

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122223\
 Data File : BF136699.D
 Acq On : 22 Dec 2023 22:24
 Operator : CG\JU
 Sample : 05765-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RT-SB-03-9.5-10.5

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-BF122023.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF136699.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.187	13	17	28	rVB2	13907	25339	1.66%	0.209%
2	3.875	300	304	312	rVB	17250	23171	1.52%	0.191%
3	4.398	389	393	402	rBV	47213	52653	3.45%	0.434%
4	5.016	493	498	508	rVB	346759	382427	25.03%	3.151%
5	5.428	564	568	579	rBV	1201476	1024637	67.07%	8.443%
6	6.428	734	738	741	rBV	1469012	1408877	92.22%	11.609%
7	6.781	795	798	802	rBV	503785	387181	25.34%	3.190%
8	7.345	889	894	897	rBV	1116199	993543	65.03%	8.186%
9	8.057	1010	1015	1018	rBV	599275	474166	31.04%	3.907%
10	8.522	1092	1094	1097	rVB	30724	26303	1.72%	0.217%
11	8.598	1103	1107	1112	rBV5	20744	35612	2.33%	0.293%
12	8.692	1117	1123	1124	rBV6	18111	28802	1.89%	0.237%
13	8.745	1129	1132	1135	rVV4	22539	24764	1.62%	0.204%
14	8.798	1139	1141	1144	rVB3	34312	30631	2.00%	0.252%
15	8.833	1144	1147	1150	rBV3	17032	22985	1.50%	0.189%
16	8.939	1161	1165	1168	rBV4	22077	34418	2.25%	0.284%
17	9.016	1175	1178	1181	rBV3	26996	36319	2.38%	0.299%
18	9.051	1181	1184	1186	rVV2	67610	57888	3.79%	0.477%
19	9.075	1186	1188	1192	rVB4	21667	23589	1.54%	0.194%
20	9.139	1194	1199	1202	rBV	1671552	1527776	100.00%	12.588%
21	9.210	1208	1211	1215	rBV2	141279	143599	9.40%	1.183%
22	9.257	1217	1219	1222	rBV4	37683	35887	2.35%	0.296%
23	9.522	1261	1264	1268	rBV	215091	183919	12.04%	1.515%
24	9.575	1271	1273	1277	rVB3	84693	96534	6.32%	0.795%
25	9.680	1288	1291	1293	rBV2	205825	214388	14.03%	1.766%
26	9.704	1293	1295	1296	rVV	87884	75048	4.91%	0.618%
27	9.722	1296	1298	1301	rVB	337315	316258	20.70%	2.606%
28	9.763	1301	1305	1307	rBV3	96360	106292	6.96%	0.876%
29	9.798	1307	1311	1312	rVV	140067	167932	10.99%	1.384%
30	9.816	1312	1314	1317	rVB	545211	419066	27.43%	3.453%
31	9.851	1317	1320	1325	rBV3	97001	159986	10.47%	1.318%
32	10.151	1369	1371	1374	rVB	98732	75086	4.91%	0.619%
33	10.186	1374	1377	1380	rVB5	84952	82141	5.38%	0.677%
34	10.222	1380	1383	1386	rBV2	300839	258937	16.95%	2.134%
35	10.610	1446	1449	1455	rBV	331965	374098	24.49%	3.082%
36	11.310	1565	1568	1570	rBV	416264	364264	23.84%	3.001%

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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-BF122023.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	11.333	1570	1572	1576	rVB	159175	140220	9.18%	1.155%
38	12.527	1772	1775	1778	rVB	202300	156664	10.25%	1.291%
39	12.757	1812	1814	1817	rVB	157309	124947	8.18%	1.030%
40	12.904	1835	1839	1842	rBV	1235099	1047316	68.55%	8.629%
41	13.963	2014	2019	2021	rBV2	424988	458096	29.98%	3.775%
42	13.986	2021	2023	2026	rVB	78766	59449	3.89%	0.490%
43	15.004	2193	2196	2204	rVB	79924	137339	8.99%	1.132%
44	15.439	2266	2270	2274	rBV	274291	317966	20.81%	2.620%

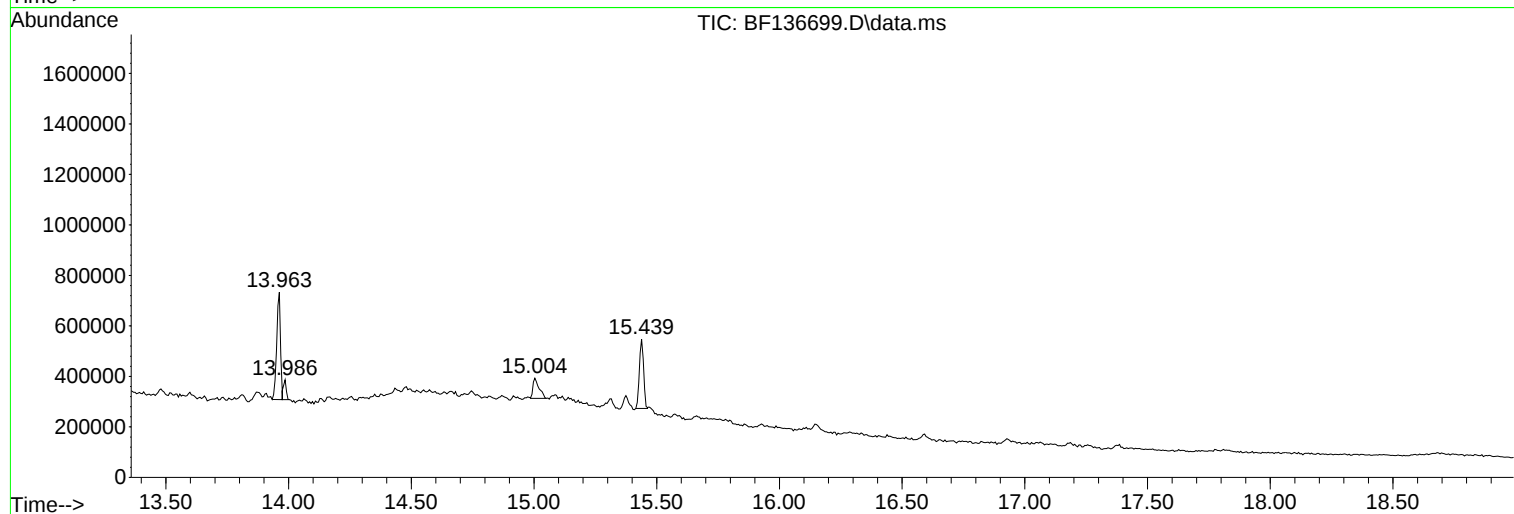
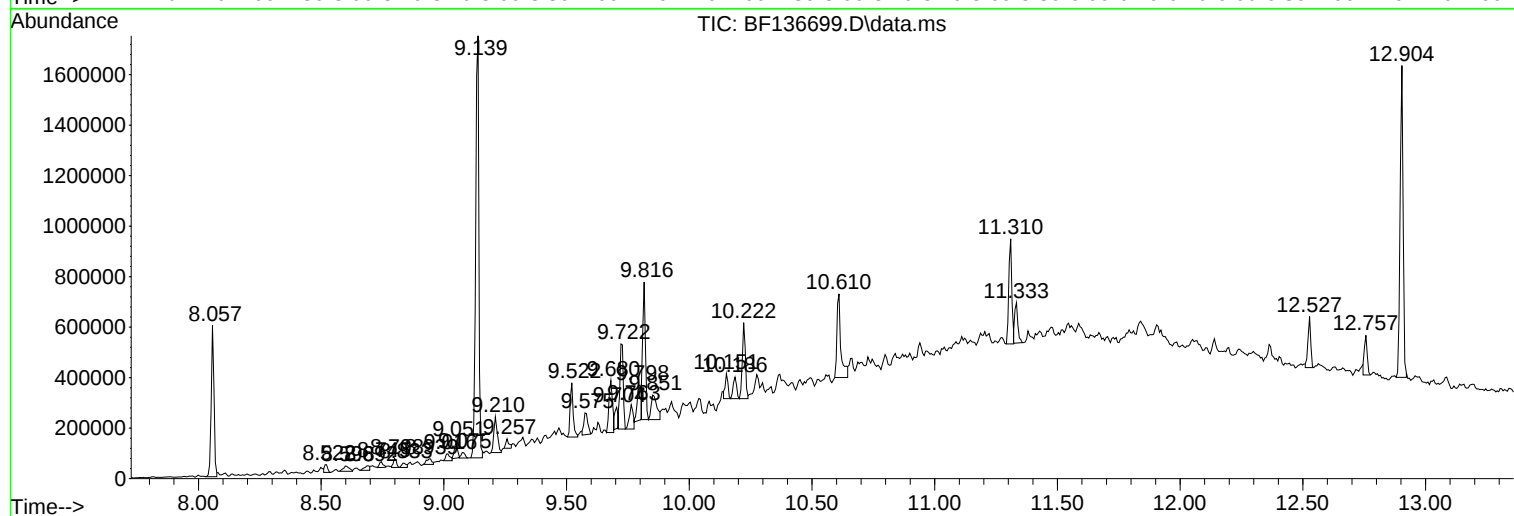
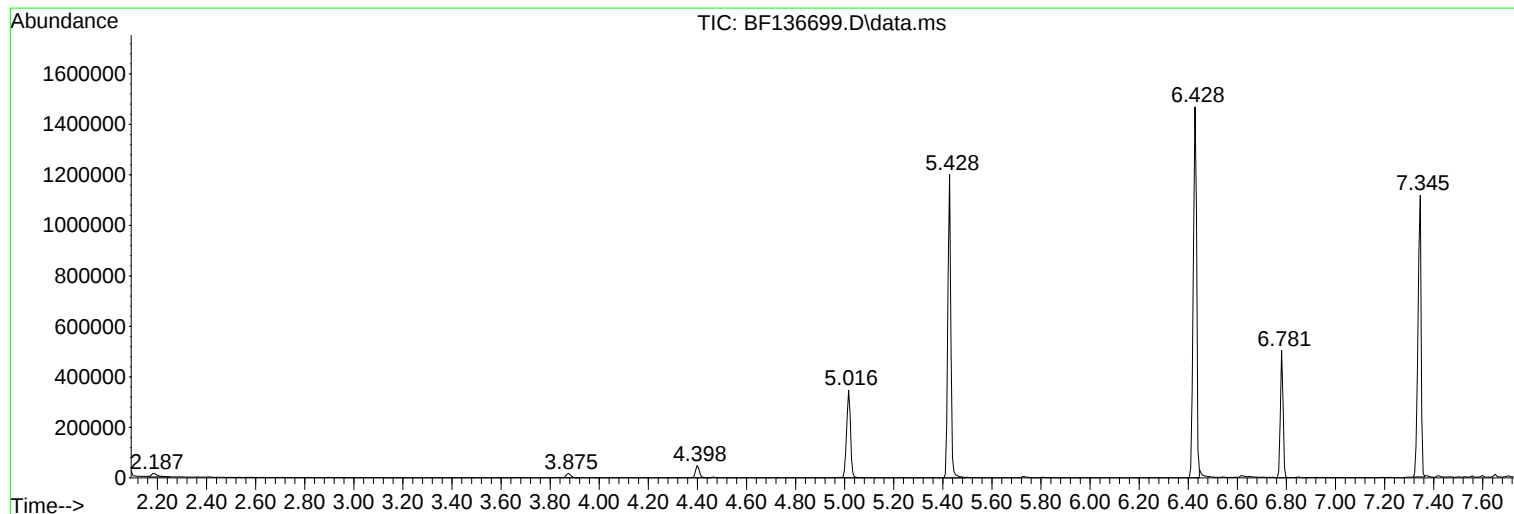
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RT-SB-03-9.5-10.5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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BNA_F
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RT-SB-03-9.5-10.5

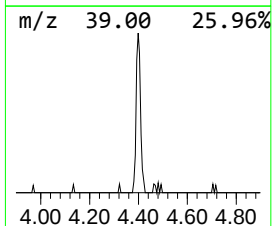
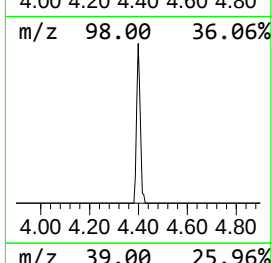
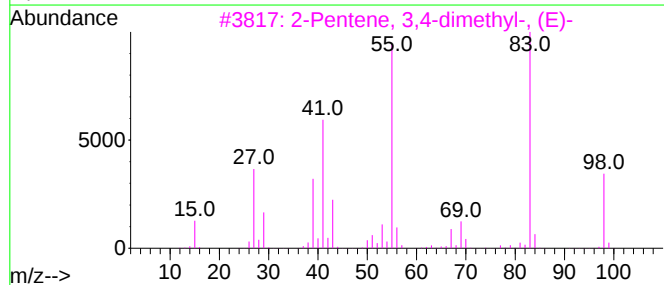
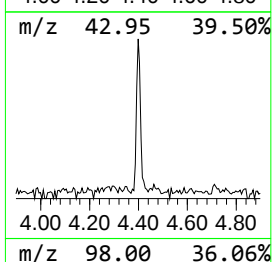
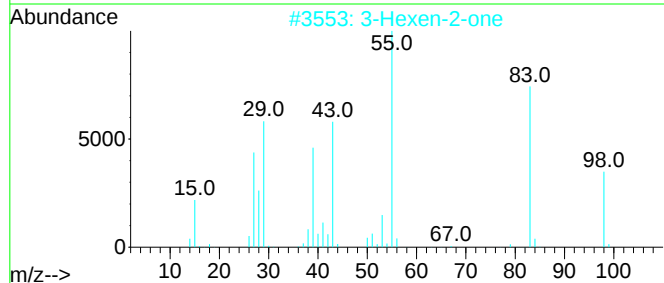
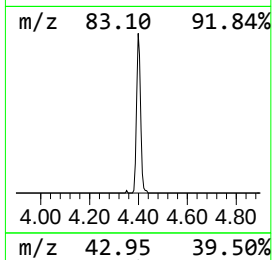
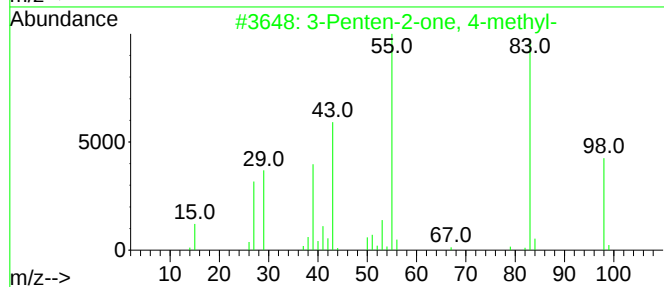
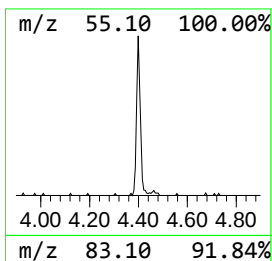
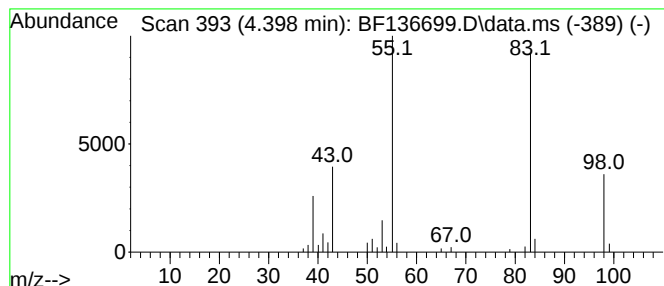
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.398	2.72 ng	52653	1,4-Dichlorobenzene-d4	6.781

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
2		3-Hexen-2-one	98	C6H10O	000763-93-9	91
3		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	86
4		2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	83
5		2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	80



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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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RT-SB-03-9.5-10.5

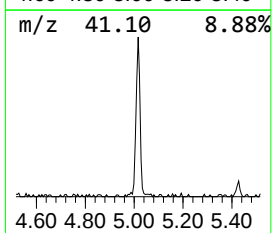
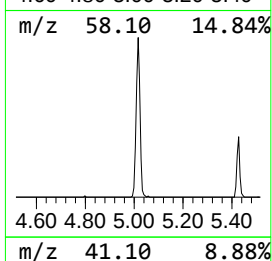
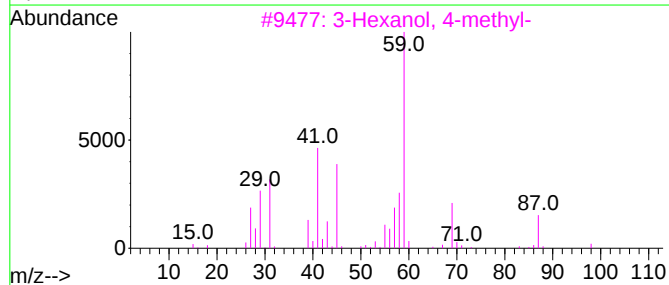
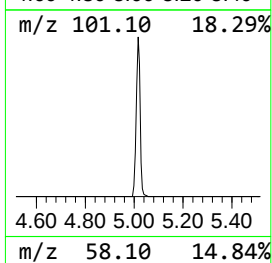
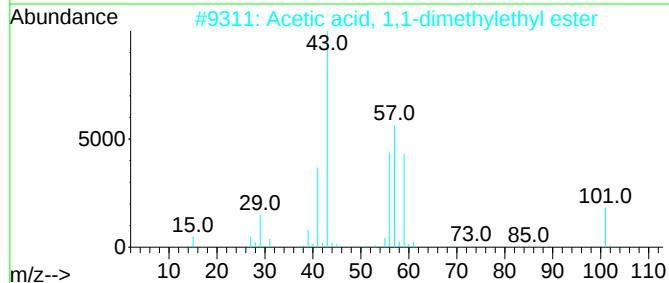
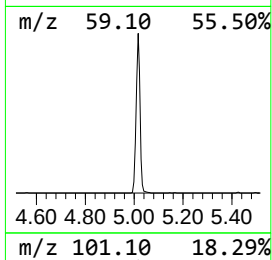
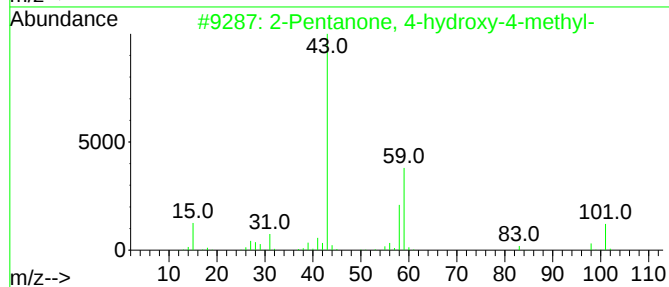
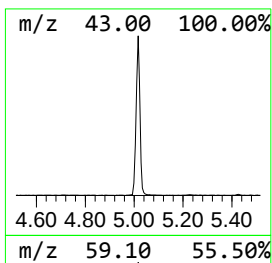
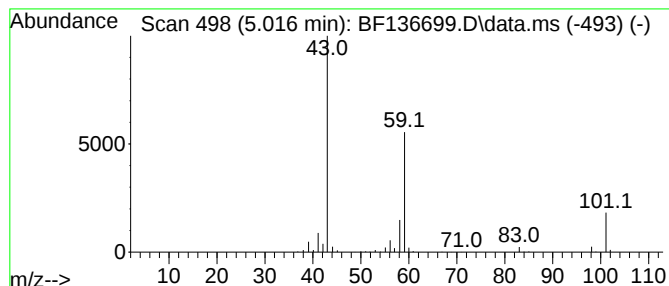
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.016	19.75 ng	382427	1,4-Dichlorobenzene-d4	6.781

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	39
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32



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Instrument :
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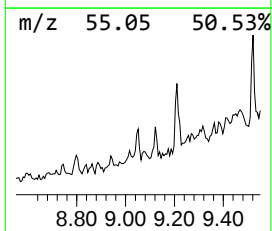
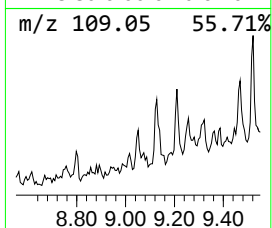
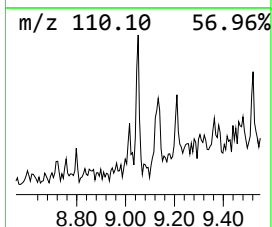
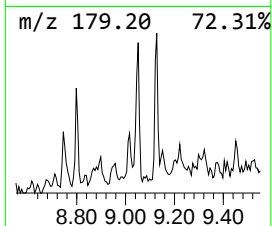
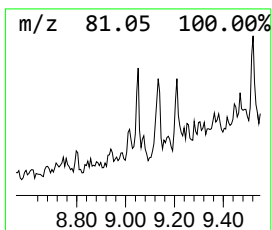
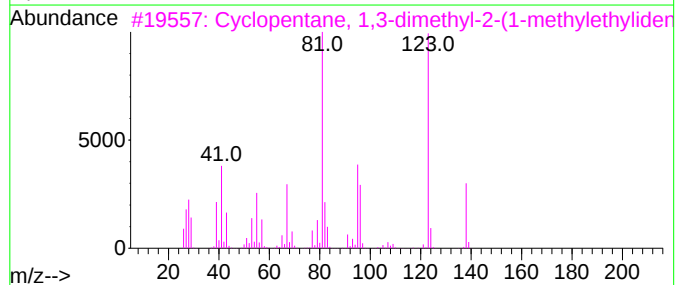
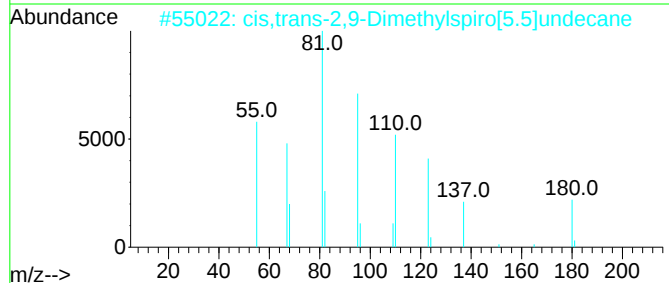
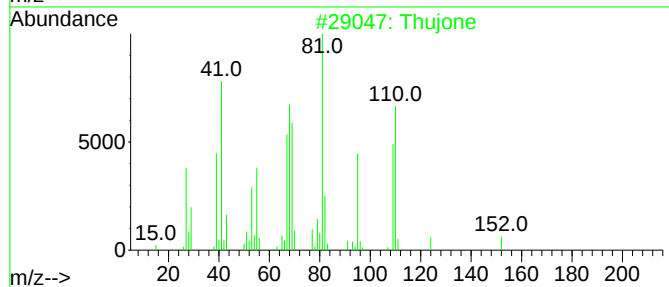
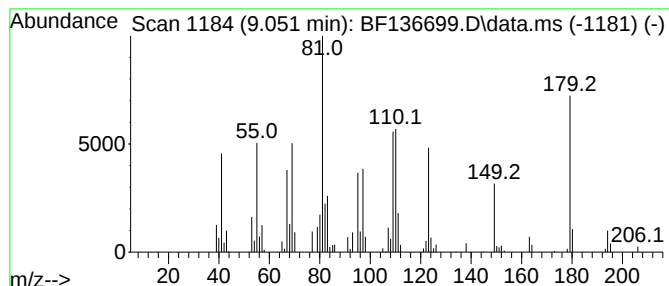
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Thujone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.051	2.76 ng	57888	Acenaphthene-d10	9.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Thujone			152	C10H16O	000546-80-5	50
2	cis,trans-2,9-Dimethylspiro[5.5]...			180	C13H24	095530-74-8	43
3	Cyclopentane, 1,3-dimethyl-2-(1-...			138	C10H18	061142-30-1	38
4	Cyclopropane, tetramethylpropyli...			138	C10H18	024519-04-8	30
5	Bicyclo[2.2.1]heptane, 1,3,3-tri...			138	C10H18	006248-88-0	25



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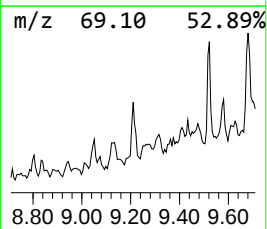
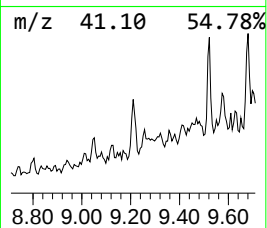
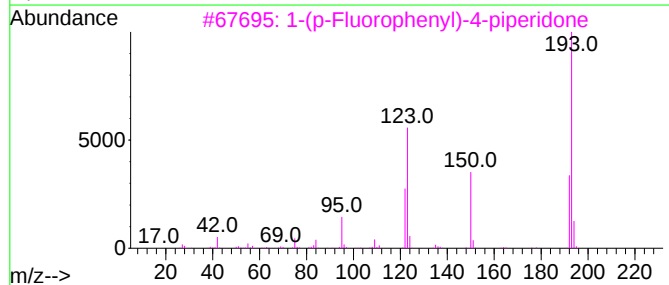
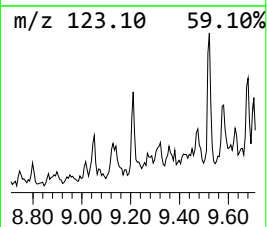
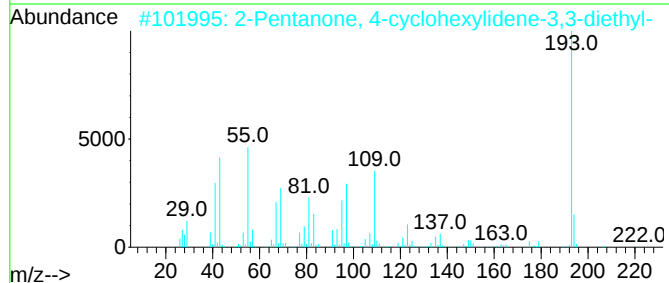
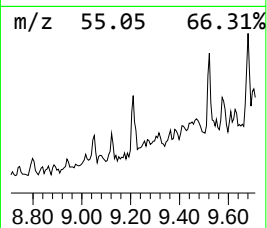
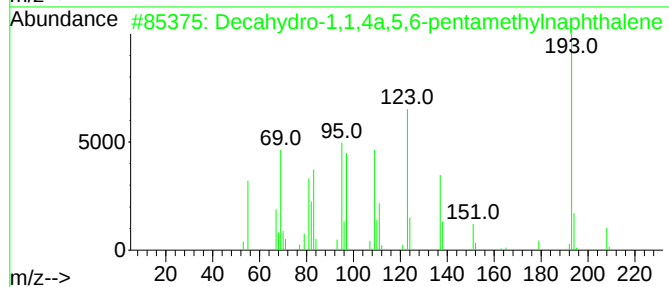
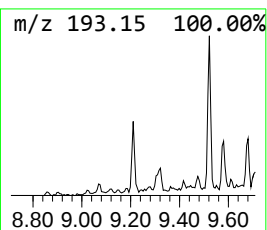
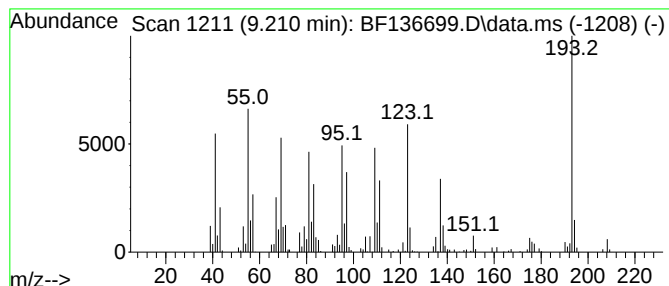
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 Decahydro-1,1,4a,5,6-pentam... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.210	6.85 ng	143599	Acenaphthene-d10	9.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	90
2			2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	52
3			1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	43
4			9-Phenanthrenamine	193	C14H11N	000947-73-9	30
5			2-Anthracenamine	193	C14H11N	000613-13-8	27



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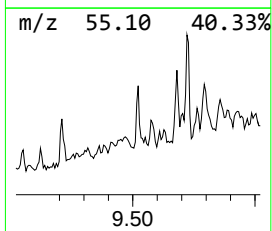
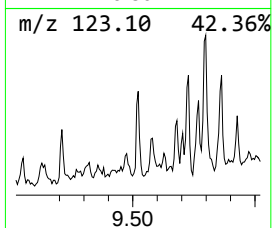
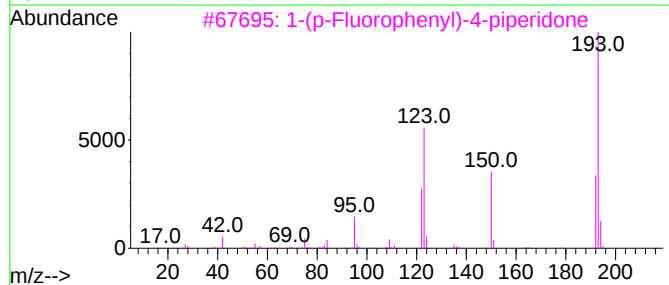
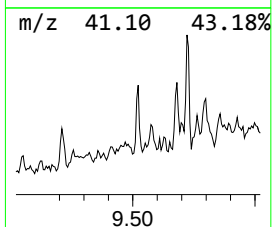
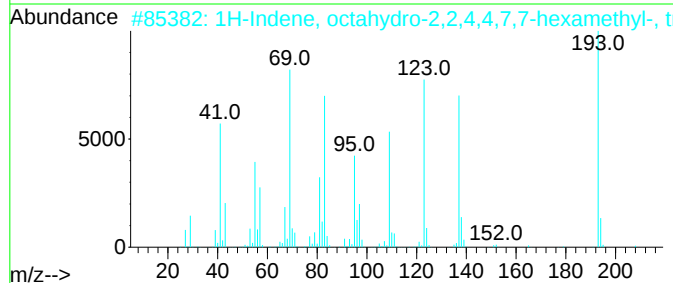
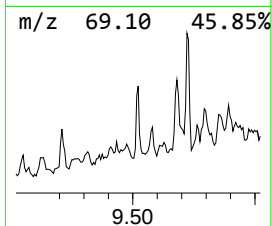
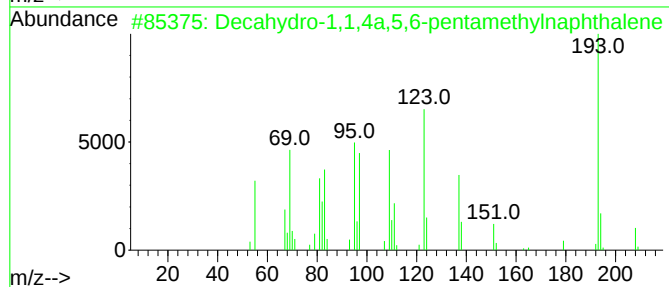
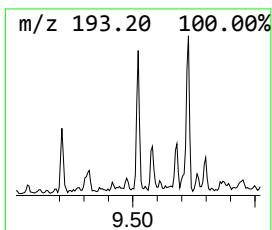
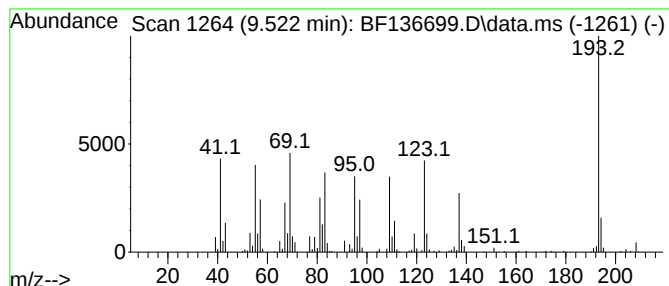
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown9.522 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.522	8.78 ng	183919	Acenaphthene-d10	9.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	52
2			1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	43
3			1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	40
4			2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	37
5			1-Anthracenamine	193	C14H11N	000610-49-1	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122223\
Data File : BF136699.D
Acq On : 22 Dec 2023 22:24
Operator : CG\JU
Sample : 05765-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RT-SB-03-9.5-10.5

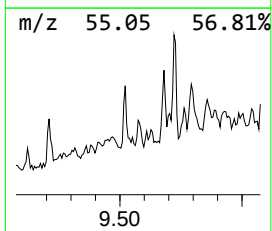
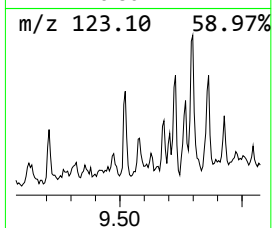
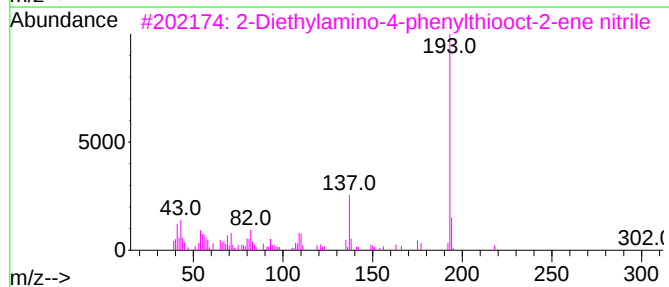
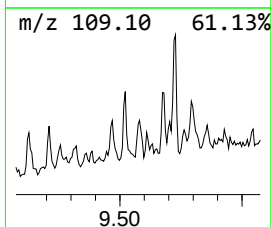
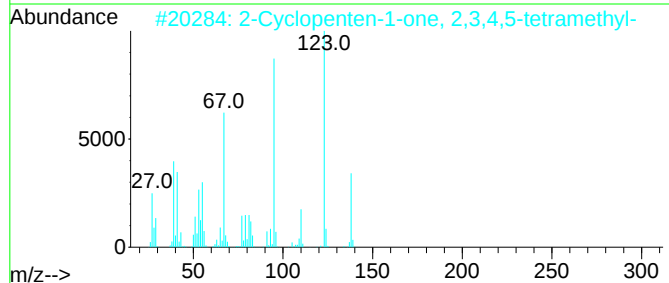
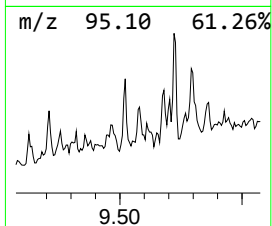
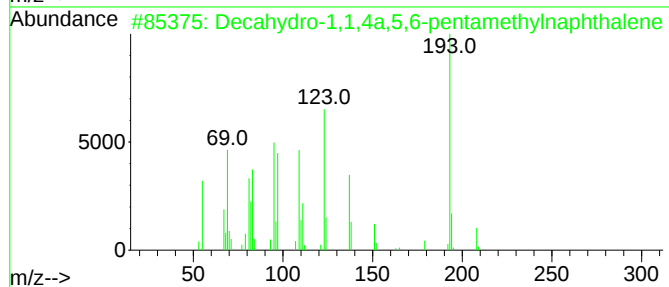
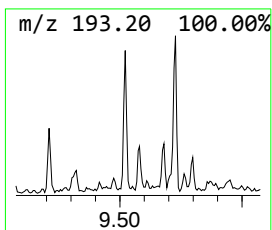
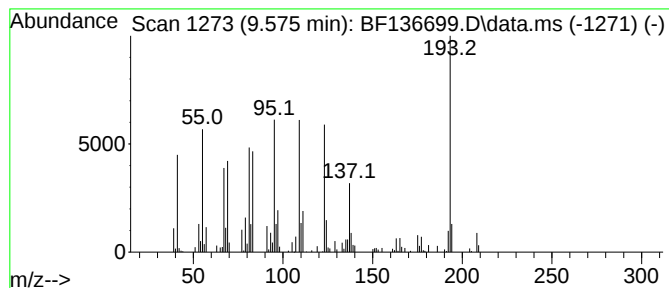
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown9.575 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.575	4.61 ng	96534	Acenaphthene-d10	9.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	89
2			2-Cyclopenten-1-one, 2,3,4,5-tet...	138	C9H14O	054458-61-6	30
3			2-Diethylamino-4-phenylthiooct-2...	302	C18H26N2S	1000090-57-0	22
4			1H-Indole, 2-phenyl-	193	C14H11N	000948-65-2	18
5			2-Anthracenamine	193	C14H11N	000613-13-8	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122223\
Data File : BF136699.D
Acq On : 22 Dec 2023 22:24
Operator : CG\JU
Sample : 05765-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
RT-SB-03-9.5-10.5

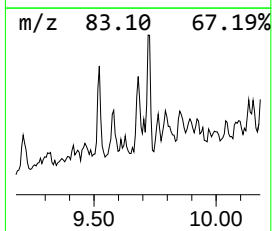
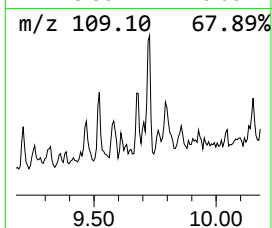
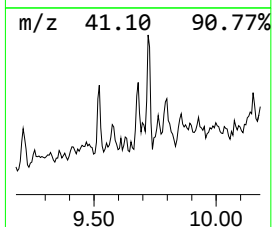
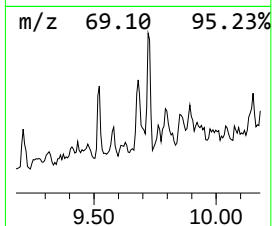
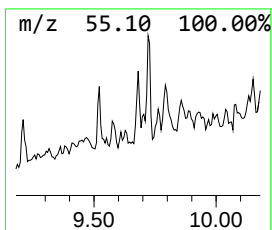
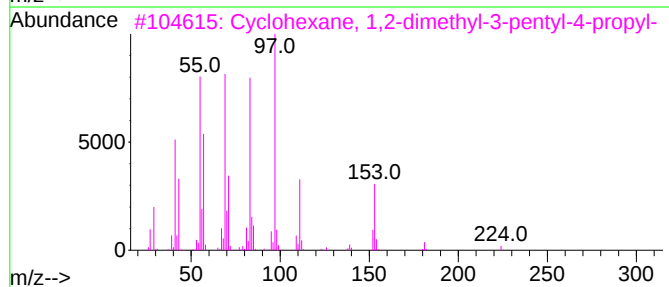
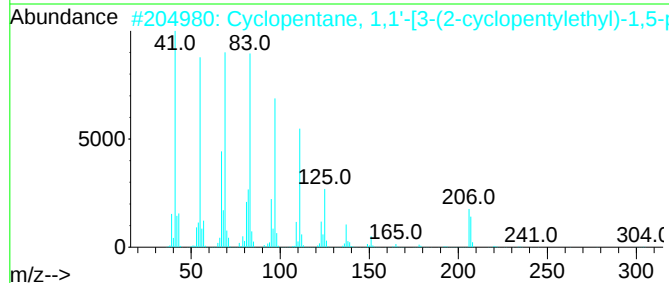
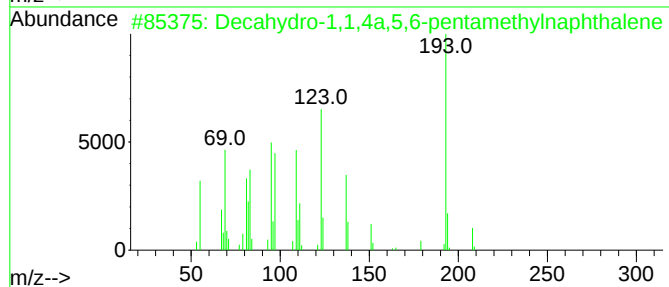
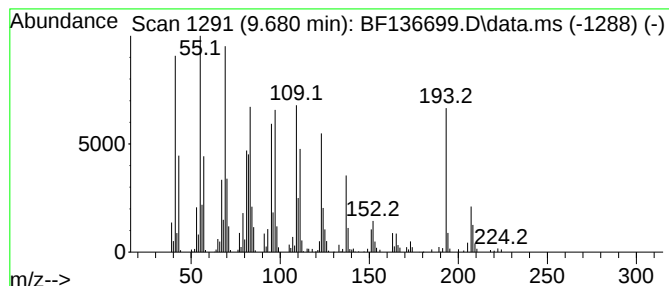
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown9.680 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.680	10.23 ng	214388	Acenaphthene-d10	9.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	93
2			Cyclopentane, 1,1'-[3-(2-cyclope...	304	C22H40	055255-85-1	38
3			Cyclohexane, 1,2-dimethyl-3-pent...	224	C16H32	062376-17-4	30
4			Octane, 1,8-dibromo-	270	C8H16Br2	004549-32-0	30
5			1-Hexyl-1-nitrocyclohexane	213	C12H23NO2	118252-09-8	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122223\
Data File : BF136699.D
Acq On : 22 Dec 2023 22:24
Operator : CG\JU
Sample : 05765-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

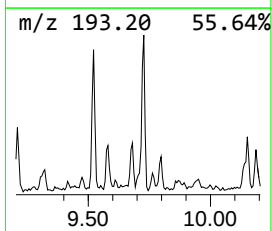
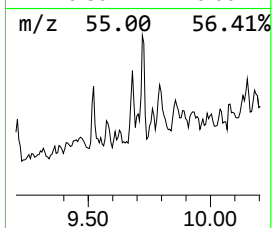
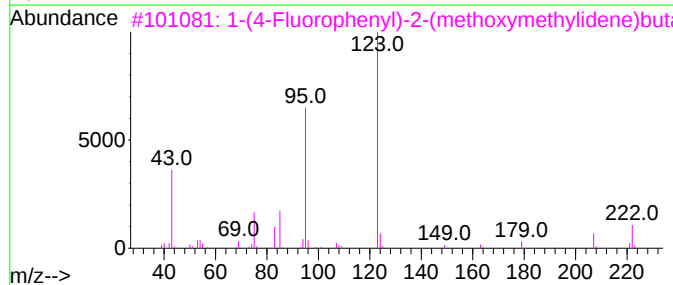
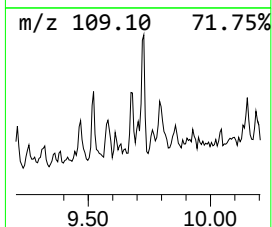
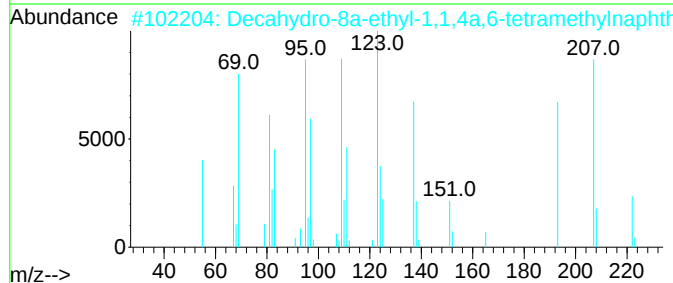
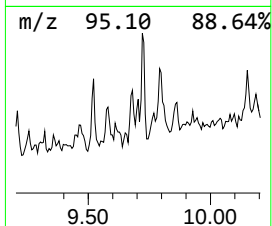
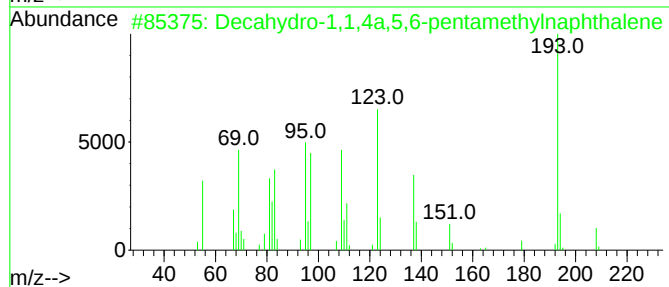
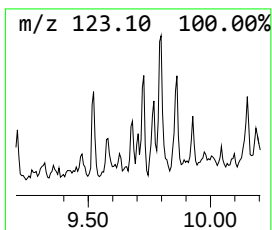
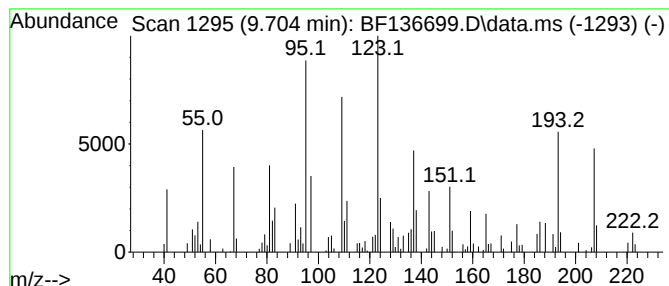
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RT-SB-03-9.5-10.5

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 unknown9.704 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.704	3.58 ng	75048	Acenaphthene-d10	9.816	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	49
2	Decahydro-8a-ethyl-1,1,4a,6-tetr...	222	C16H30	1000100-23-6	49
3	1-(4-Fluorophenyl)-2-(methoxymet...	222	C12H11FO3	1000388-14-4	22
4	1-(2,4-Diethoxyphenyl)propan-1-one	222	C13H18O3	1000410-44-8	20
5	Borinic acid, diethyl-, 3,3,5-tr...	208	C13H25BO	057387-76-5	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122223\
Data File : BF136699.D
Acq On : 22 Dec 2023 22:24
Operator : CG\JU
Sample : 05765-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
RT-SB-03-9.5-10.5

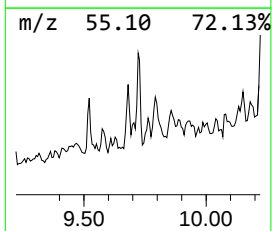
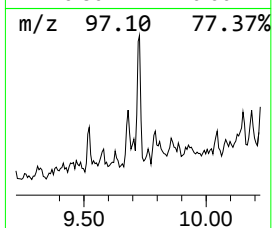
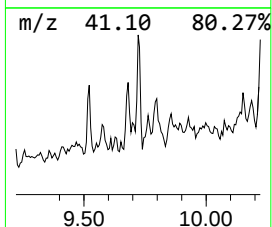
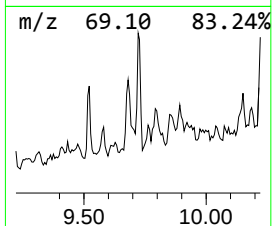
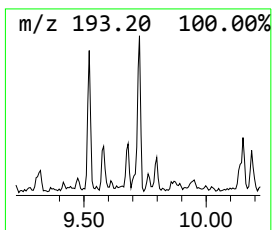
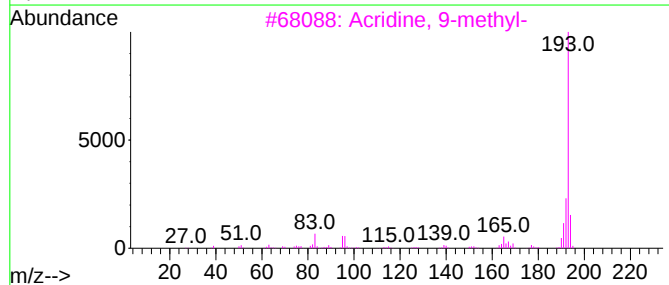
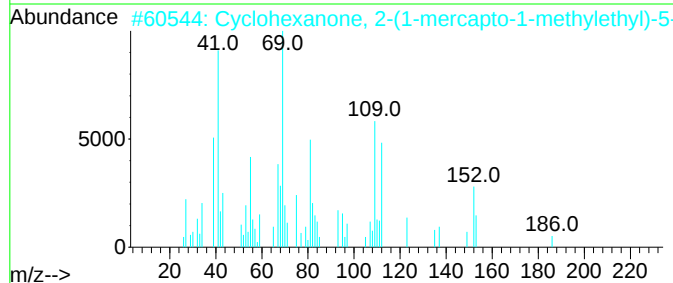
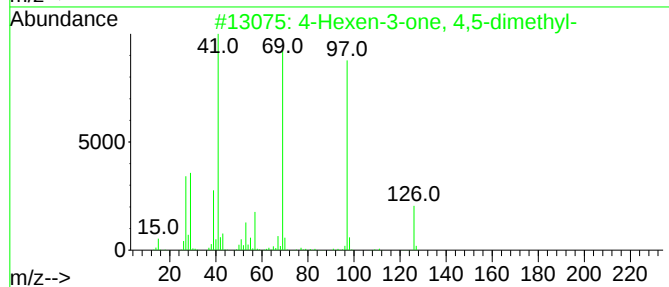
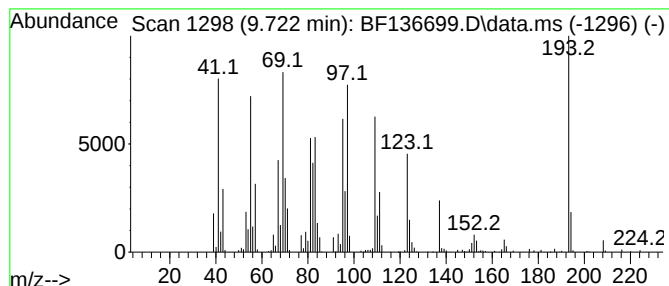
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown9.722 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.722	15.09 ng	316258	Acenaphthene-d10	9.816

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Hexen-3-one, 4,5-dimethyl-	126	C8H14O	017325-90-5	25	
2	Cyclohexanone, 2-(1-mercapto-1-methyl-5-oxo-5-phenylethyl)-	186	C10H18OS	033281-91-3	22	
3	Acridine, 9-methyl-	193	C14H11N	000611-64-3	14	
4	5-Hexen-1-ol, 2-(2-methyl-1-propenyl)-	154	C10H18O	018675-19-9	14	
5	Cyclopentanecarboxylic acid, 2-penten-3-yl-	324	C21H40O2	1010280-59-0	14	



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122223\
Data File : BF136699.D
Acq On : 22 Dec 2023 22:24
Operator : CG\JU
Sample : 05765-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

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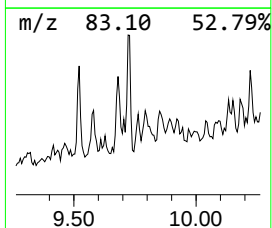
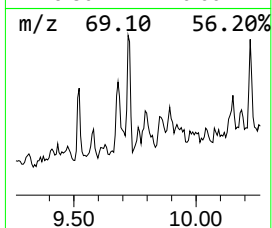
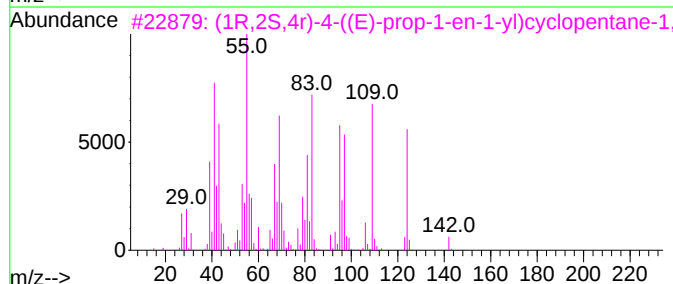
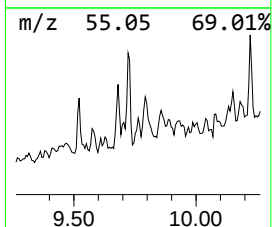
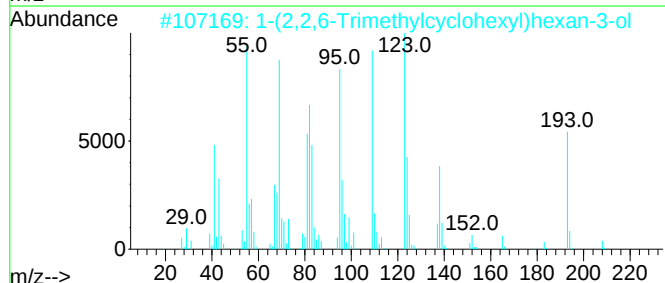
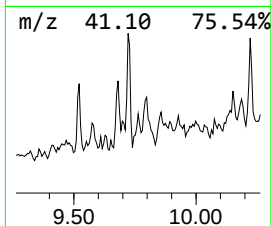
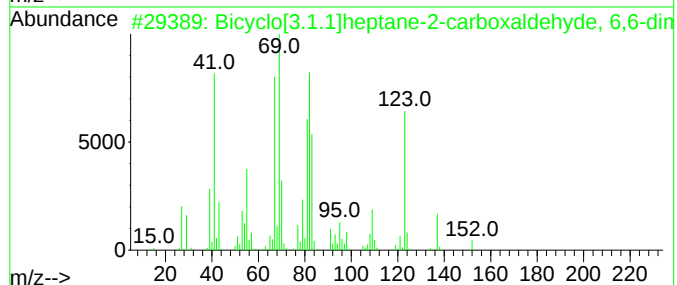
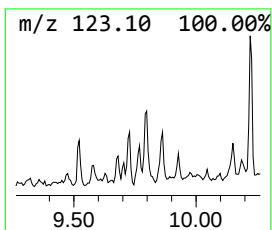
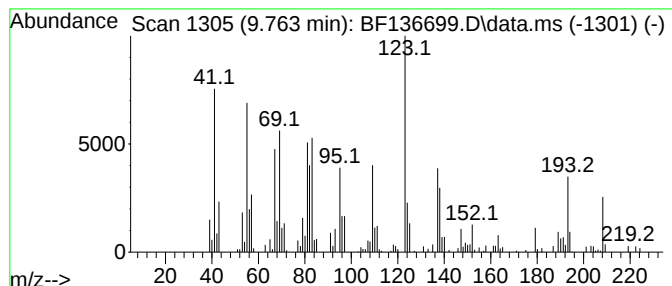
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 unknown9.763 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.763	5.07 ng	106292	Acenaphthene-d10	9.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[3.1.1]heptane-2-carboxal...	152	C10H16O	004764-14-1	49
2			1-(2,2,6-Trimethylcyclohexyl)hex...	226	C15H30O	070788-30-6	49
3			(1R,2S,4r)-4-((E)-prop-1-en-1-yl...	142	C8H14O2	1010481-89-9	38
4			1,3-Benzenediamine, 4-methoxy-	138	C7H10N2O	000615-05-4	35
5			Cyclopropane, tetramethylpropyli...	138	C10H18	024519-04-8	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122223\
Data File : BF136699.D
Acq On : 22 Dec 2023 22:24
Operator : CG\JU
Sample : 05765-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

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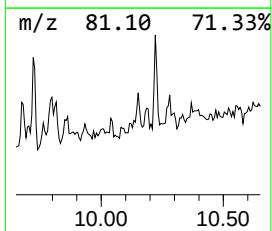
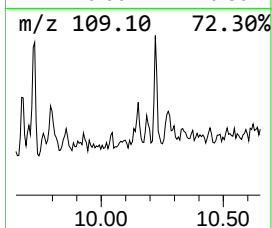
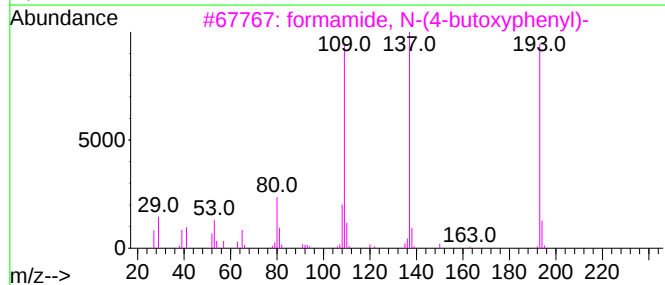
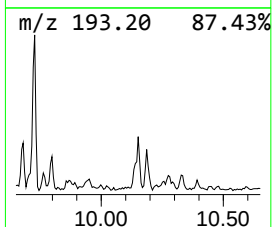
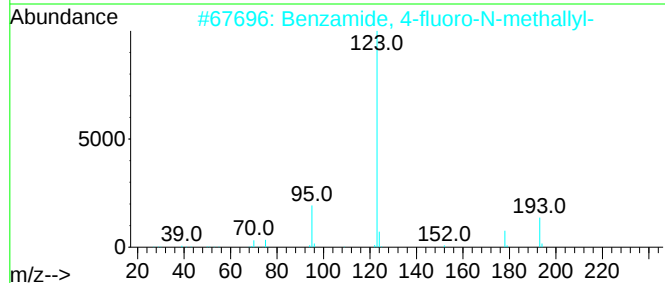
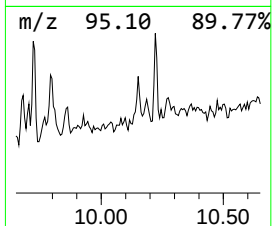
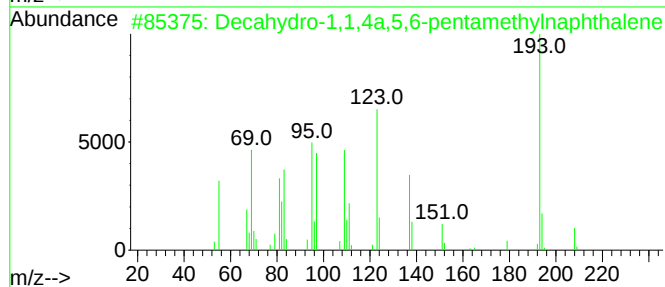
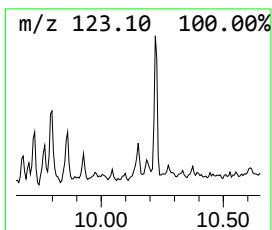
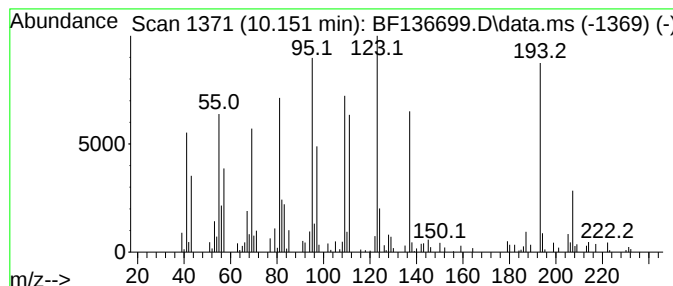
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 unknown10.151 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.151	3.58 ng	75086	Acenaphthene-d10	9.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	64
2			Benzamide, 4-fluoro-N-methallyl-	193	C11H12FNO	1000340-31-9	43
3			formamide, N-(4-butoxyphenyl)-	193	C11H15NO2	1000400-28-0	30
4			6-Acetamido-1,4-benzodioxane	193	C10H11NO3	063546-19-0	22
5			3-Octyne, 2,2,7-trimethyl-	152	C11H20	055402-13-6	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122223\
Data File : BF136699.D
Acq On : 22 Dec 2023 22:24
Operator : CG\JU
Sample : 05765-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RT-SB-03-9.5-10.5

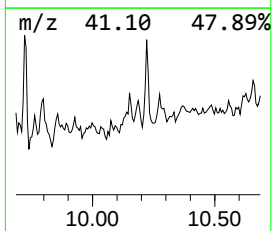
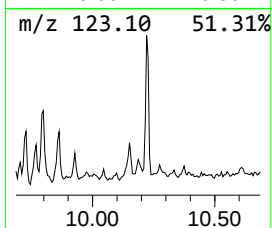
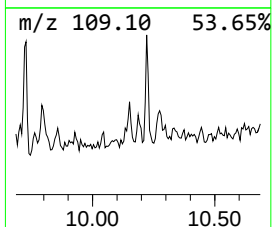
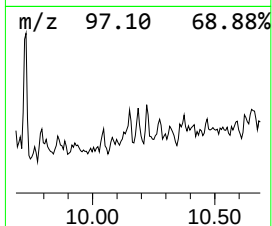
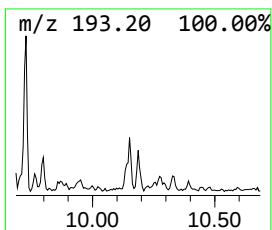
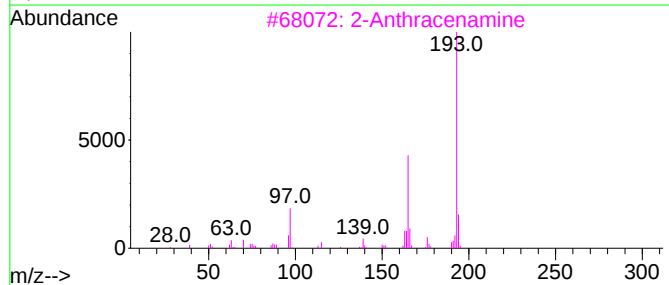
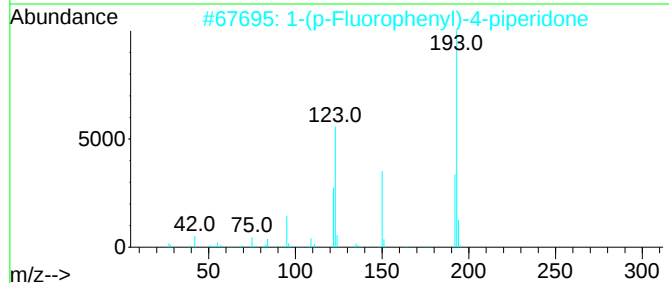
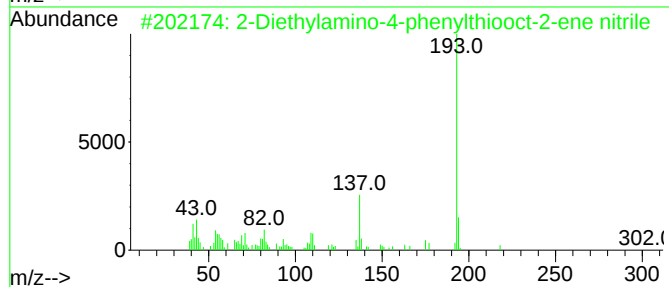
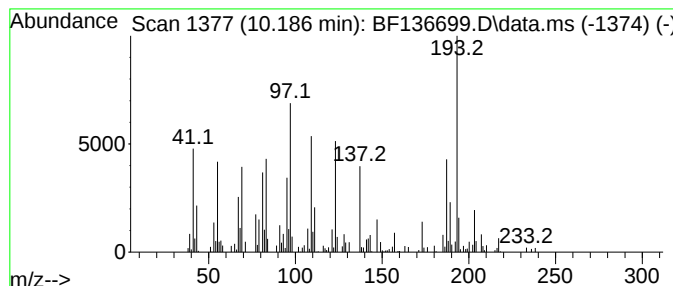
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown10.186 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.186	3.92 ng	82141	Acenaphthene-d10	9.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Diethylamino-4-phenylthiooct-2...	302	C18H26N2S	1000090-57-0	27		
2	1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	27		
3	2-Anthracenamine	193	C14H11N	000613-13-8	22		
4	(1-Acetyloxy-3-hydroxy-6,8a-dime...	502	C27H38O7Si	1000494-92-2	14		
5	9-Phenanthrenamine	193	C14H11N	000947-73-9	11		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122223\
Data File : BF136699.D
Acq On : 22 Dec 2023 22:24
Operator : CG\JU
Sample : 05765-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RT-SB-03-9.5-10.5

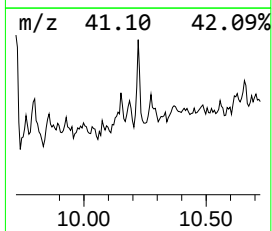
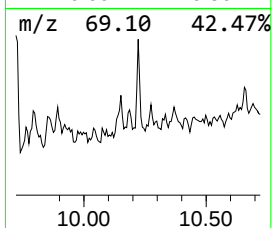
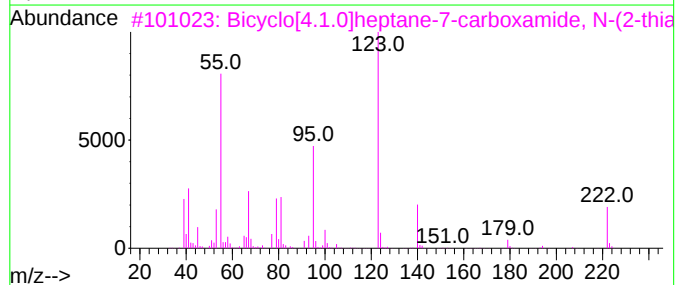
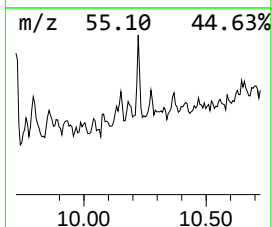
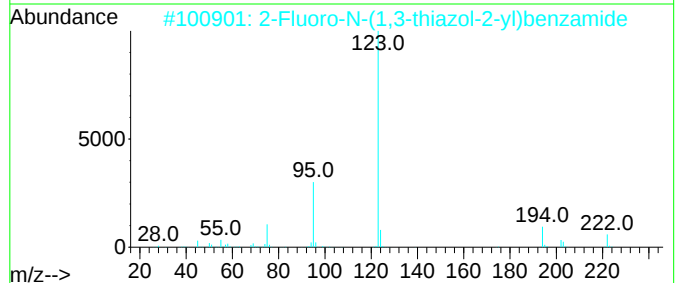
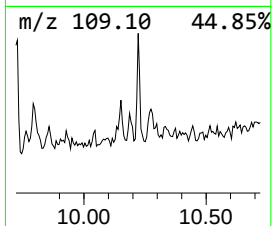
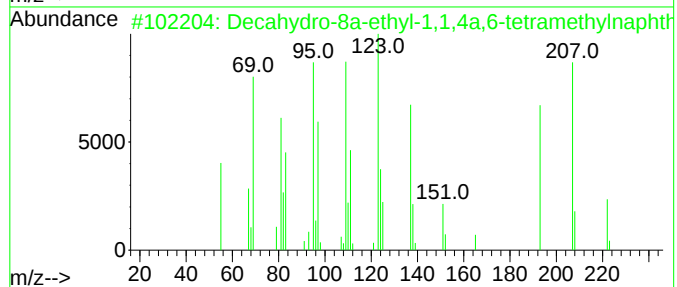
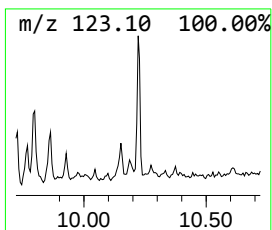
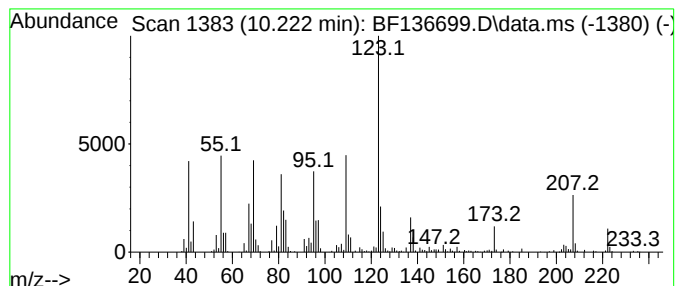
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown10.222 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.222	12.36 ng	258937	Acenaphthene-d10	9.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decahydro-8a-ethyl-1,1,4a,6-tetr...	222	C16H30	1000100-23-6	43
2			2-Fluoro-N-(1,3-thiazol-2-yl)ben...	222	C10H7FN2OS	000319-38-0	38
3			Bicyclo[4.1.0]heptane-7-carboxam...	222	C11H14N2OS	329912-72-3	38
4			Ethanone, 1-[3-[2-methyl-2-(5-me...	222	C13H18O3	080114-28-9	38
5			1-Butanone, 1-bicyclo[4.1.0]hept...	166	C11H18O	054764-62-4	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122223\
Data File : BF136699.D
Acq On : 22 Dec 2023 22:24
Operator : CG\JU
Sample : 05765-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RT-SB-03-9.5-10.5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-one,...	4.398	2.7	ng	52653	1	6.781	387181	20.0
2-Pentanone, 4-...	5.016	19.8	ng	382427	1	6.781	387181	20.0
Thujone	9.051	2.8	ng	57888	3	9.816	419066	20.0
Decahydro-1,1,4...	9.210	6.8	ng	143599	3	9.816	419066	20.0
unknown9.522	9.522	8.8	ng	183919	3	9.816	419066	20.0
unknown9.575	9.575	4.6	ng	96534	3	9.816	419066	20.0
unknown9.680	9.680	10.2	ng	214388	3	9.816	419066	20.0
unknown9.704	9.704	3.6	ng	75048	3	9.816	419066	20.0
unknown9.722	9.722	15.1	ng	316258	3	9.816	419066	20.0
unknown9.763	9.763	5.1	ng	106292	3	9.816	419066	20.0
unknown10.151	10.151	3.6	ng	75086	3	9.816	419066	20.0
unknown10.186	10.186	3.9	ng	82141	3	9.816	419066	20.0
unknown10.222	10.222	12.4	ng	258937	3	9.816	419066	20.0