

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032818\
 Data File : BF104014.D
 Acq On : 28 Mar 2018 17:11
 Operator : JU/SJ
 Sample : J2087-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JRS-RB201719RT-CV-2530

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-BF032818.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	2.293	30	35	43	rVB	82989	110987	3.96%	0.392%
2	5.210	527	531	541	rBV	55653	92564	3.30%	0.327%
3	5.598	593	597	633	rBV	1352108	2497773	89.13%	8.814%
4	6.598	763	767	787	rBV	1586692	2552321	91.07%	9.006%
5	6.740	787	791	820	rVB	1961497	2802530	100.00%	9.889%
6	6.963	826	829	849	rVB	540751	751847	26.83%	2.653%
7	7.116	852	855	877	rVB	1632288	1894860	67.61%	6.686%
8	7.528	921	925	944	rBV	1219933	1643092	58.63%	5.798%
9	8.245	1043	1047	1053	rBV	878676	928157	33.12%	3.275%
10	8.581	1099	1104	1107	rBV6	29557	51143	1.82%	0.180%
11	8.622	1107	1111	1113	rVV5	27221	43325	1.55%	0.153%
12	8.675	1113	1120	1122	rVV6	36178	89003	3.18%	0.314%
13	8.698	1122	1124	1128	rVB4	47634	57022	2.03%	0.201%
14	8.769	1133	1136	1138	rBV3	30891	36884	1.32%	0.130%
15	8.857	1148	1151	1153	rBV3	40225	57733	2.06%	0.204%
16	8.981	1170	1172	1174	rVB	65354	48551	1.73%	0.171%
17	9.010	1174	1177	1180	rBV3	58397	77433	2.76%	0.273%
18	9.075	1186	1188	1190	rBV2	39873	38870	1.39%	0.137%
19	9.116	1192	1195	1197	rBV4	52038	58056	2.07%	0.205%
20	9.228	1211	1214	1219	rBV3	148967	188474	6.73%	0.665%
21	9.316	1224	1229	1234	rBV	2499526	2585020	92.24%	9.122%
22	9.386	1238	1241	1245	rBV	302328	326057	11.63%	1.151%
23	9.422	1245	1247	1248	rBV	66629	52214	1.86%	0.184%
24	9.628	1279	1282	1283	rBV2	116167	129709	4.63%	0.458%
25	9.704	1292	1295	1298	rBV3	1080749	1048118	37.40%	3.698%
26	9.733	1298	1300	1302	rVB2	172446	134239	4.79%	0.474%
27	9.757	1302	1304	1307	rBV2	220637	196260	7.00%	0.693%
28	9.792	1308	1310	1311	rBV2	145061	109508	3.91%	0.386%
29	9.863	1318	1322	1324	rBV2	739293	731729	26.11%	2.582%
30	9.910	1327	1330	1333	rVB2	1164740	1096969	39.14%	3.871%
31	9.951	1335	1337	1339	rBV2	202605	210183	7.50%	0.742%
32	9.980	1339	1342	1344	rVV	529482	536368	19.14%	1.893%
33	10.004	1344	1346	1348	rVB2	741931	581065	20.73%	2.050%
34	10.245	1384	1387	1391	rBV3	345789	437964	15.63%	1.545%

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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 >
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35	10.316	1397	1399	1400	rBV2	317533	266432	9.51%	0.940%
36	10.410	1412	1415	1418	rVB2	1159089	1076601	38.42%	3.799%
37	10.804	1479	1482	1485	rBV3	836868	1064560	37.99%	3.756%
38	10.910	1498	1500	1503	rVB3	887508	796596	28.42%	2.811%
39	13.092	1867	1871	1876	rVB	1416142	1509204	53.85%	5.325%
40	14.151	2048	2051	2055	rBV	597445	614994	21.94%	2.170%
41	15.692	2309	2313	2326	rVB	483745	815349	29.09%	2.877%

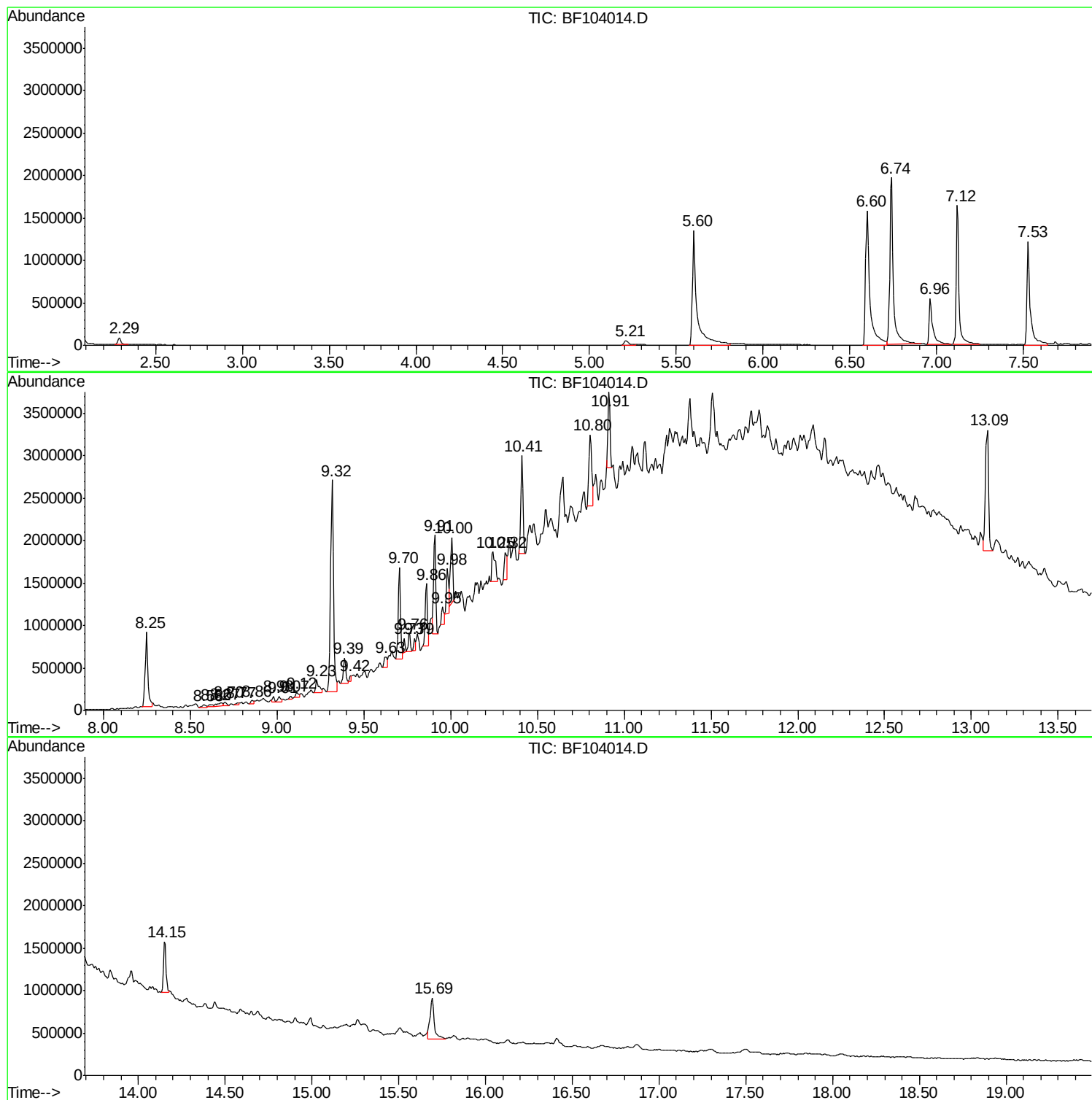
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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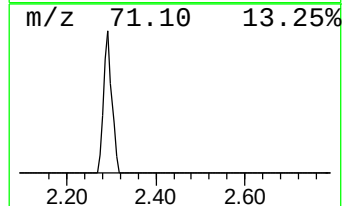
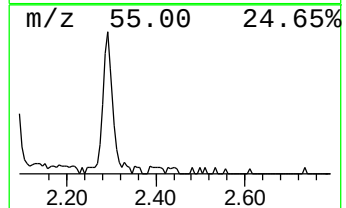
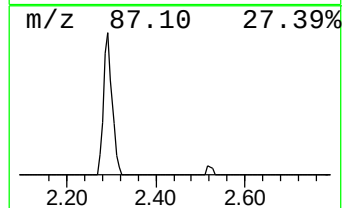
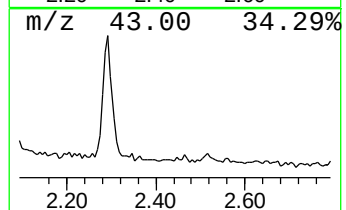
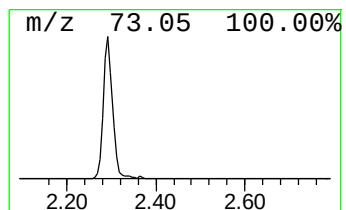
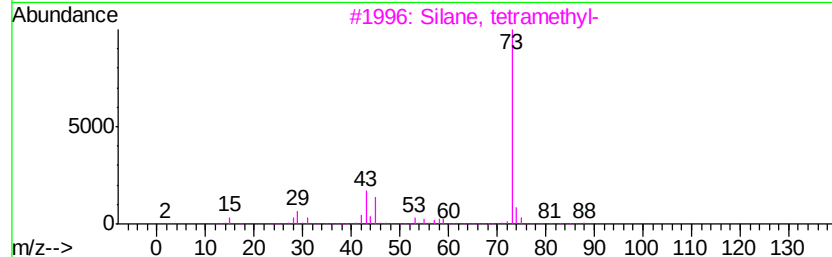
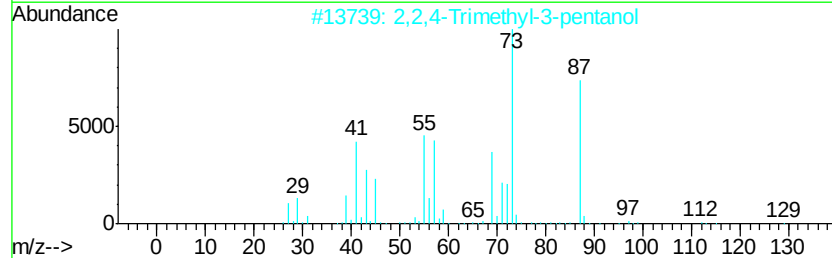
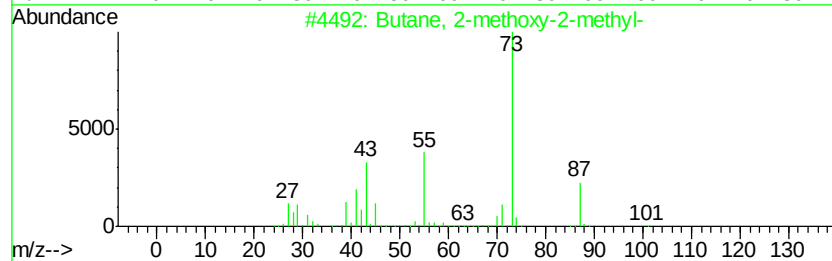
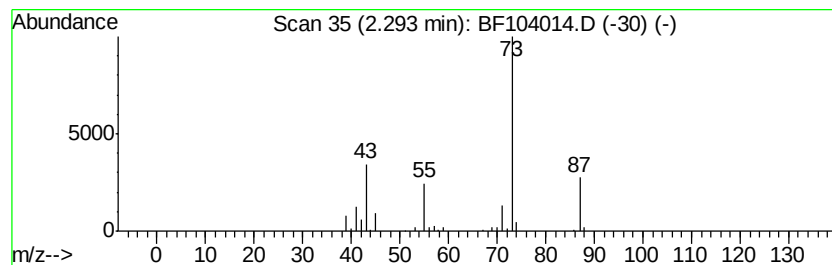
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.29	2.95 ng	110987	1,4-Dichlorobenzene-d4	6.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	23
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5		3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17



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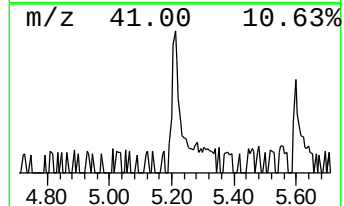
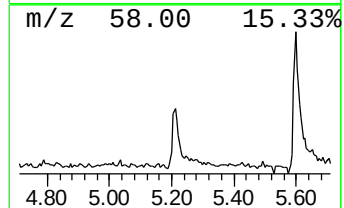
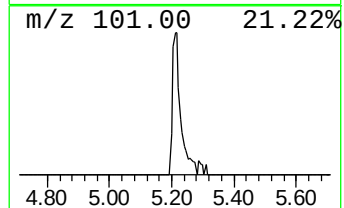
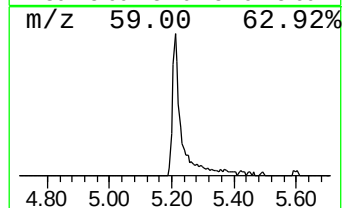
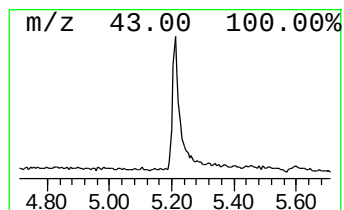
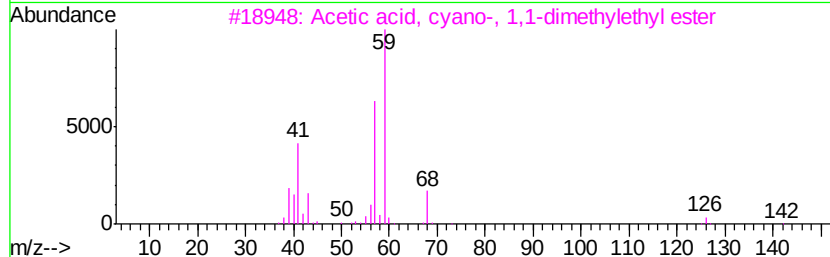
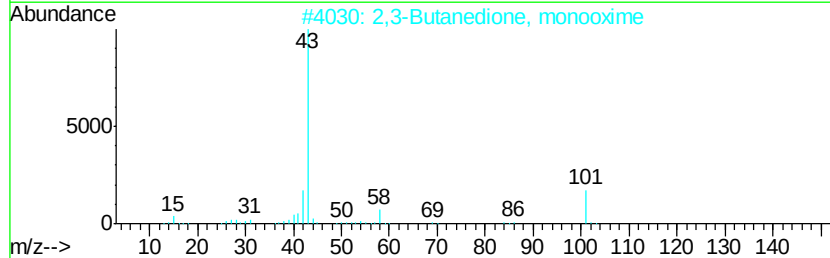
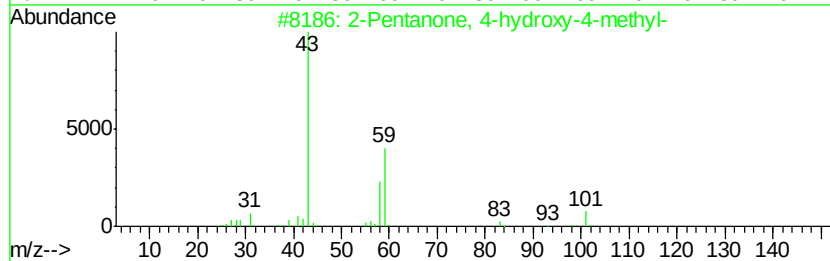
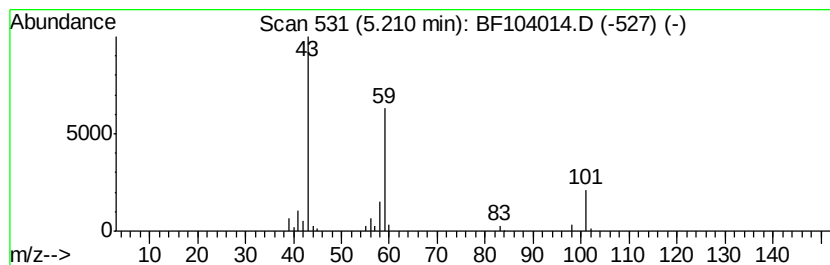
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	2.46 ng	92564	1,4-Dichlorobenzene-d4	6.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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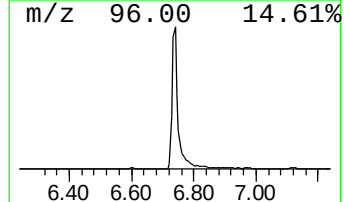
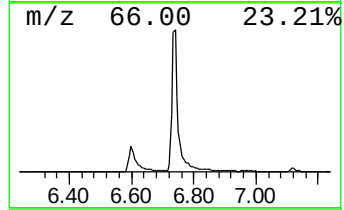
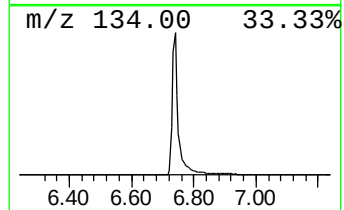
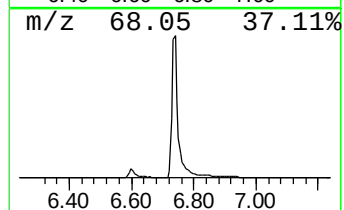
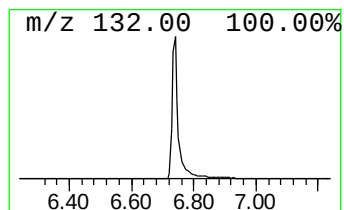
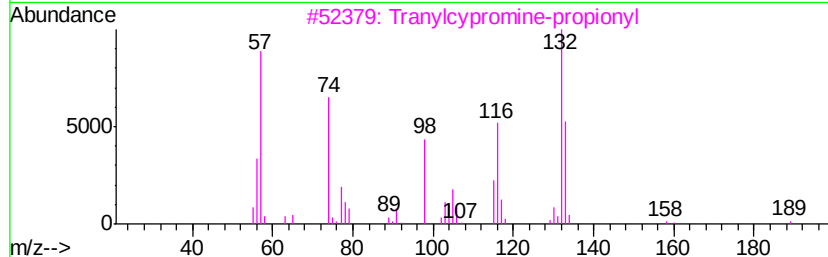
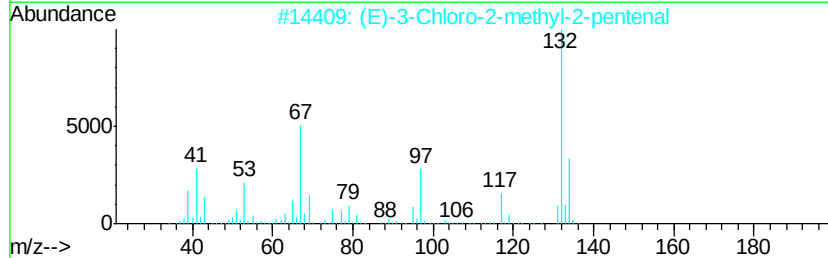
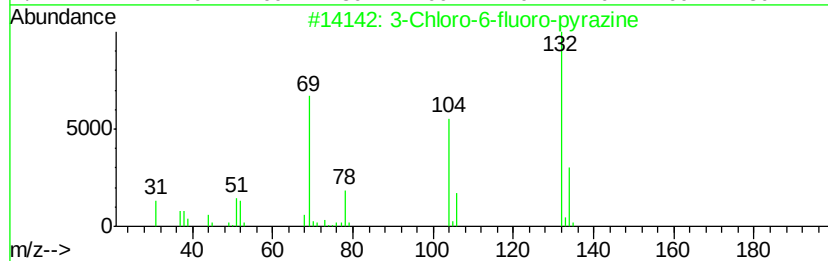
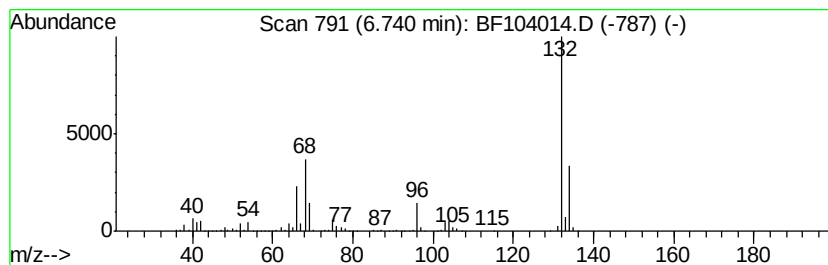
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	74.55 ng	2802530	1,4-Dichlorobenzene-d4	6.96

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	25
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Xenon	132	Xe	007440-63-3	9



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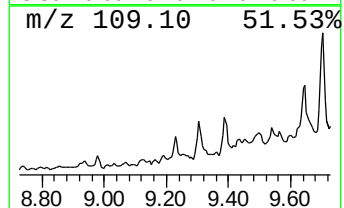
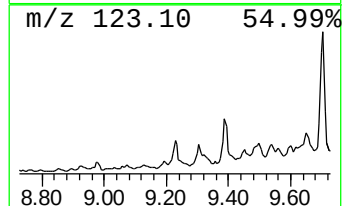
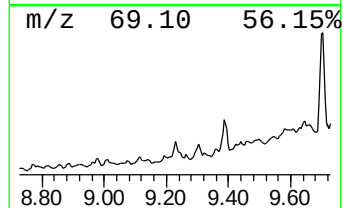
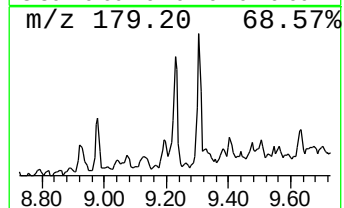
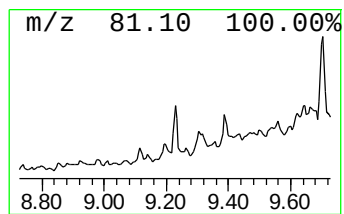
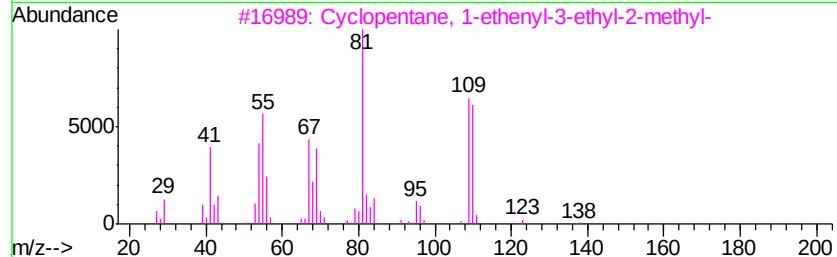
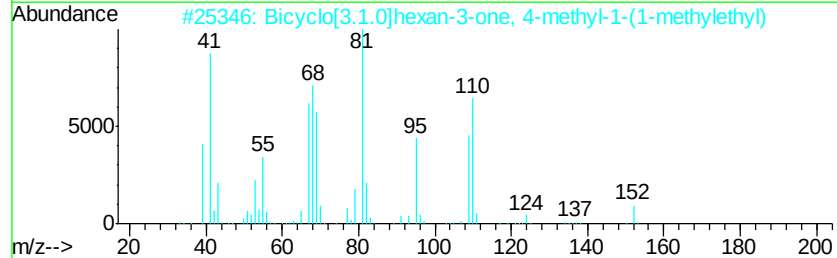
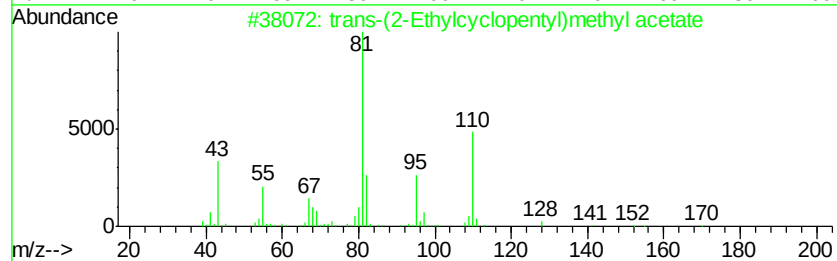
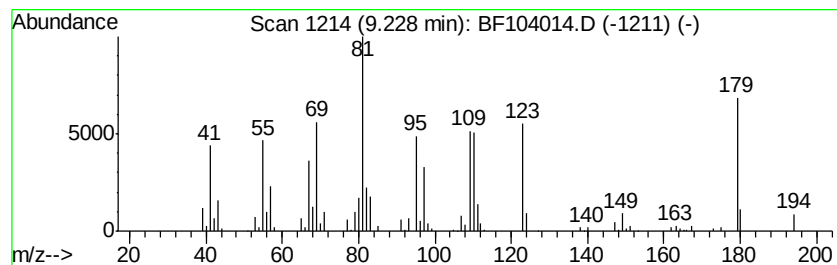
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Peak Number 5 unknown9.23 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.23	6.49 ng	188474	Acenaphthene-d10	10.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	trans-(2-Ethylcyclopentyl)methyl...	170	C10H18O2	1000099-13-9	35
2		Bicyclo[3.1.0]hexan-3-one, 4-met...	152	C10H16O	001125-12-8	32
3		Cyclopentane, 1-ethenyl-3-ethyl-...	138	C10H18	055170-98-4	32
4		4-Ethylcyclohexene	110	C8H14	003742-42-5	30
5		1H-Imidazole, 2-ethyl-4-methyl-	110	C6H10N2	000931-36-2	30



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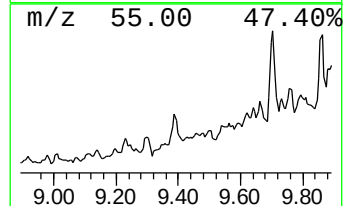
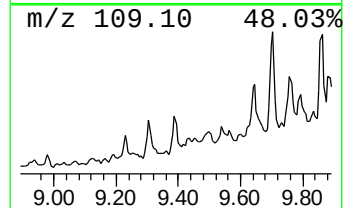
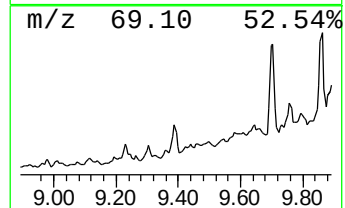
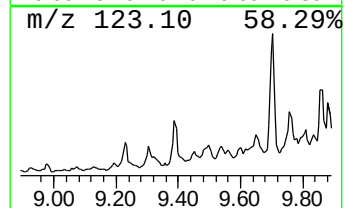
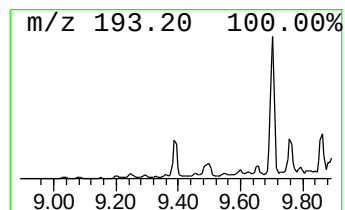
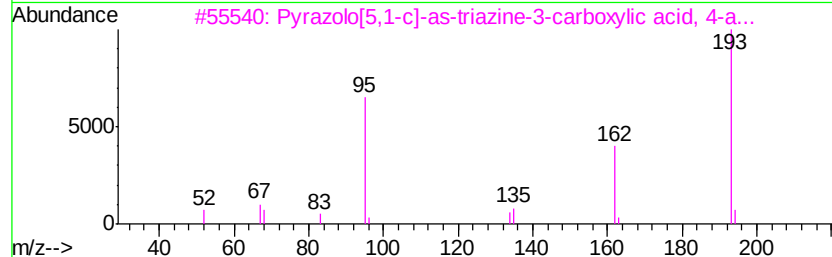
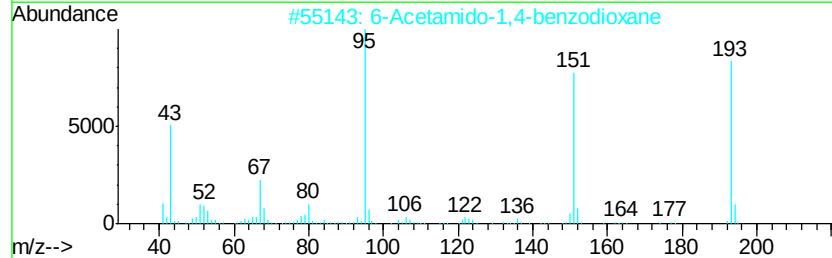
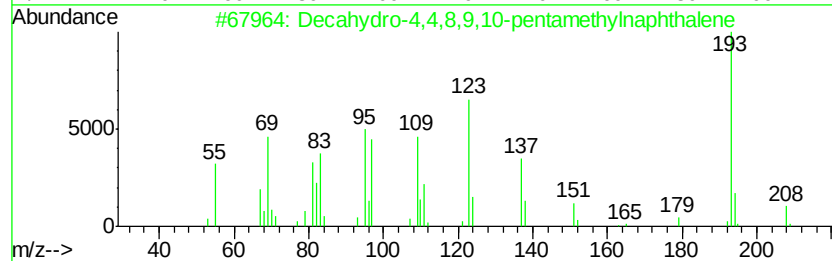
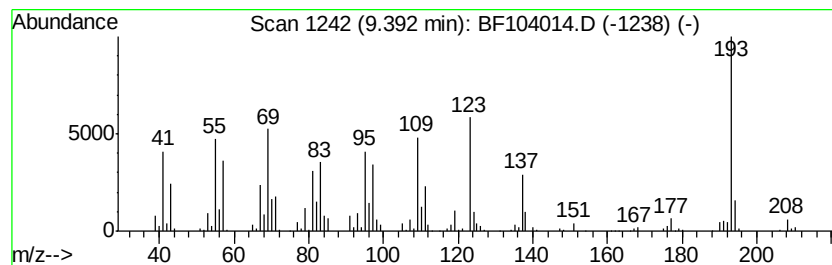
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Peak Number 6 Decahydro-4,4,8,9,10-pentam... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.39	11.22 ng	326057	Acenaphthene-d10	10.00
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Decahydro-4,4,8,9,10-pentamethyl...	208 C15H28	080655-44-3	80
2	6-Acetamido-1,4-benzodioxane	193 C10H11N03	063546-19-0	38
3	Pvrazolo[5,1-c]-as-triazine-3-ca...	193 C6H7N7O	016111-78-7	35
4	Pyrimido(5,4-e)-as-triazine-5,7(...	193 C7H7N5O2	000084-82-2	27
5	2-Anthracenamine	193 C14H11N	000613-13-8	27



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Data File : BF104014.D
Acq On : 28 Mar 2018 17:11
Operator : JU/SJ
Sample : J2087-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
JRS-RB201719RT-CV-2530

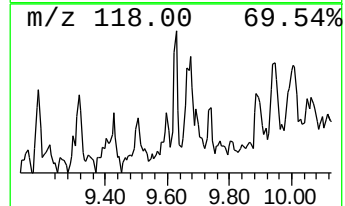
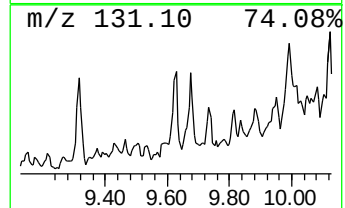
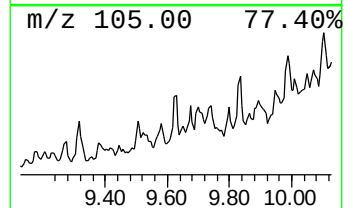
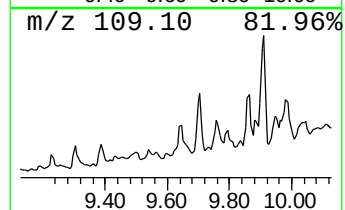
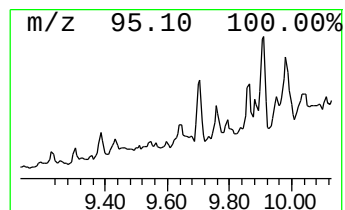
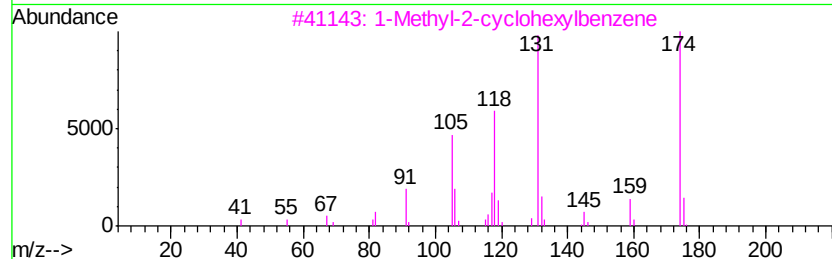
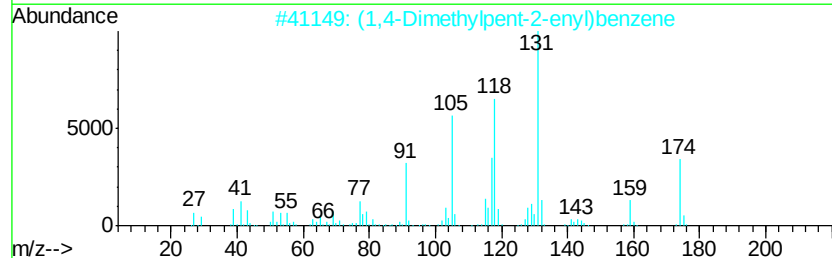
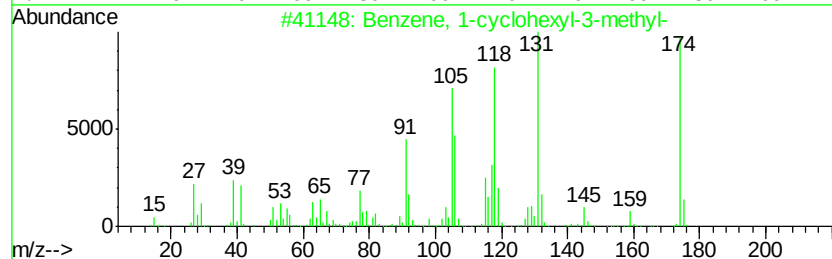
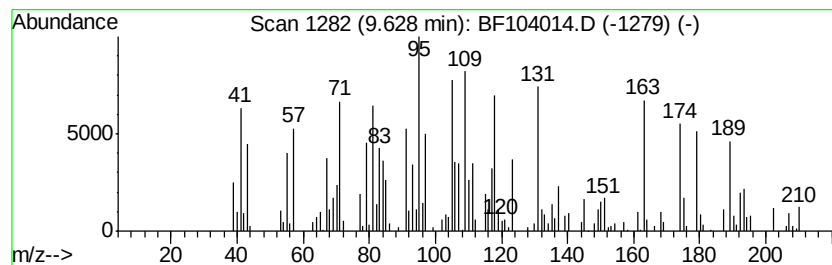
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown9.63 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.63	4.46 ng	129709	Acenaphthene-d10	10.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-cyclohexyl-3-methyl-	174	C13H18	004575-46-6	48
2			(1,4-Dimethylpent-2-enyl)benzene	174	C13H18	1000184-97-9	35
3			1-Methyl-2-cyclohexylbenzene	174	C13H18	004501-35-3	35
4			3-Phenyl-4,4-dimethyl-isoxazol-5...	189	C11H11N02	051942-54-2	25
5			4-Methylimidazole-5-butyric acid	168	C8H12N2O2	1000129-14-7	14



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032818\
Data File : BF104014.D
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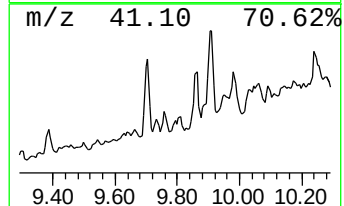
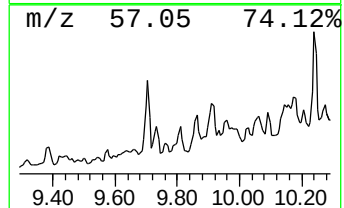
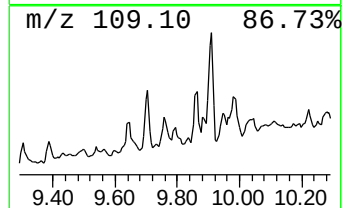
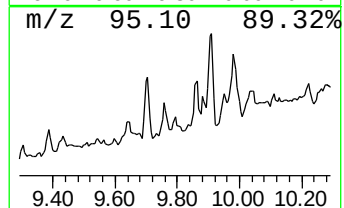
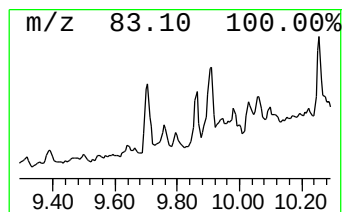
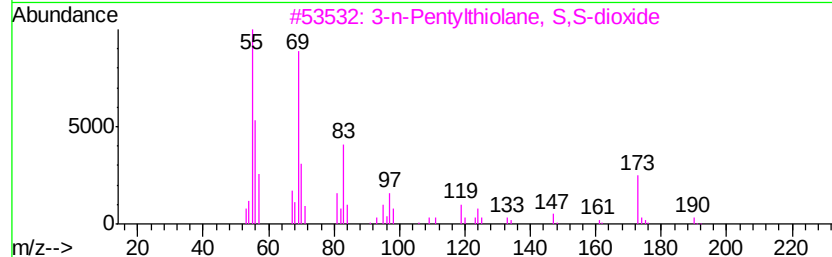
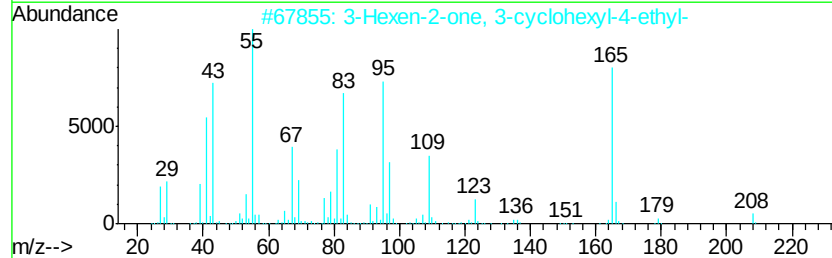
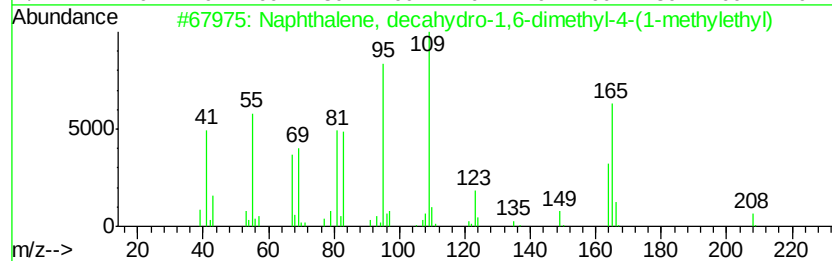
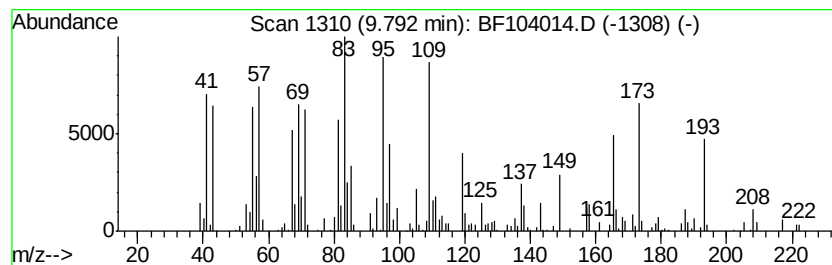
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Naphthalene, decahydro-1,6-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.79	3.77 ng	109508	Acenaphthene-d10	10.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, decahydro-1,6-dimet...	208	C15H28	029788-41-8	50
2		3-Hexen-2-one, 3-cyclohexyl-4-et...	208	C14H24O	1000150-19-1	43
3		3-n-Pentylthiolane, S,S-dioxide	190	C9H18O2S	071053-06-0	27
4		2,6-Heptadienal, 2,4-dimethyl-	138	C9H14O	085136-08-9	25
5		Bicyclo[3.1.1]heptan-3-one, 2,6,...	152	C10H16O	000547-60-4	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032818\
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ClientSampleId :
JRS-RB201719RT-CV-2530

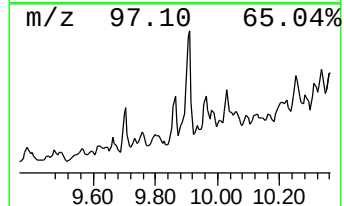
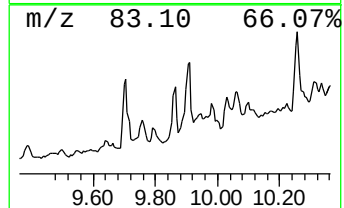
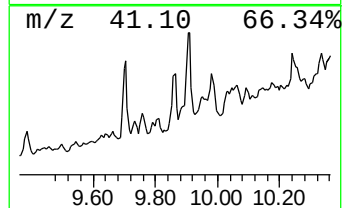
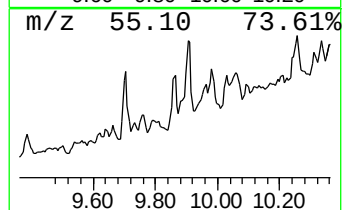
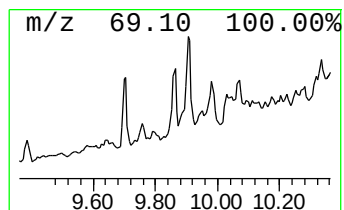
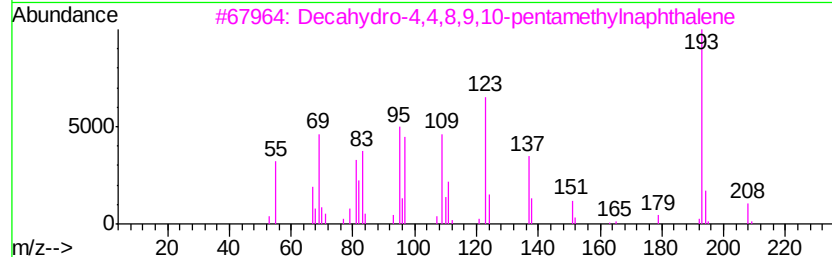
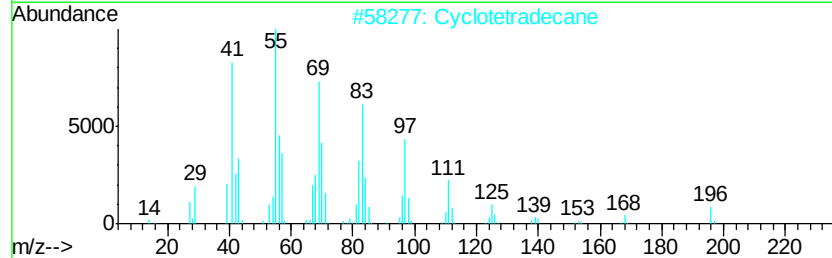
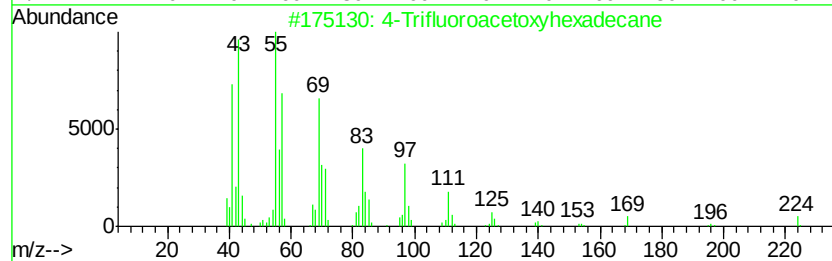
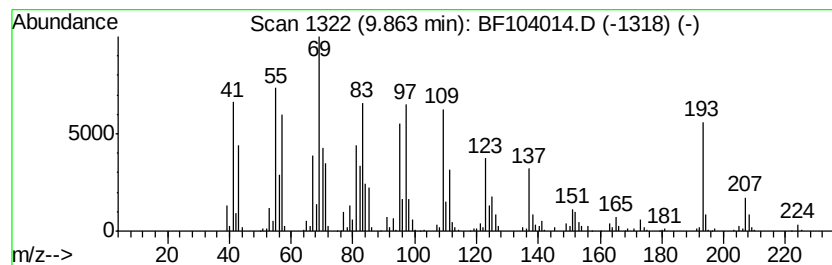
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown9.86 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.86	25.19 ng	731729	Acenaphthene-d10	10.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Trifluoroacetoxylhexadecane	338	C18H33F3O2	1000215-97-5	49
2		Cyclotetradecane	196	C14H28	000295-17-0	45
3		Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	42
4		Silane, chlorodiethyl(2-methylpe...	222	C10H23ClOSi	1000363-79-1	38
5		Pentafluoropropionic acid, tetra...	360	C17H29F5O2	006222-06-6	35



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032818\
Data File : BF104014.D
Acq On : 28 Mar 2018 17:11
Operator : JU/SJ
Sample : J2087-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
JRS-RB201719RT-CV-2530

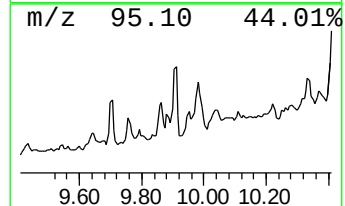
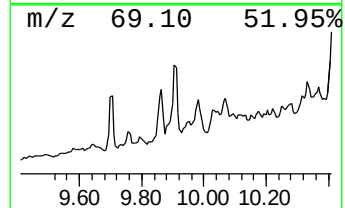
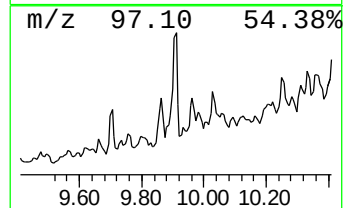
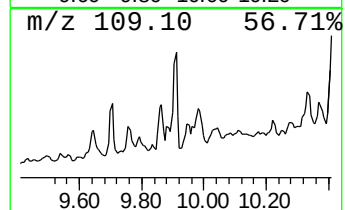
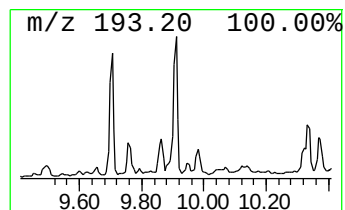
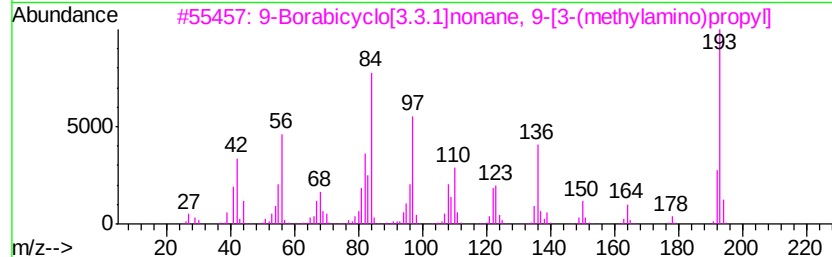
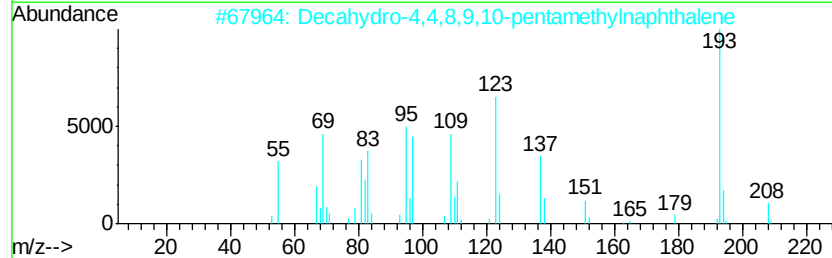
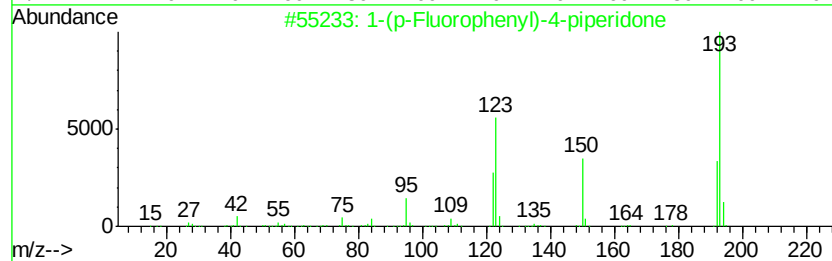
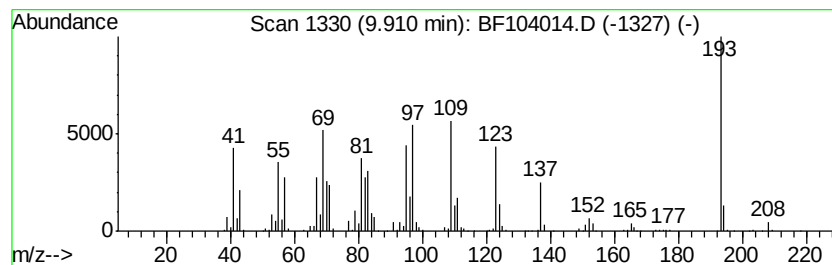
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown9.91 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.91	37.76 ng	1096970	Acenaphthene-d10	10.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	43
2		Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	42
3		9-Borabicyclo[3.3.1]nonane, 9-[3...	193	C12H24BN	1000162-56-0	38
4		Naphthalene, decahydro-1,4a-dime...	208	C15H28	030824-81-8	25
5		[1,1'-Biphenyl]-4-acetonitrile	193	C14H11N	031603-77-7	18



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032818\
Data File : BF104014.D
Acq On : 28 Mar 2018 17:11
Operator : JU/SJ
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Instrument :
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ClientSampleId :
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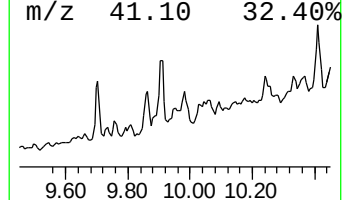
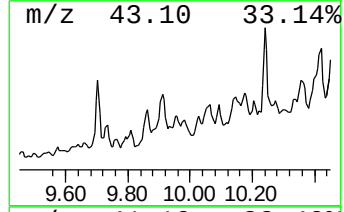
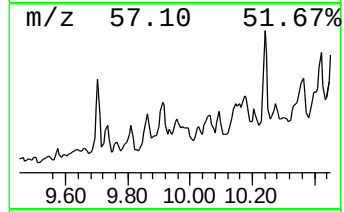
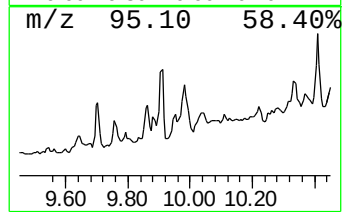
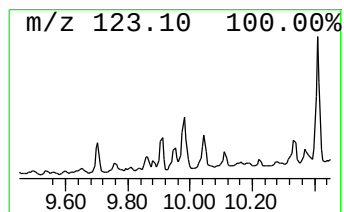
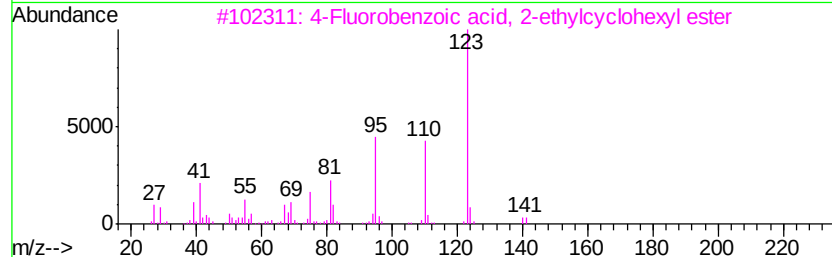
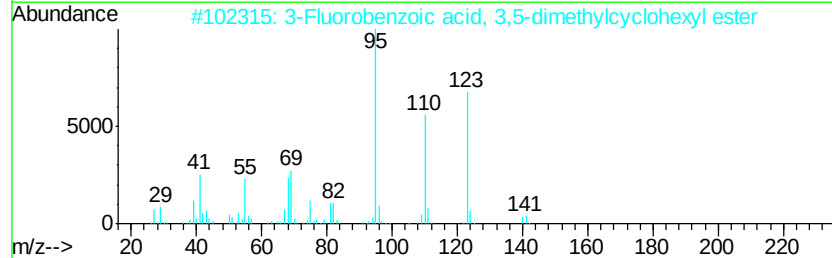
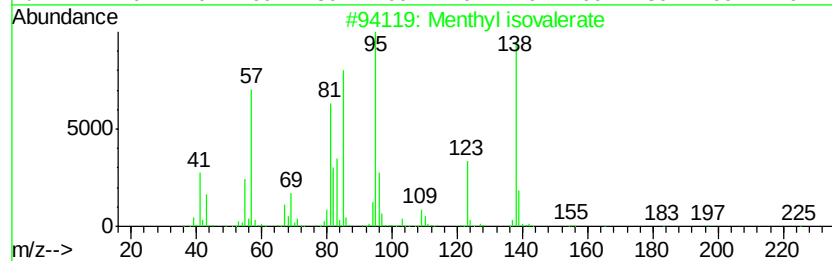
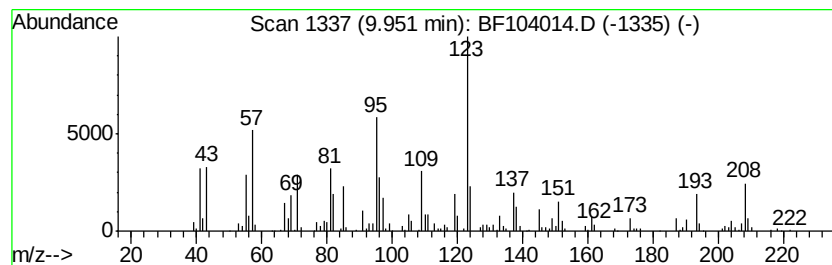
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown9.95 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.95	7.23 ng	210183	Acenaphthene-d10	10.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Menthyl isovalerate	240	C15H28O2	016409-46-4	43
2		3-Fluorobenzoic acid, 3,5-dimeth...	250	C15H19F02	1000279-01-8	38
3		4-Fluorobenzoic acid, 2-ethylcyc...	250	C15H19F02	1000279-06-0	35
4		3,3,5,5-Tetramethylcyclohexanol	156	C10H20O	002650-40-0	27
5		2-Pyrimidinamine, 4,6-dimethyl-	123	C6H9N3	000767-15-7	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032818\
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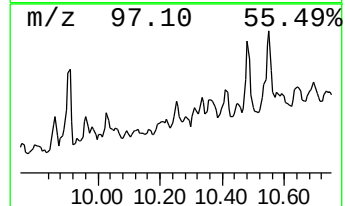
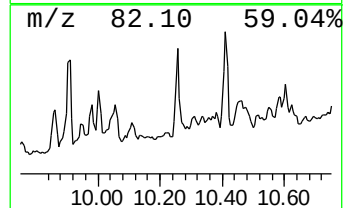
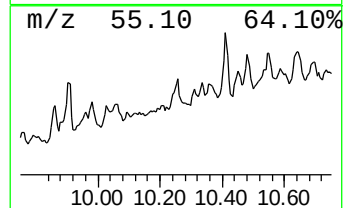
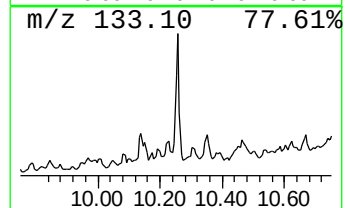
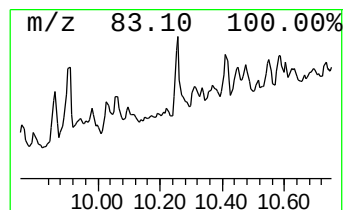
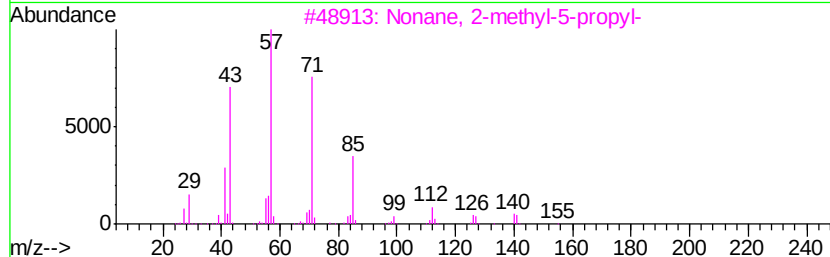
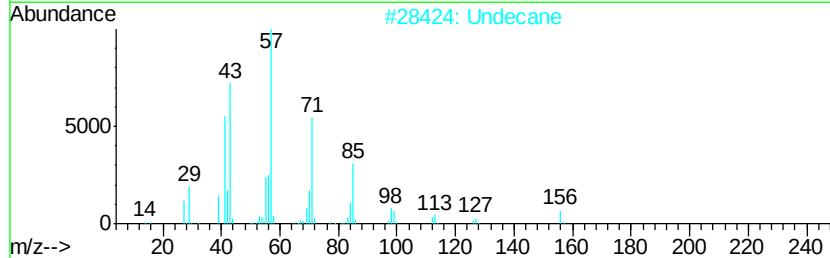
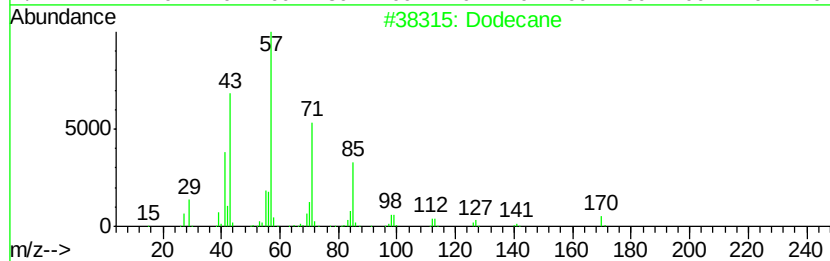
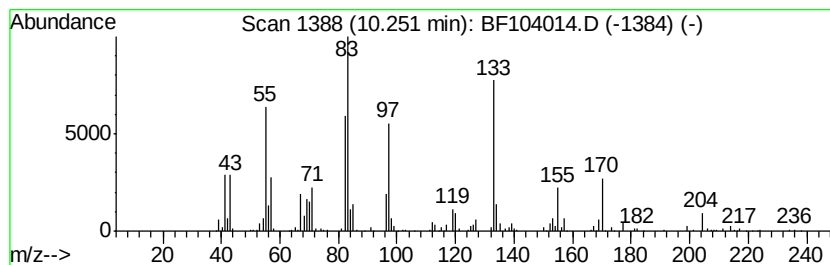
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Dodecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.25	15.07 ng	437964	Acenaphthene-d10	10.00
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Dodecane	170 C12H26	000112-40-3	74
2	Undecane	156 C11H24	001120-21-4	49
3	Nonane, 2-methyl-5-propyl-	184 C13H28	031081-17-1	47
4	Decane, 3-bromo-	220 C10H21Br	030571-71-2	46
5	Heptacosane	380 C27H56	000593-49-7	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032818\
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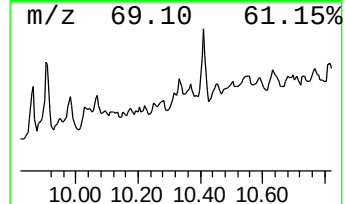
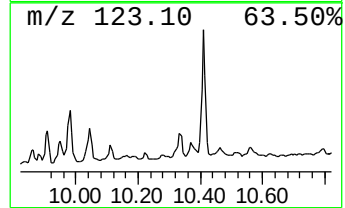
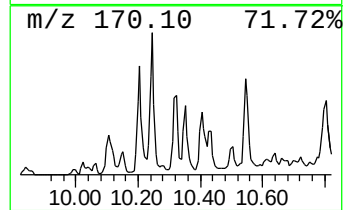
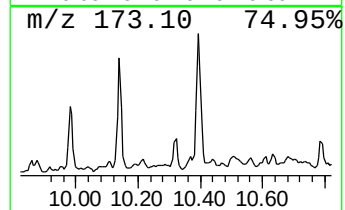
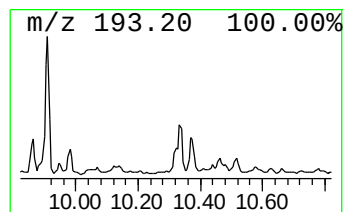
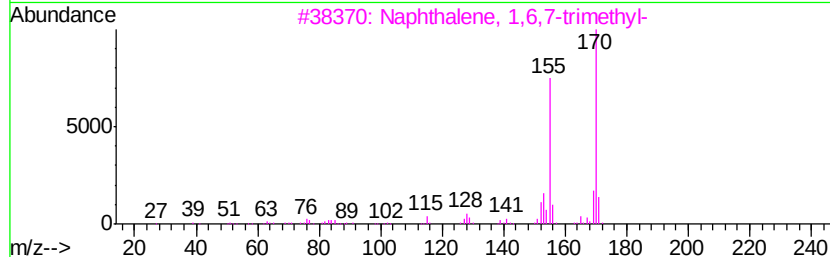
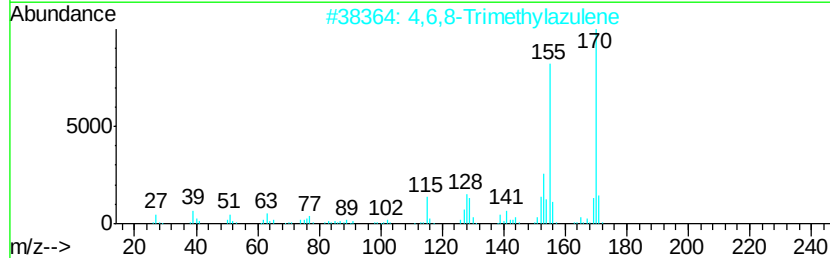
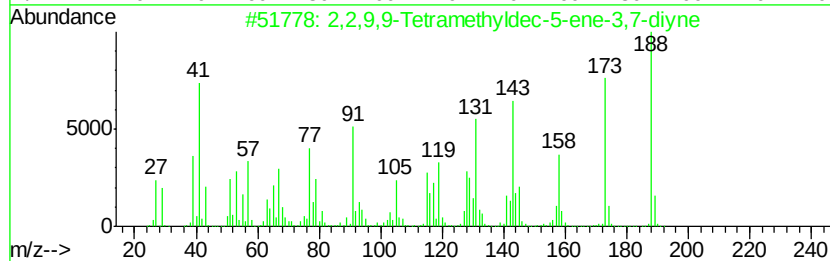
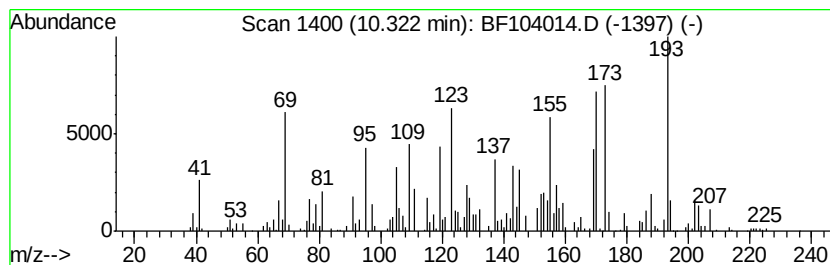
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown10.32 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.32	9.17 ng	266432	Acenaphthene-d10	10.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2,9,9-Tetramethyldec-5-ene-3,7...	188	C14H20	102745-35-7	47		
2	4,6,8-Trimethylazulene	170	C13H14	000941-81-1	35		
3	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	35		
4	Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	35		
5	1H,3H-Thieno[3,4-c]thiophene, 4,...	170	C8H10S2	015441-54-0	25		



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032818\
Data File : BF104014.D
Acq On : 28 Mar 2018 17:11
Operator : JU/SJ
Sample : J2087-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
JRS-RB201719RT-CV-2530

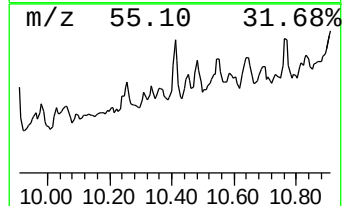
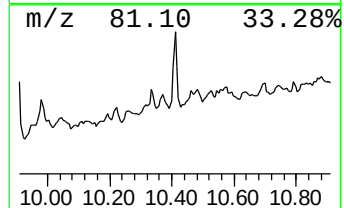
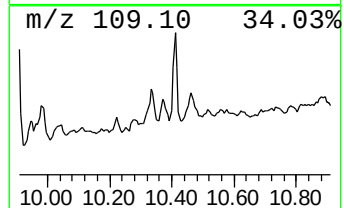
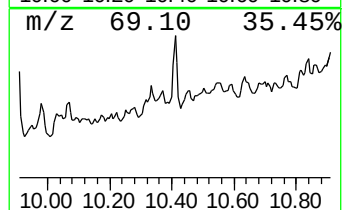
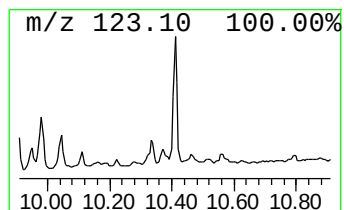
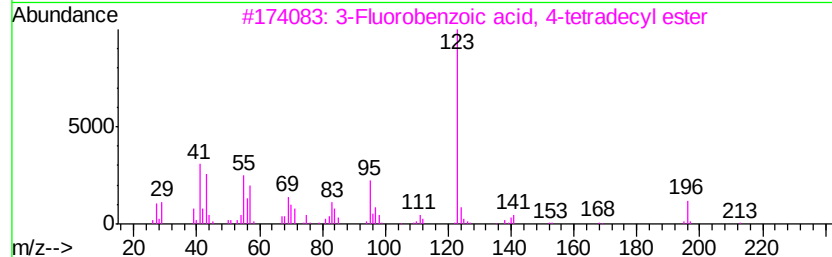
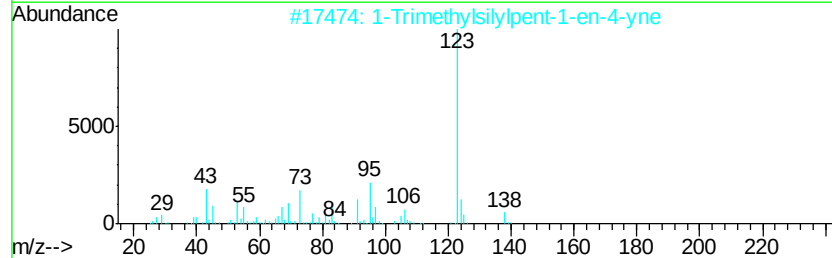
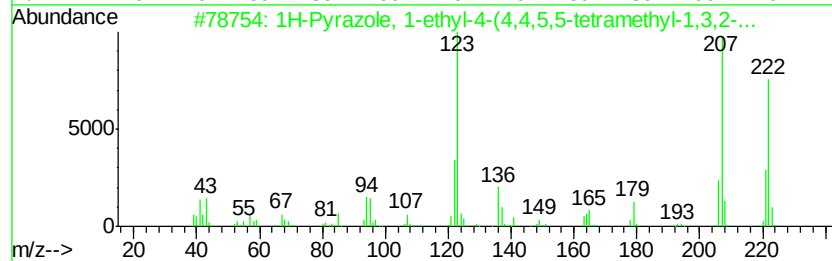
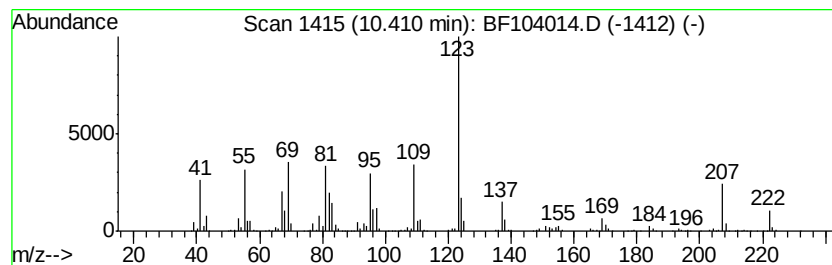
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown10.41 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.41	37.06 ng	1076600	Acenaphthene-d10	10.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Pyrazole, 1-ethyl-4-(4,4,5,5-...	222	C11H19BN2O2	1000319-69-6	53
2		1-Trimethylsilylpent-1-en-4-yne	138	C8H14Si	079516-25-9	52
3		3-Fluorobenzoic acid, 4-tetradec...	336	C21H33F02	1000280-60-6	50
4		Pentanamide, N-(4-methoxyphenyl)-	207	C12H17NO2	1000306-89-3	49
5		2-Fluorobenzoic acid, 4-pentadec...	350	C22H35F02	1000280-61-7	47



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032818\
Data File : BF104014.D
Acq On : 28 Mar 2018 17:11
Operator : JU/SJ
Sample : J2087-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

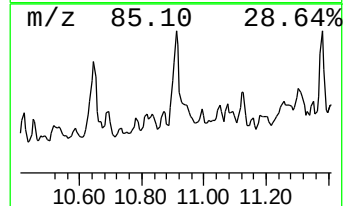
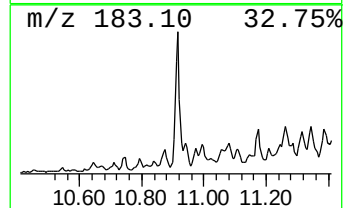
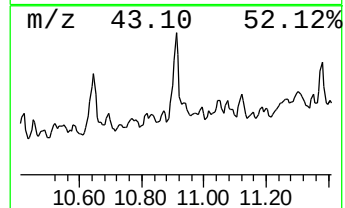
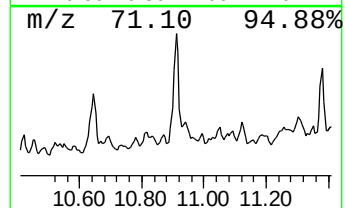
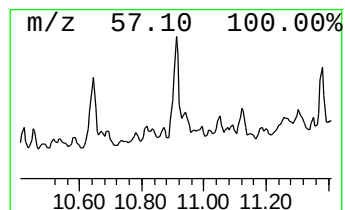
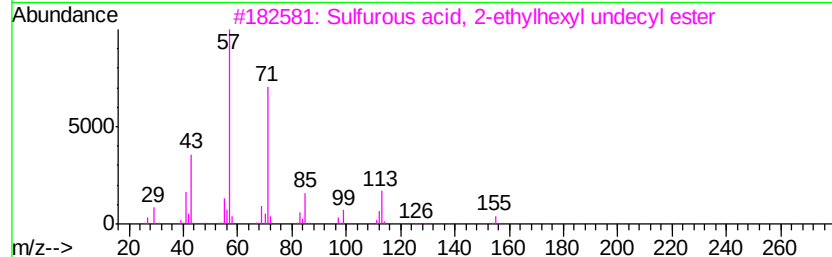
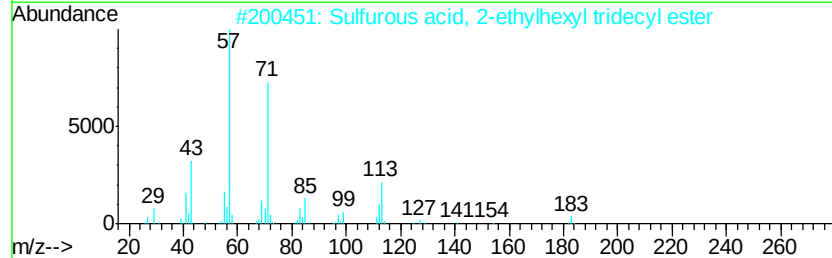
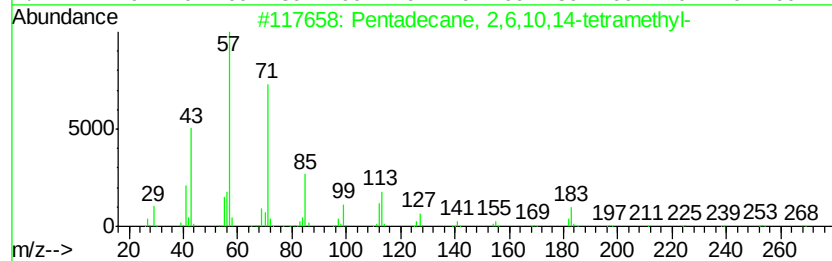
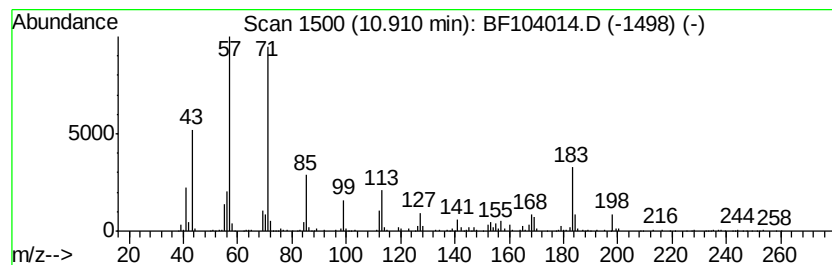
Instrument :
BNA_F
ClientSampleId :
JRS-RB201719RT-CV-2530

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Pentadecane, 2,6,10,14-tetr... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.91	20.00 ng	796596	Phenanthrene-d10	11.51
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Pentadecane, 2,6,10,14-tetramethyl-	268 C19H40	001921-70-6 86
2		Sulfurous acid, 2-ethylhexyl tri...	376 C21H44O3S	1000309-19-6 64
3		Sulfurous acid, 2-ethylhexyl und...	348 C19H40O3S	1000309-19-4 58
4		Dodecane, 2,6,11-trimethyl-	212 C15H32	031295-56-4 53
5		Sulfurous acid, dodecyl 2-ethylh...	362 C20H42O3S	1000309-19-5 53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032818\
Data File : BF104014.D
Acq On : 28 Mar 2018 17:11
Operator : JU/SJ
Sample : J2087-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
JRS-RB201719RT-CV-2530

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032818.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
Butane, 2-methoxy...	2.29	3.0	ng	110987	1	6.96	751847	20.0
2-Pentanone, 4-hy...	5.21	2.5	ng	92564	1	6.96	751847	20.0
unknown6.74	6.74	74.5	ng	2802530	1	6.96	751847	20.0
unknown9.23	9.23	6.5	ng	188474	3	10.00	581065	20.0
Decahydro-4,4,8,9...	9.39	11.2	ng	326057	3	10.00	581065	20.0
unknown9.63	9.63	4.5	ng	129709	3	10.00	581065	20.0
Naphthalene, deca...	9.79	3.8	ng	109508	3	10.00	581065	20.0
unknown9.86	9.86	25.2	ng	731729	3	10.00	581065	20.0
unknown9.91	9.91	37.8	ng	1096970	3	10.00	581065	20.0
unknown9.95	9.95	7.2	ng	210183	3	10.00	581065	20.0
Dodecane	10.25	15.1	ng	437964	3	10.00	581065	20.0
unknown10.32	10.32	9.2	ng	266432	3	10.00	581065	20.0
unknown10.41	10.41	37.1	ng	1076600	3	10.00	581065	20.0
Pentadecane, 2,6,...	10.91	20.0	ng	796596	4	11.51	796596	20.0