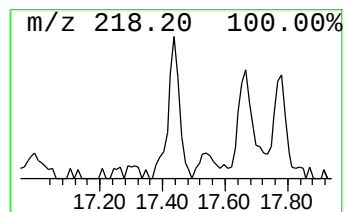
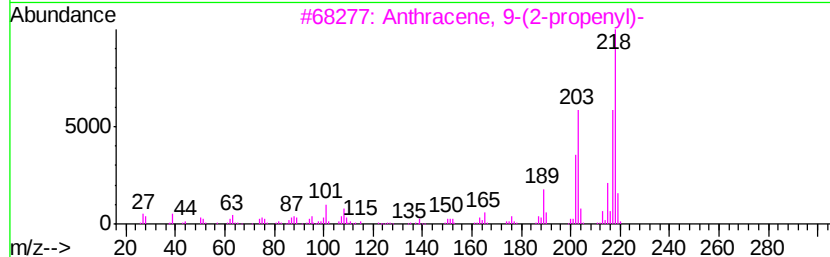
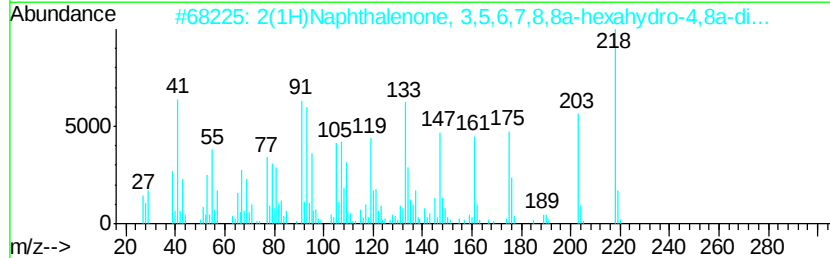
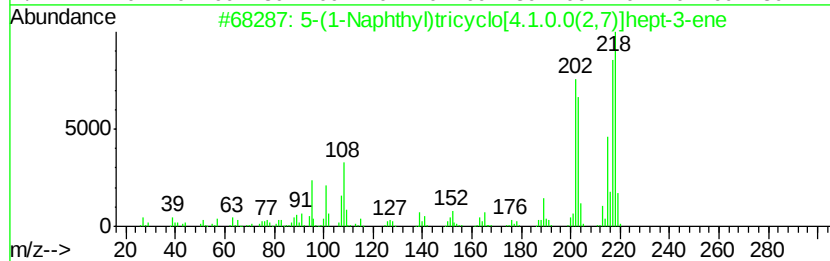
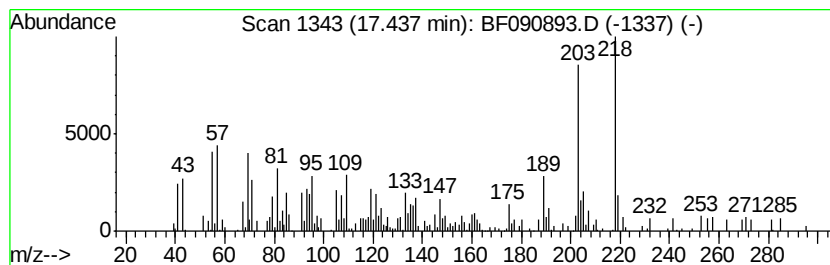
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 5-(1-Naphthyl)tricyclo[4.1.... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.44	5.64 ng	211000	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-(1-Naphthyl)tricyclo[4.1.0.0(2...]	218	C17H14	107729-05-5	60
2		2(1H)Naphthalenone, 3,5,6,7,8,8a...	218	C15H22O	1000188-66-5	49
3		Anthracene, 9-(2-propenyl)-	218	C17H14	023707-65-5	43
4		1-Ethoxy-4-nitroisoquinoline	218	C11H10N2O3	103863-04-3	43
5		8-Amino-6,7-dimethoxy-4-methylqu...	218	C12H14N2O2	1000213-07-2	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF102716\
Data File : BF090893.D
Acq On : 28 Oct 2016 3:24
Operator : UM/SJ
Sample : H5430-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EME-BLEND-007

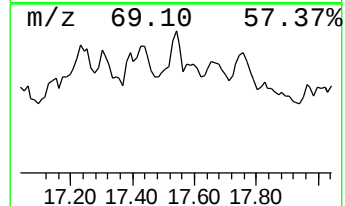
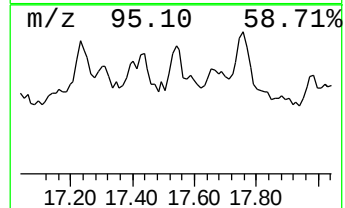
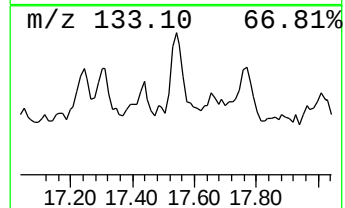
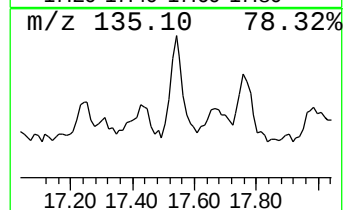
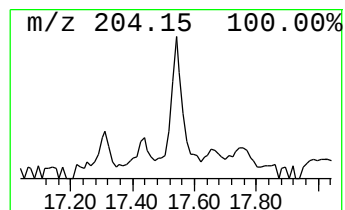
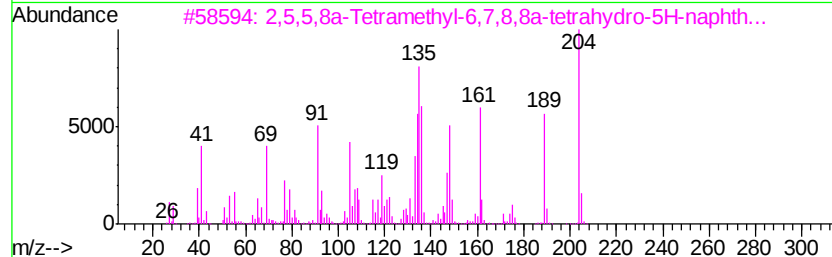
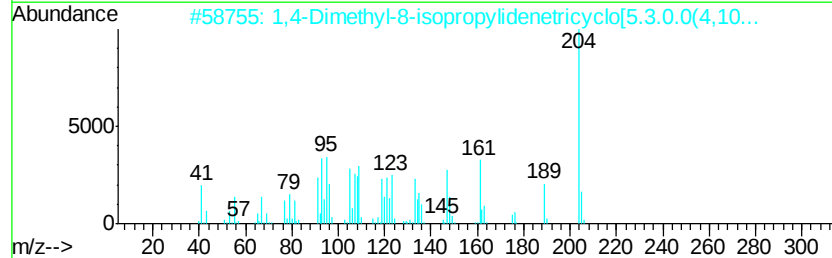
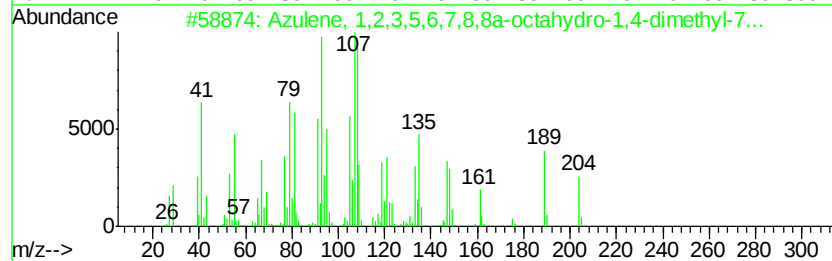
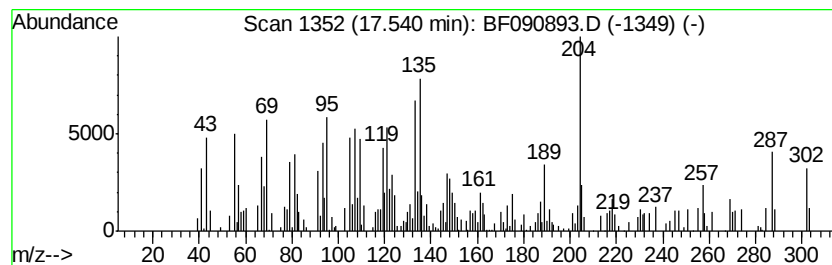
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Azulene, 1,2,3,5,6,7,8,8a-o... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.54	4.42 ng	165283	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Azulene, 1,2,3,5,6,7,8,8a-octahy...	204	C15H24	003691-11-0	64
2		1,4-Dimethyl-8-isopropylidenetri...	204	C15H24	1000140-07-7	52
3		2,5,5,8a-Tetramethyl-6,7,8,8a-te...	204	C14H200	124957-09-1	50
4		Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	004630-07-3	49
5		Coumarin, 3,4-dihydro-4,4,6,8-te...	204	C13H1602	249629-60-5	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102716\
Data File : BF090893.D
Acq On : 28 Oct 2016 3:24
Operator : UM/SJ
Sample : H5430-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

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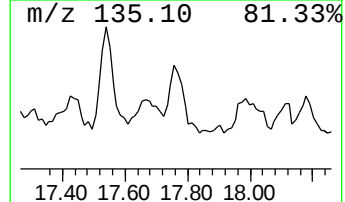
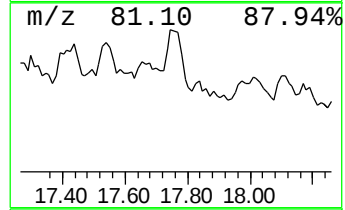
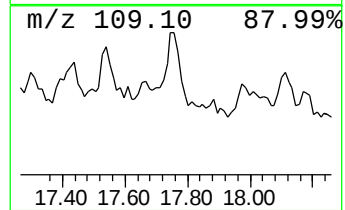
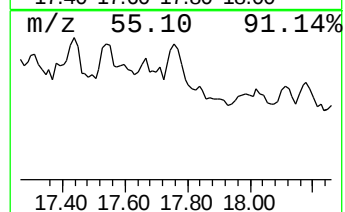
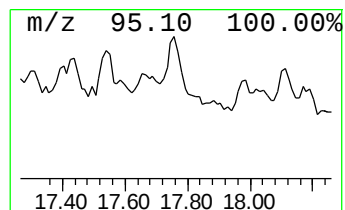
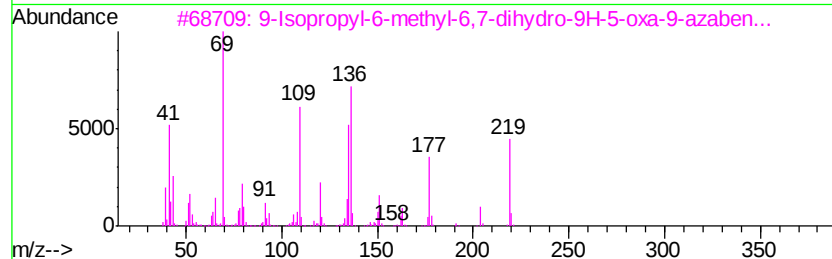
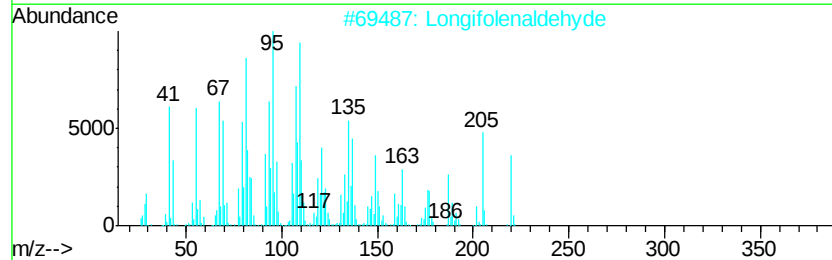
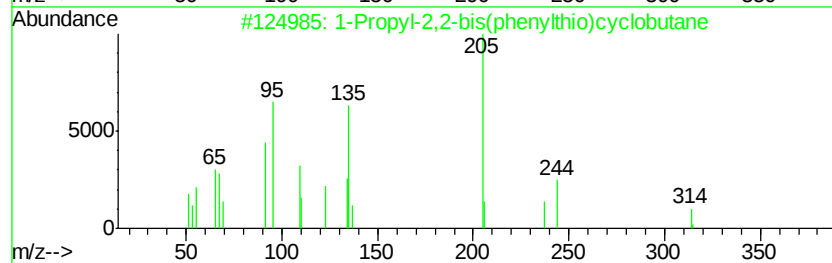
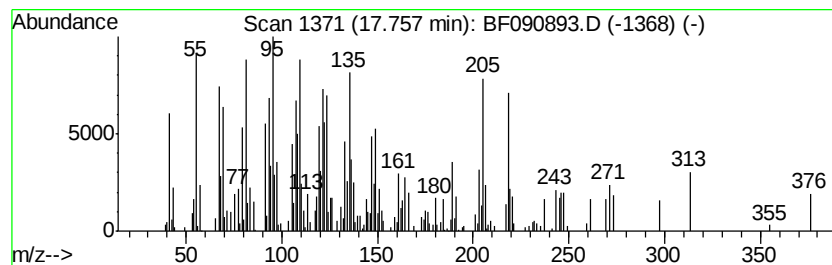
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown17.76 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.76	7.78 ng	290748	Perylene-d12	15.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Propyl-2,2-bis(phenylthio)cyclobutane	314	C19H22S2	122732-89-2	38
2			Longifolenaldehyde	220	C15H24O	019890-84-7	35
3			9-Isopropyl-6-methyl-6,7-dihydro-9H-5-oxa-9-azabenzo[5,6-b]pyridine	219	C13H17N02	134076-78-1	35
4			6.beta.Bicyclo[4.3.0]nonane, 5.beta.	332	C15H25I	1000195-85-9	27
5			Androst-5-en-3-ol, 4,4-dimethyl-	302	C21H34O	007673-17-8	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF102716\
Data File : BF090893.D
Acq On : 28 Oct 2016 3:24
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Misc :
ALS Vial : 26 Sample Multiplier: 1

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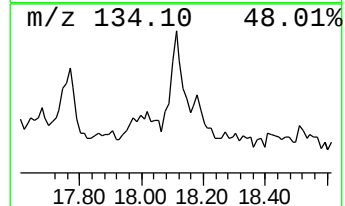
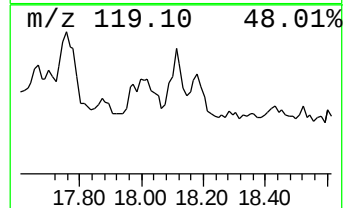
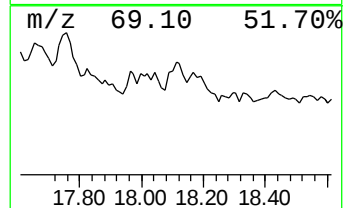
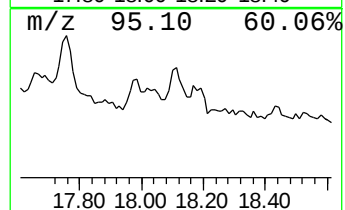
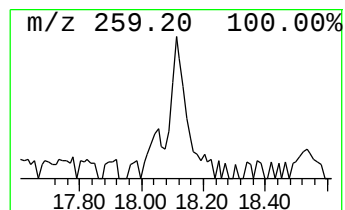
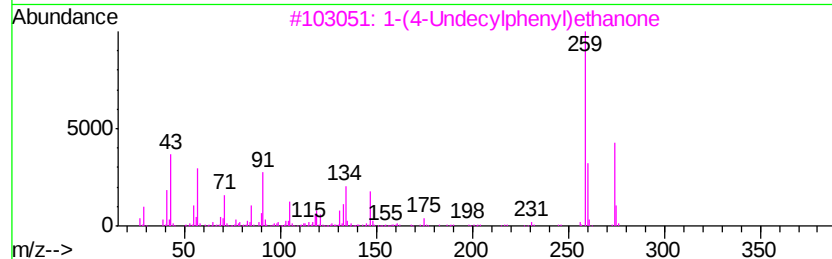
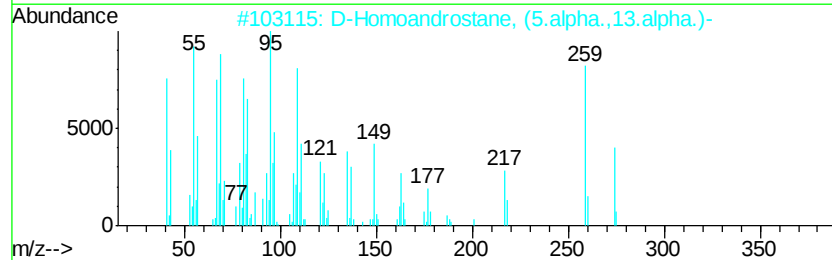
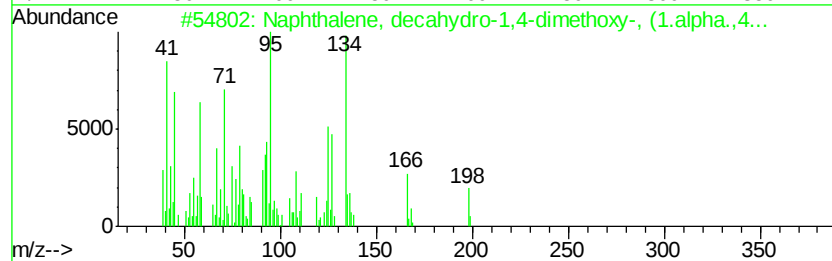
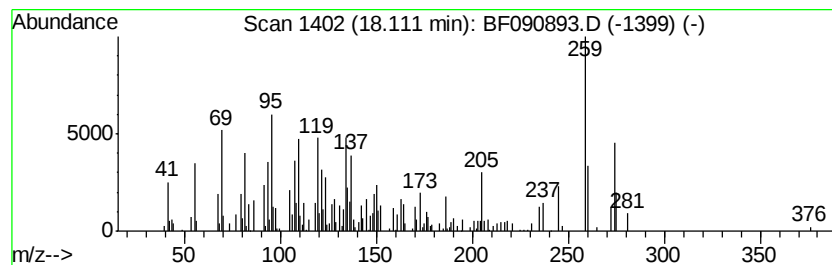
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Naphthalene, decahydro-1,4-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.11	2.63 ng	98292	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, decahydro-1,4-dimet...	198	C12H22O2	042067-56-1	56
2		D-Homoandrostane, (5.alpha.,13.a...	274	C20H34	054482-31-4	49
3		1-(4-Undecylphenyl)ethanone	274	C19H30O	1000185-92-4	30
4		Androstan-6-one, (5.alpha.)-	274	C19H30O	003676-06-0	25
5		1,3-Diethyl-2-hydroxybenzothiazo...	274	C14H14N2O2S	1000227-15-3	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF102716\
Data File : BF090893.D
Acq On : 28 Oct 2016 3:24
Operator : UM/SJ
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Misc :
ALS Vial : 26 Sample Multiplier: 1

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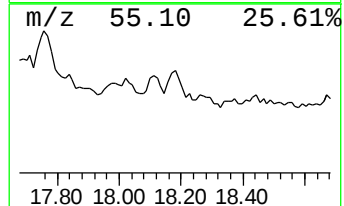
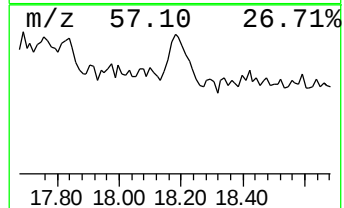
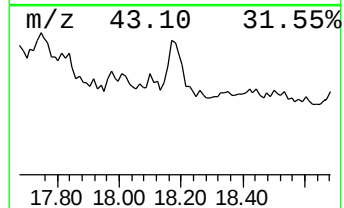
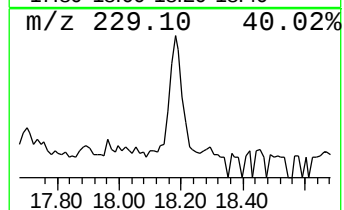
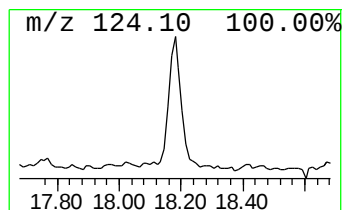
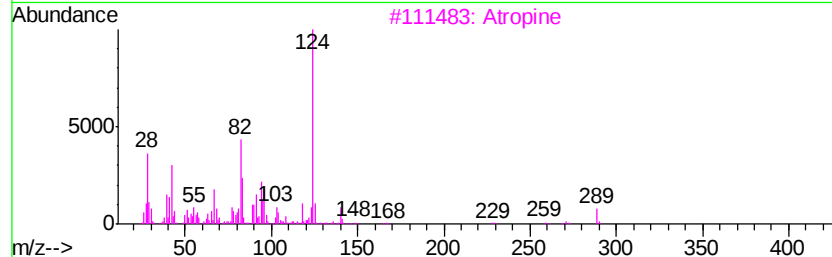
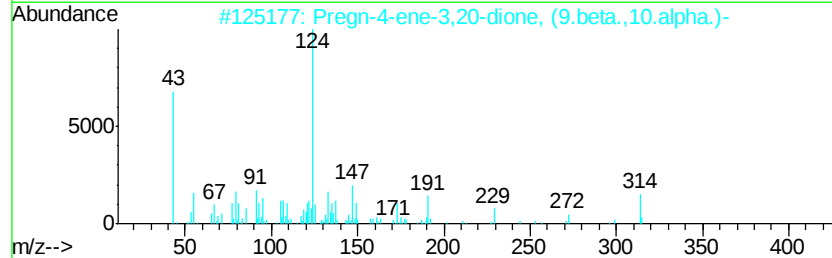
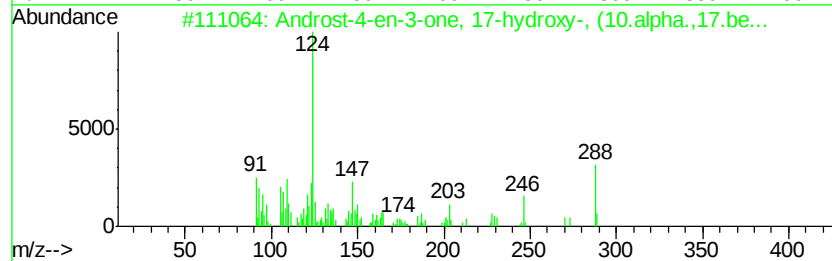
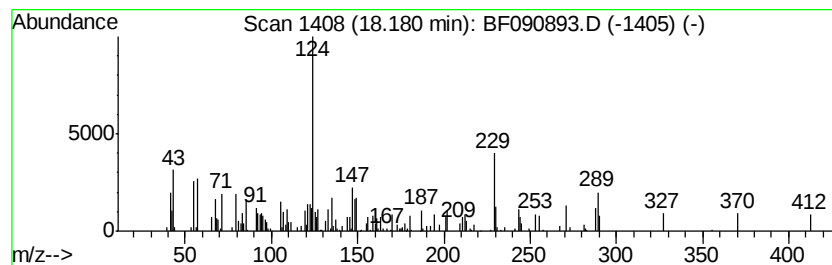
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown18.18 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.18	4.86 ng	181576	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Androst-4-en-3-one, 17-hydroxy-, ...	288	C19H28O2	000604-39-7	46
2		Pregn-4-ene-3,20-dione, (9.beta....	314	C21H30O2	002755-10-4	38
3		Atropine	289	C17H23NO3	000051-55-8	38
4		Tropine 2,4-dichlorophenylacetate	327	C16H19Cl2NO2	1000212-39-1	38
5		Hyoscyamine	289	C17H23NO3	000101-31-5	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102716\
Data File : BF090893.D
Acq On : 28 Oct 2016 3:24
Operator : UM/SJ
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Misc :
ALS Vial : 26 Sample Multiplier: 1

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ClientSampleId :
EME-BLEND-007

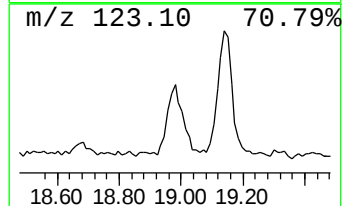
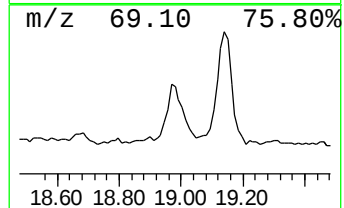
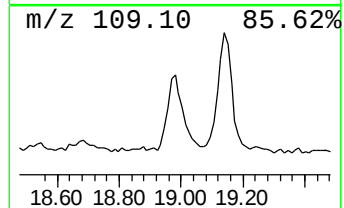
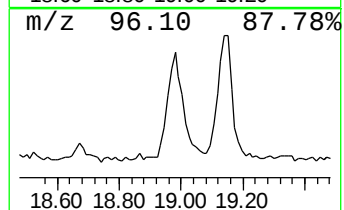
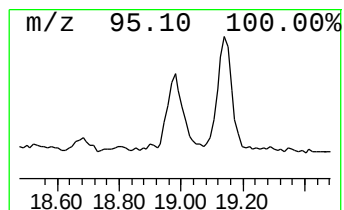
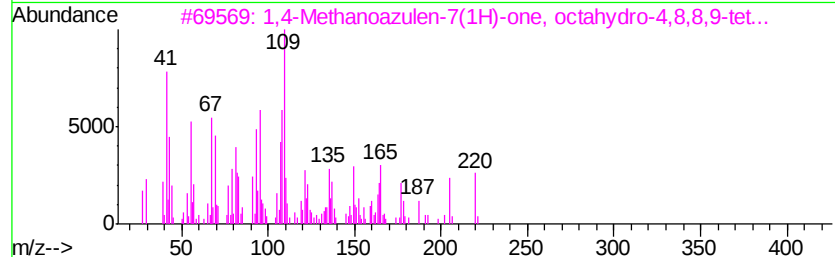
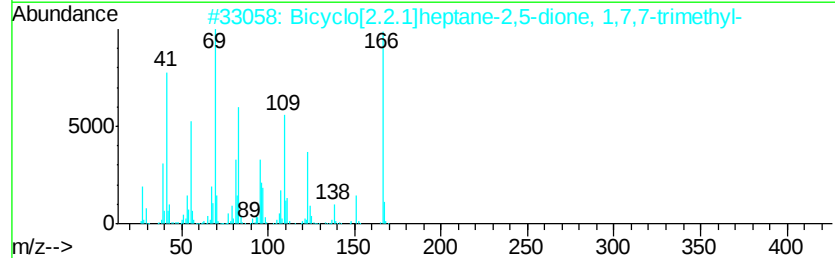
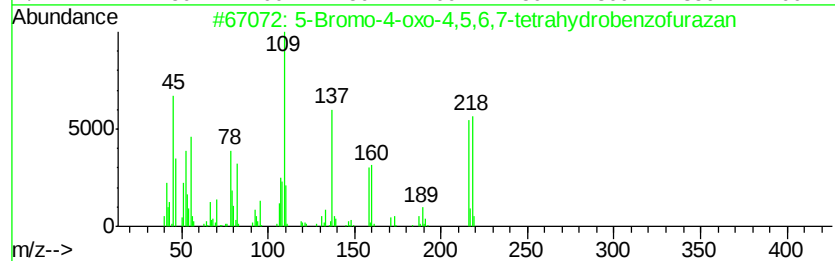
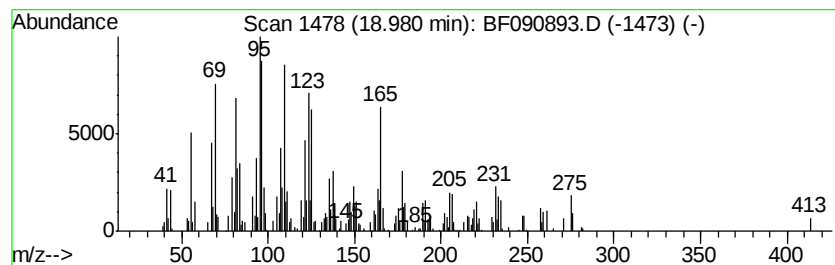
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 5-Bromo-4-oxo-4,5,6,7-tetra... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.98	12.53 ng	468541	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	90
2		Bicyclo[2.2.1]heptane-2,5-dione,...	166	C10H14O2	004230-32-4	45
3		1,4-Methanoazulen-7(1H)-one, oct...	220	C15H24O	018319-28-3	44
4		Bicyclo[2.2.2]octan-1-amine	125	C8H15N	001193-42-6	42
5		Naphthalene, decahydro-1,2-dimet...	166	C12H22	003604-14-6	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF102716\
Data File : BF090893.D
Acq On : 28 Oct 2016 3:24
Operator : UM/SJ
Sample : H5430-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EME-BLEND-007

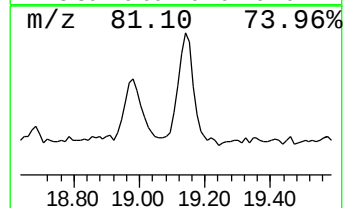
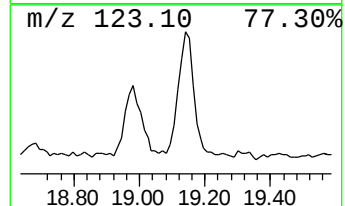
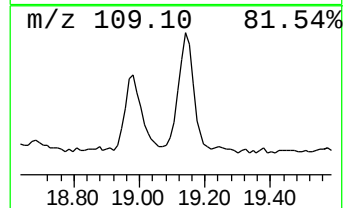
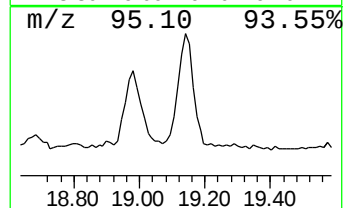
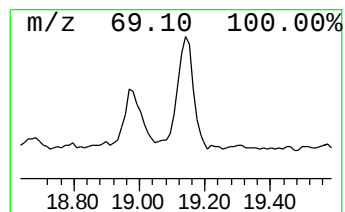
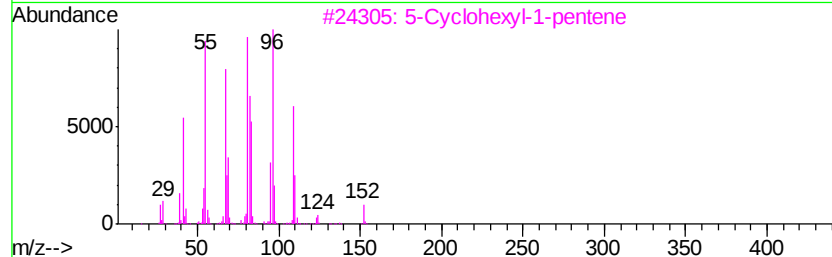
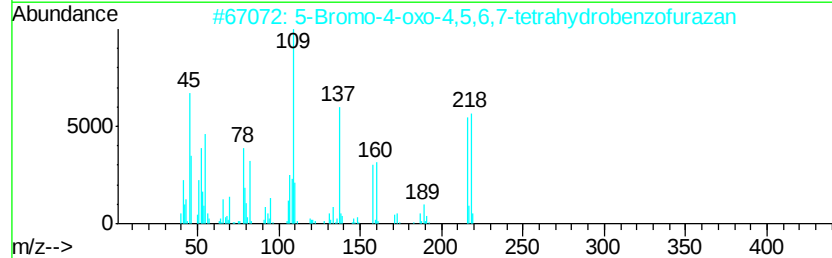
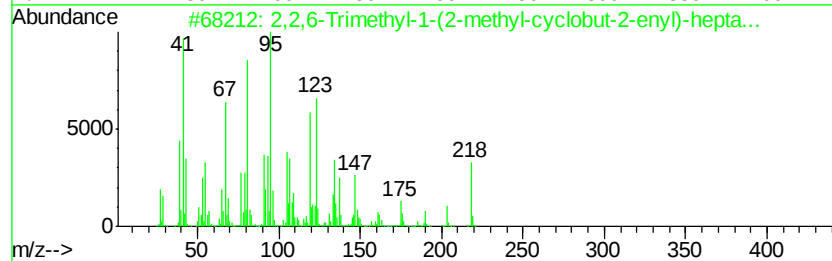
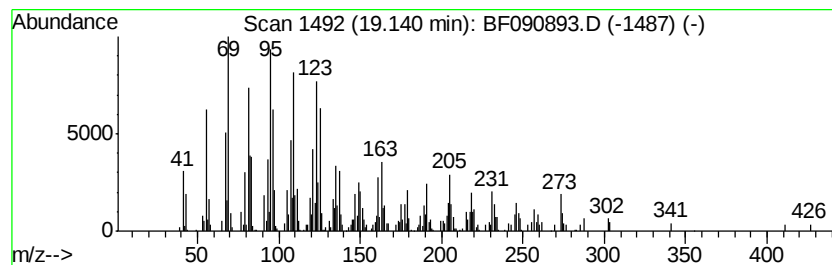
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 2,2,6-Trimethyl-1-(2-methyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.14	20.80 ng	777896	Perylene-d12	15.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2,6-Trimethyl-1-(2-methyl-cycl...	218	C15H22O	1000188-72-8	64
2		5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	59
3		5-Cyclohexyl-1-pentene	152	C11H20	005729-54-4	56
4		3-Cyclopentylpropionic acid, but...	194	C12H18O2	1000292-46-4	38
5		Germaera-1(10),4,11(13)-trien-12...	232	C15H20O2	000553-21-9	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102716\
 Data File : BF090893.D
 Acq On : 28 Oct 2016 3:24
 Operator : UM/SJ
 Sample : H5430-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EME-BLEND-007

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102616.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...	4.93	11.9	ng	551937	1	6.77	924452	20.0
unknown6.54	6.54	66.7	ng	3082930	1	6.77	924452	20.0
Heptafluorobutyri...	13.79	3.4	ng	164507	5	13.93	956971	20.0
Heptafluorobutano...	14.42	2.4	ng	112599	5	13.93	956971	20.0
Vitamin E	16.03	3.1	ng	115433	6	15.35	747819	20.0
unknown17.24	17.24	6.2	ng	229840	6	15.35	747819	20.0
5-(1-Naphthyl)tri...	17.44	5.6	ng	211000	6	15.35	747819	20.0
Azulene, 1,2,3,5,...	17.54	4.4	ng	165283	6	15.35	747819	20.0
unknown17.76	17.76	7.8	ng	290748	6	15.35	747819	20.0
Naphthalene, deca...	18.11	2.6	ng	98292	6	15.35	747819	20.0
unknown18.18	18.18	4.9	ng	181576	6	15.35	747819	20.0
5-Bromo-4-oxo-4,5...	18.98	12.5	ng	468541	6	15.35	747819	20.0
2,2,6-Trimethyl-1...	19.14	20.8	ng	777896	6	15.35	747819	20.0