

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101823\
 Data File : BF135876.D
 Acq On : 19 Oct 2023 06:00
 Operator : CG\JU
 Sample : 04740-05
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW4

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-BF101323.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF135876.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.963	485	489	499	rBV	41672	60968	2.77%	0.510%
2	5.040	499	502	510	rVB	20976	28360	1.29%	0.237%
3	5.381	557	560	581	rBV	779596	976427	44.31%	8.163%
4	6.398	730	733	747	rBV	446517	606144	27.51%	5.067%
5	6.739	787	791	800	rBV	459255	395224	17.94%	3.304%
6	7.310	884	888	891	rBV	1495359	1283569	58.25%	10.731%
7	7.586	932	935	946	rBV	172803	178608	8.11%	1.493%
8	8.016	1005	1008	1014	rBV	574874	480964	21.83%	4.021%
9	8.086	1018	1020	1028	rVV	46355	58988	2.68%	0.493%
10	8.186	1033	1037	1040	rBV	725895	608470	27.61%	5.087%
11	8.757	1131	1134	1139	rBV	19805	33892	1.54%	0.283%
12	9.098	1188	1192	1195	rBV	2686082	2203485	100.00%	18.421%
13	9.316	1226	1229	1233	rBV	63647	48125	2.18%	0.402%
14	9.775	1304	1307	1315	rVB	550756	442290	20.07%	3.698%
15	10.310	1394	1398	1404	rVB2	42723	45661	2.07%	0.382%
16	10.575	1439	1443	1454	rBV	1156757	1064632	48.32%	8.900%
17	10.657	1454	1457	1462	rVV	104610	103813	4.71%	0.868%
18	10.698	1462	1464	1469	rVB3	24964	28417	1.29%	0.238%
19	11.010	1514	1517	1523	rBV	51133	68551	3.11%	0.573%
20	11.274	1558	1562	1570	rVB	465572	400731	18.19%	3.350%
21	11.474	1592	1596	1600	rBV	41373	41977	1.91%	0.351%
22	11.810	1650	1653	1665	rBV2	92541	117964	5.35%	0.986%
23	12.521	1769	1774	1784	rBV4	40603	92671	4.21%	0.775%
24	12.598	1784	1787	1795	rVV2	80363	93283	4.23%	0.780%
25	12.751	1810	1813	1823	rBV2	27475	48778	2.21%	0.408%
26	12.874	1829	1834	1837	rBV	1670215	1691211	76.75%	14.139%
27	13.827	1992	1996	2003	rBV6	22652	47118	2.14%	0.394%
28	13.939	2012	2015	2020	rBV	382300	362635	16.46%	3.032%
29	14.774	2154	2157	2162	rVB	43402	46439	2.11%	0.388%
30	15.421	2262	2267	2274	rBV	236897	302206	13.71%	2.526%

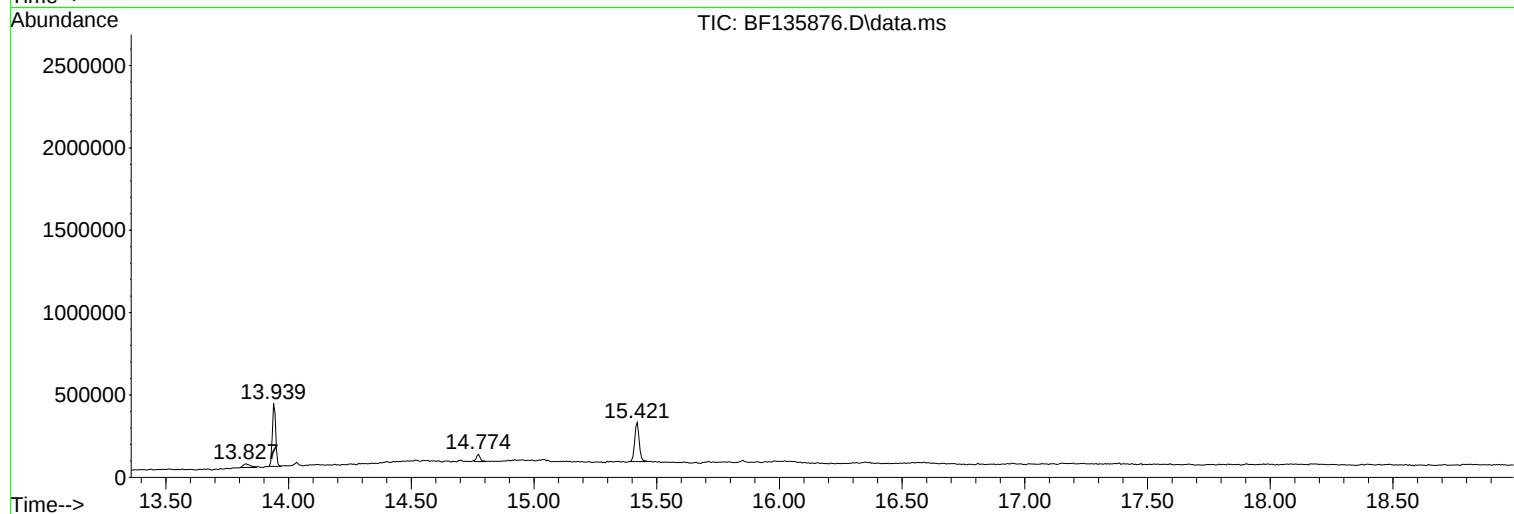
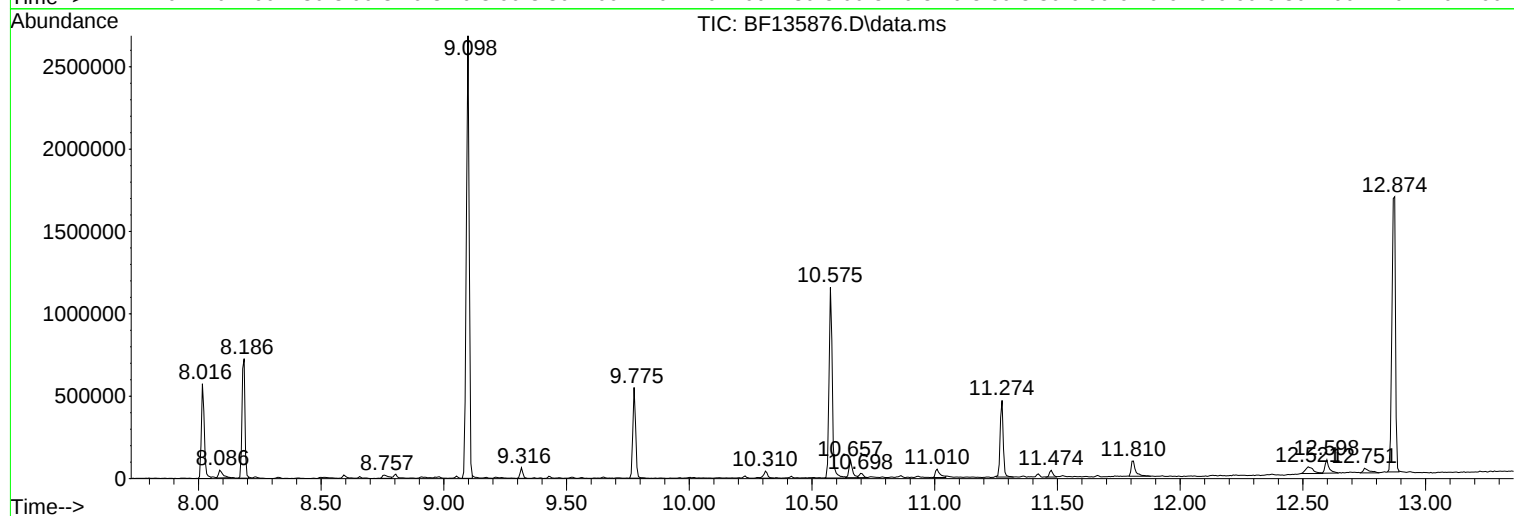
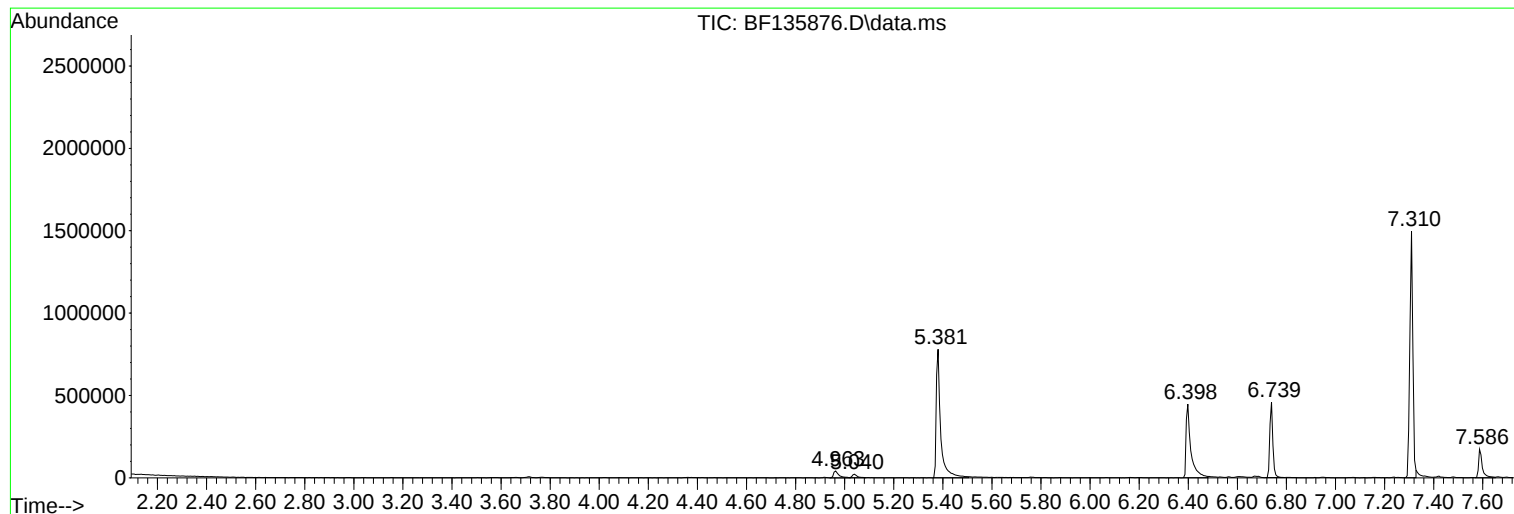
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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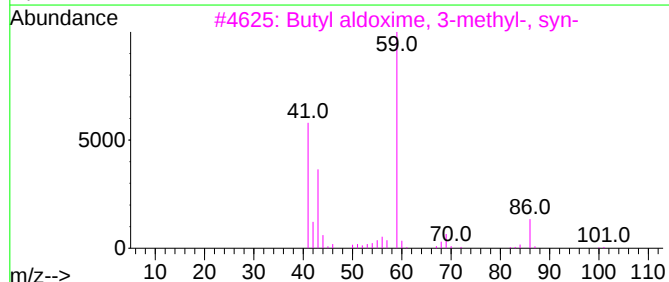
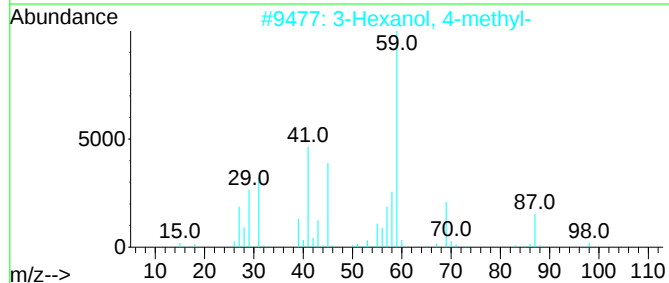
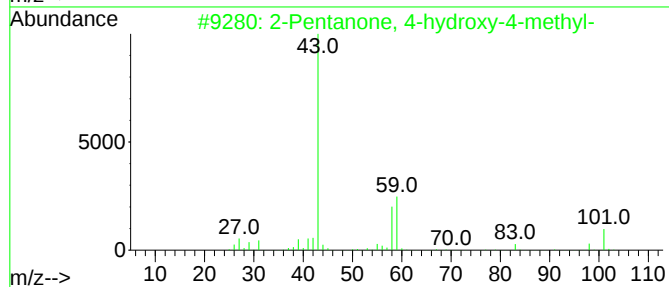
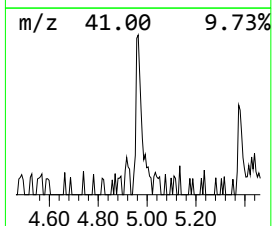
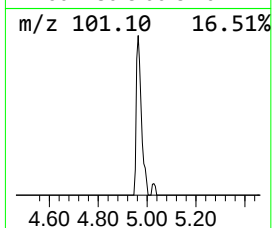
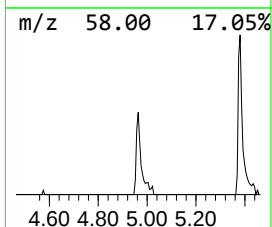
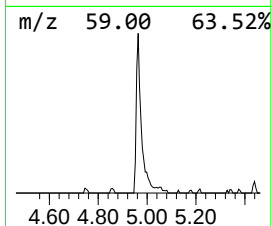
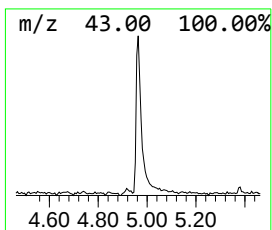
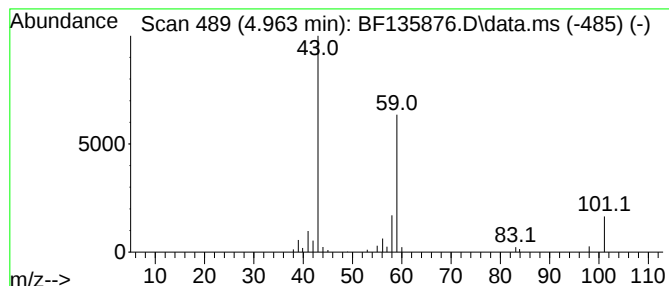
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.963	3.09 ng	60968	1,4-Dichlorobenzene-d4	6.739

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53		
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33		
3	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	32		
4	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28		
5	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	25		



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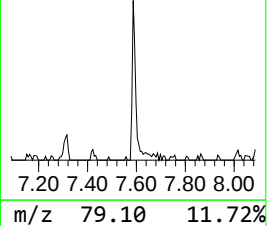
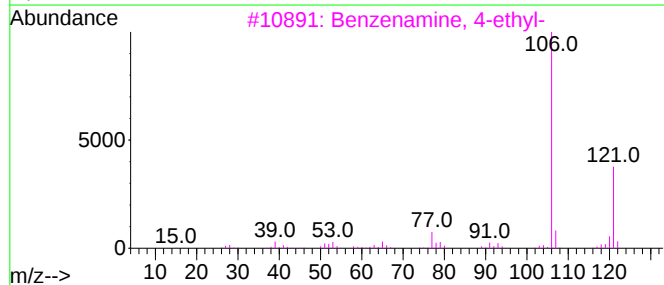
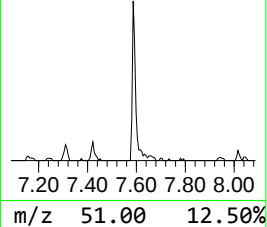
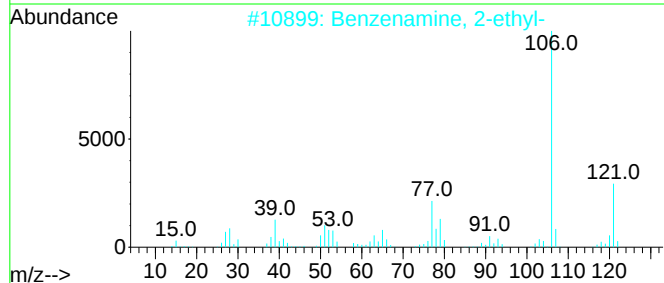
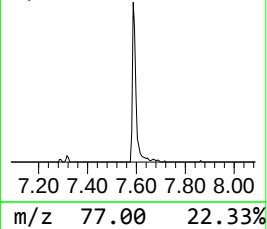
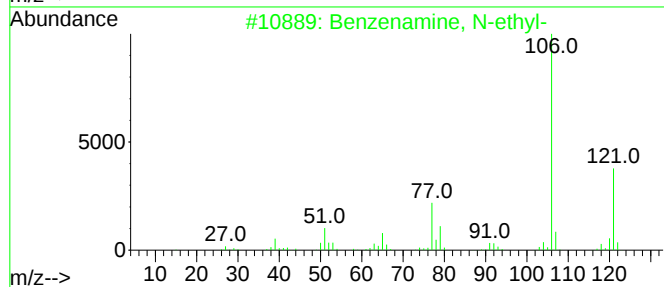
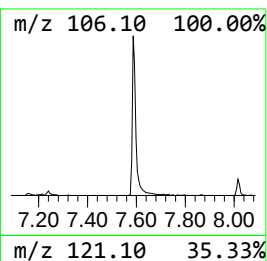
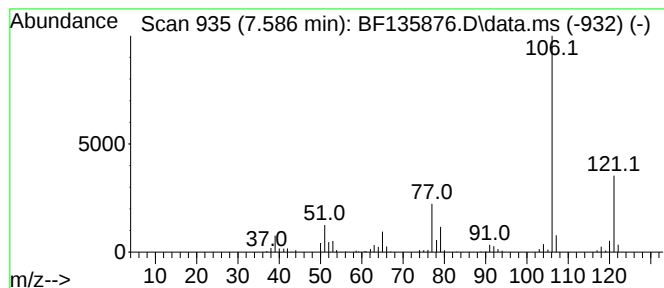
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Benzenamine, N-ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.586	7.43 ng	178608	Naphthalene-d8	8.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, N-ethyl-	121	C8H11N	000103-69-5	95
2			Benzenamine, 2-ethyl-	121	C8H11N	000578-54-1	90
3			Benzenamine, 4-ethyl-	121	C8H11N	000589-16-2	80
4			4-Isopropylpyridine	121	C8H11N	000696-30-0	80
5			Aniline, N-methyl-	107	C7H9N	000100-61-8	72



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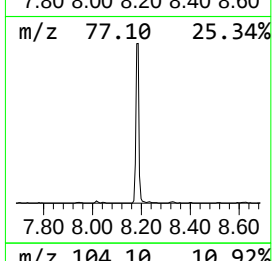
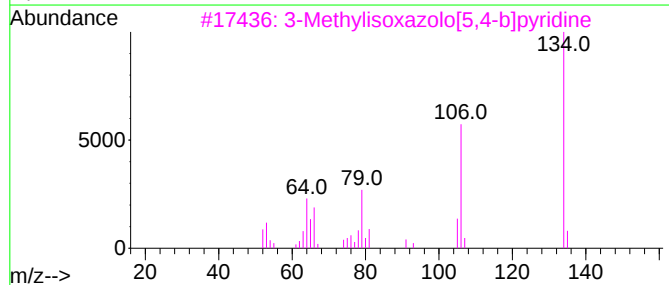
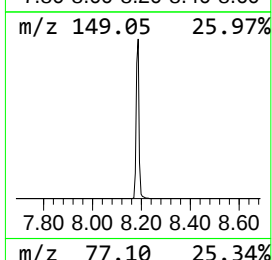
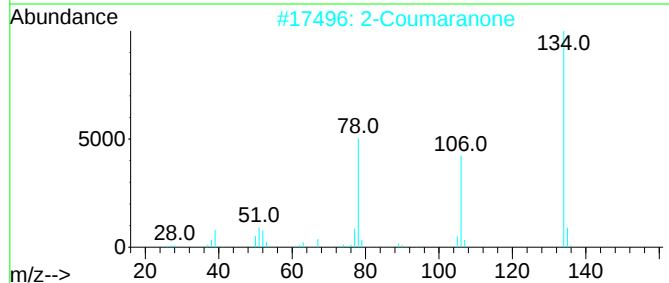
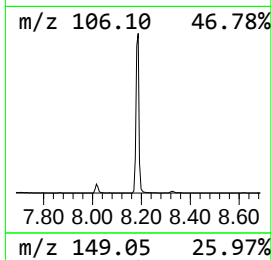
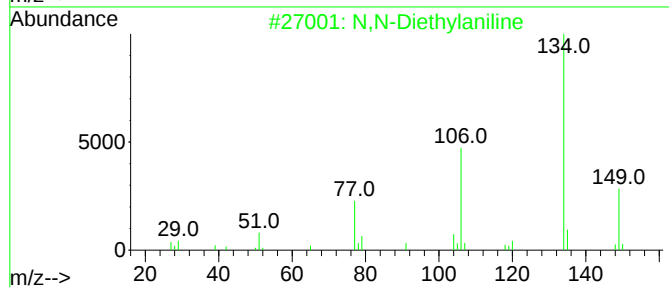
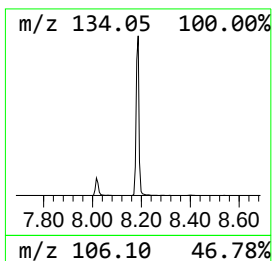
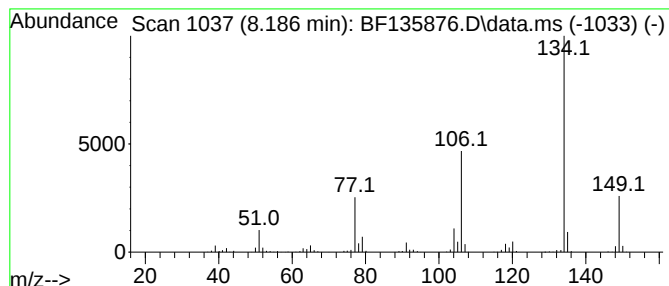
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 N,N-Diethylaniline Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.186	25.30 ng	608470	Naphthalene-d8	8.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N,N-Diethylaniline	149	C10H15N	000091-66-7	97
2			2-Coumaranone	134	C8H6O2	000553-86-6	58
3			3-Methylisoxazolo[5,4-b]pyridine	134	C7H6N2O	058035-50-0	53
4			3-Pyridinecarbonitrile, 1,4-dihy...	134	C7H6N2O	000767-98-6	50
5			Benzenamine, 3-(1,1-dimethylethyl)-	149	C10H15N	005369-19-7	50



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Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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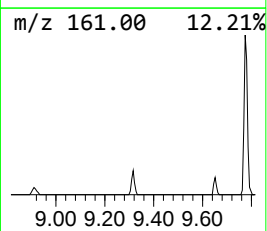
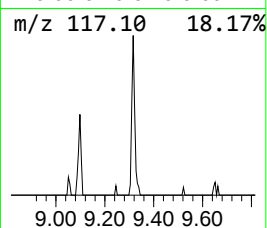
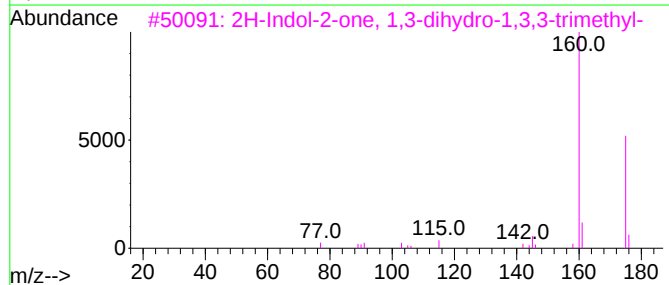
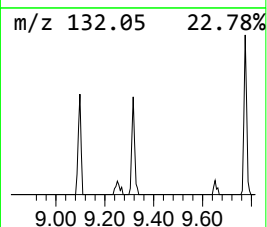
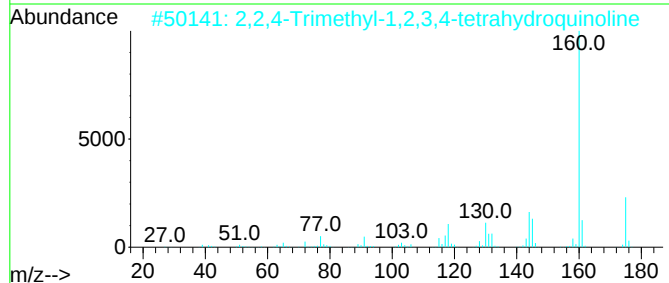
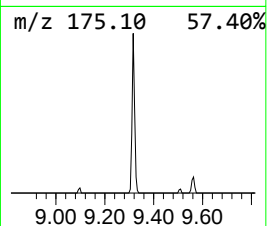
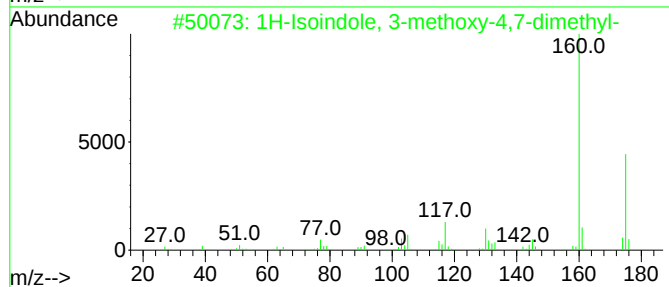
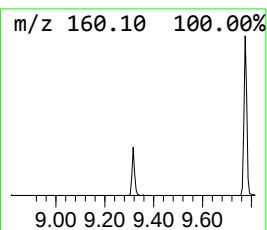
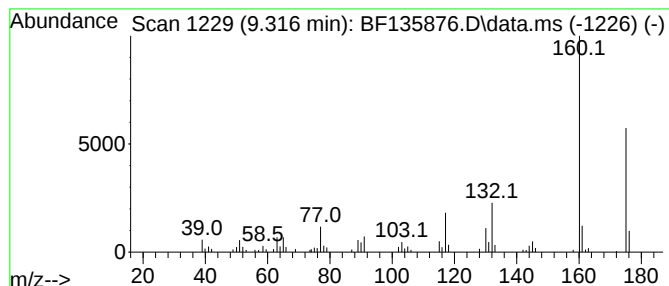
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 1H-Isoindole, 3-methoxy-4,7... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.316	2.18 ng	48125	Acenaphthene-d10	9.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Isoindole, 3-methoxy-4,7-dime...	175	C11H13NO	100813-60-3	87
2			2,2,4-Trimethyl-1,2,3,4-tetrahyd...	175	C12H17N	004497-58-9	80
3			2H-Indol-2-one, 1,3-dihydro-1,3,...	175	C11H13NO	020200-86-6	68
4			1H-Isoindole-1,3(2H)-dione, N-et...	175	C10H9NO2	005022-29-7	64
5			2,6-Diethylphenyl isocyanate	175	C11H13NO	020458-99-5	59



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ClientSampleId :
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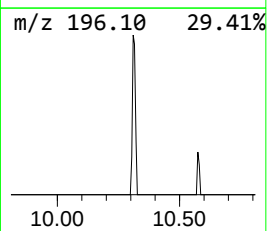
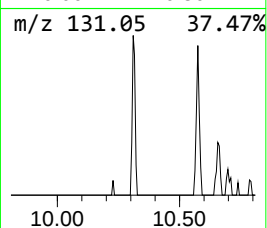
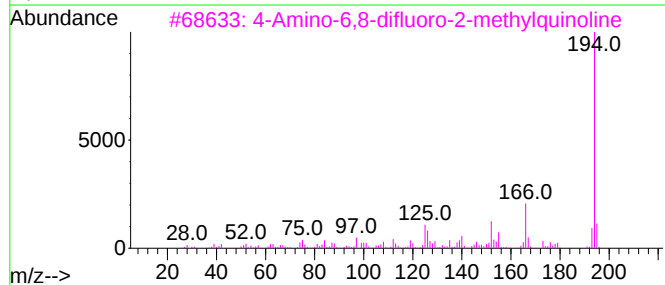
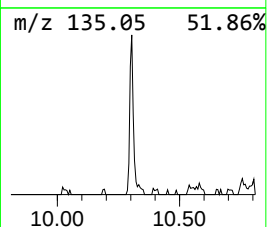
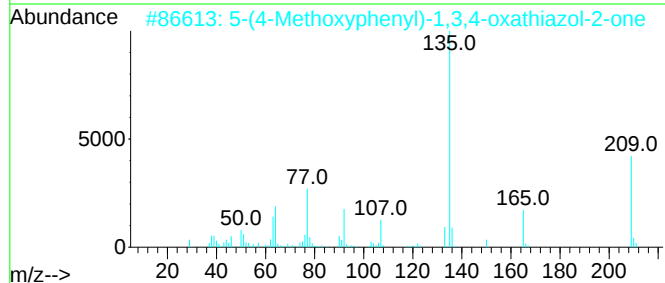
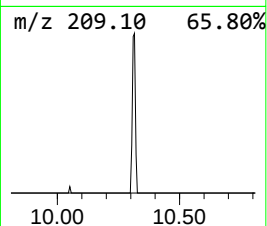
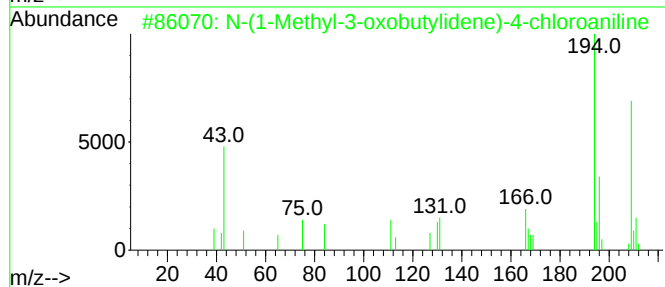
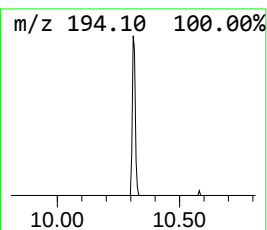
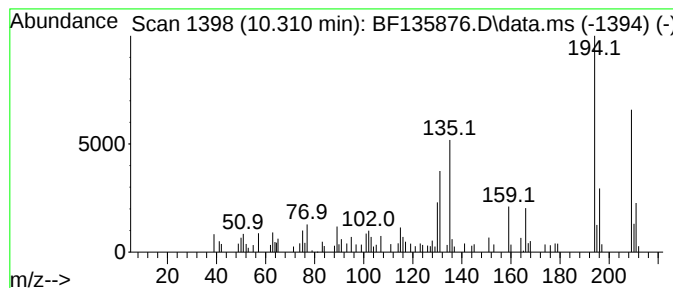
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Peak Number 5 N-(1-Methyl-3-oxobutylidene... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.310	2.06 ng	45661	Acenaphthene-d10	9.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N-(1-Methyl-3-oxobutylidene)-4-c...	209	C11H12ClNO	050519-24-9	64
2			5-(4-Methoxyphenyl)-1,3,4-oxathi...	209	C9H7NO3S	052533-09-2	35
3			4-Amino-6,8-difluoro-2-methylqui...	194	C10H8F2N2	288151-32-6	25
4			2-Chlorobenzo[d]thiazole-6-carbo...	194	C8H3ClN2S	080945-83-1	25
5			N-(3,4-Dimethoxy-alpha-methylben...	209	C11H15NO3	083834-92-8	25



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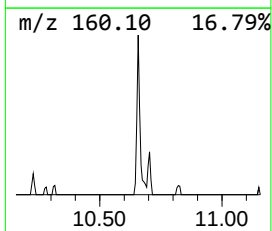
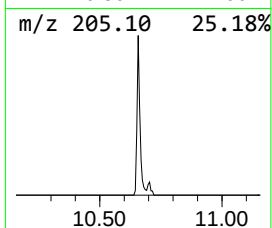
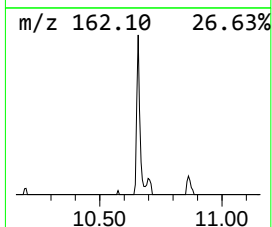
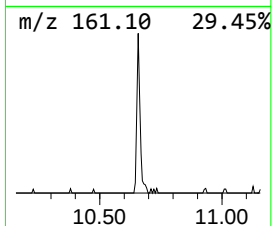
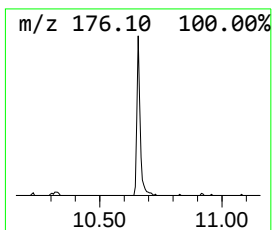
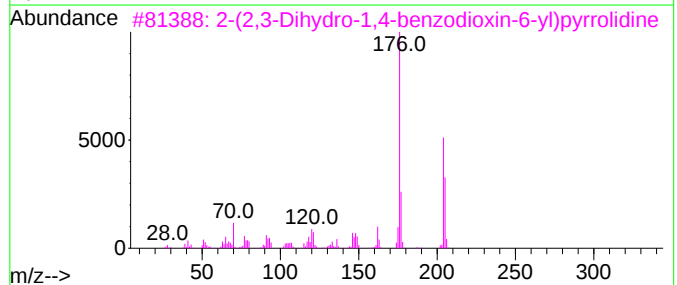
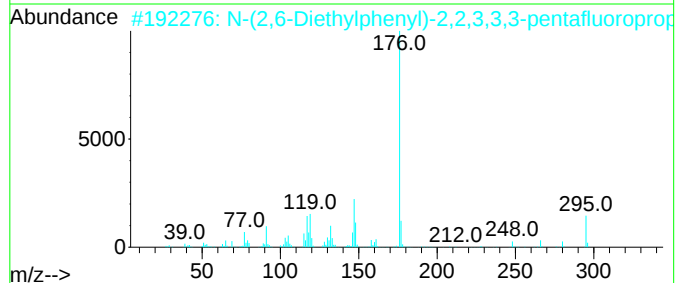
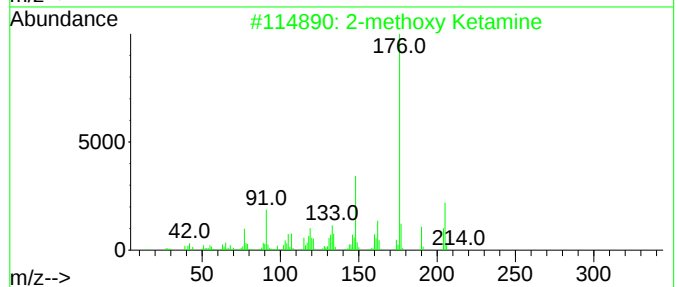
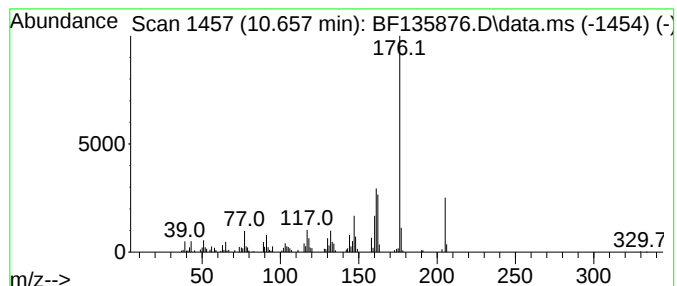
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2-methoxy Ketamine Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.657	5.18 ng	103813	Phenanthrene-d10	11.275

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-methoxy Ketamine	233	C14H19NO2	007063-51-6	53
2			N-(2,6-Diethylphenyl)-2,2,3,3,3-...	295	C13H14F5NO	1000467-22-6	47
3			2-(2,3-Dihydro-1,4-benzodioxin-6-...	205	C12H15NO2	524674-08-6	46
4			2-Cyanomethyl-2-methyl-8-methoxy...	216	C13H16N2O	1000287-23-5	45
5			4-Methoxy-7-methylindan-1-one	176	C11H12O2	103988-25-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101823\
Data File : BF135876.D
Acq On : 19 Oct 2023 06:00
Operator : CG\JU
Sample : 04740-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW4

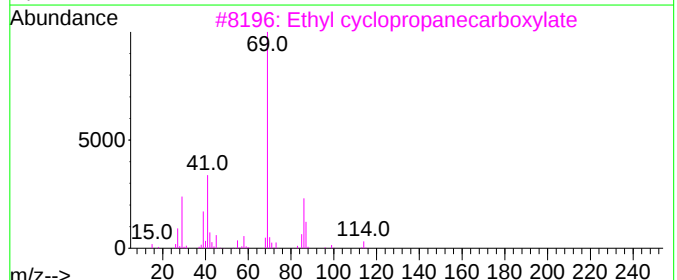
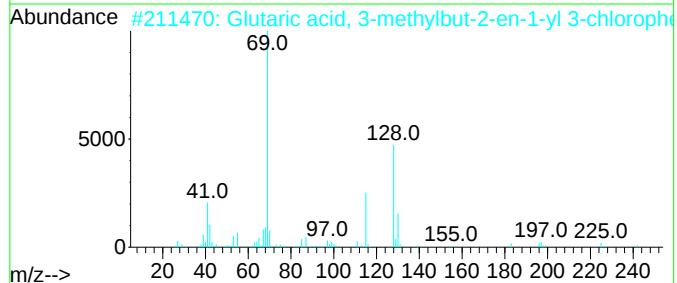
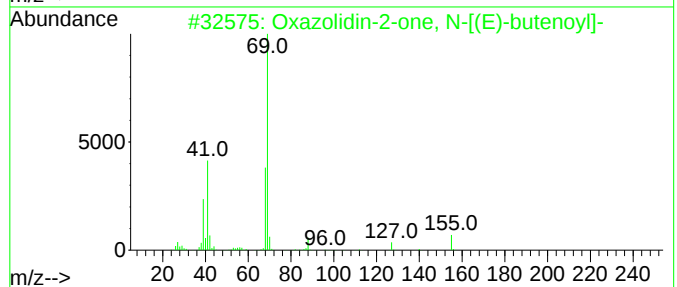
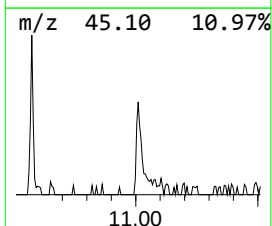
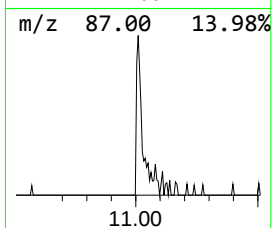
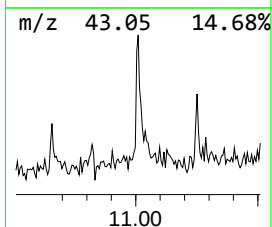
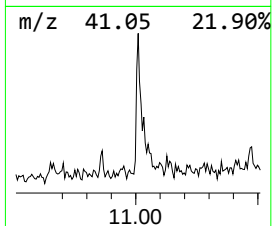
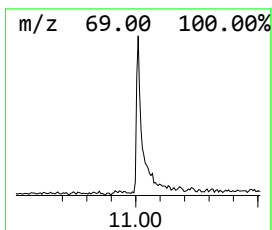
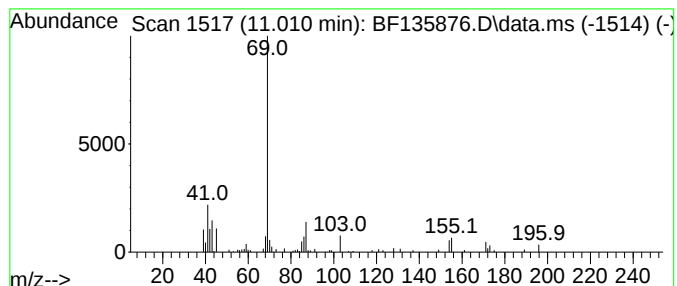
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Oxazolidin-2-one, N-[(E)-bu... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.010	3.42 ng	68551	Phenanthrene-d10	11.275

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxazolidin-2-one, N-[(E)-butenoyl]-	155	C7H9NO3	1010156-45-5	52
2			Glutaric acid, 3-methylbut-2-en-...	310	C16H19ClO4	1000390-45-9	38
3			Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	38
4			Cyclopropanecarboxylic acid, 2-m...	170	C10H18O2	1000354-65-8	36
5			Succinic acid, 3-methylbut-2-en-...	296	C15H17ClO4	1000389-79-8	33



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101823\
Data File : BF135876.D
Acq On : 19 Oct 2023 06:00
Operator : CG\JU
Sample : 04740-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW4

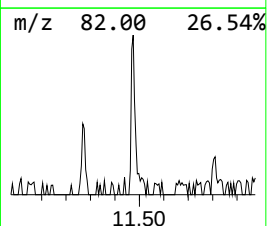
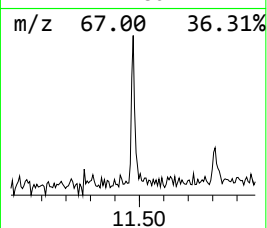
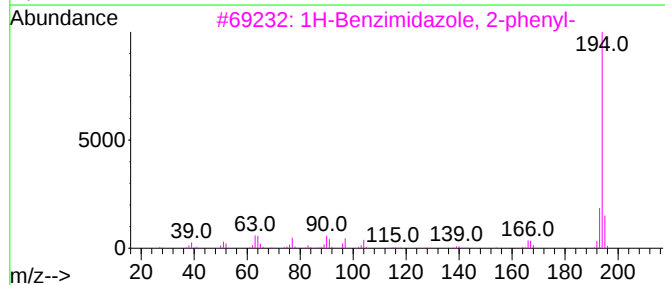
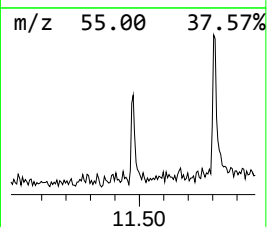
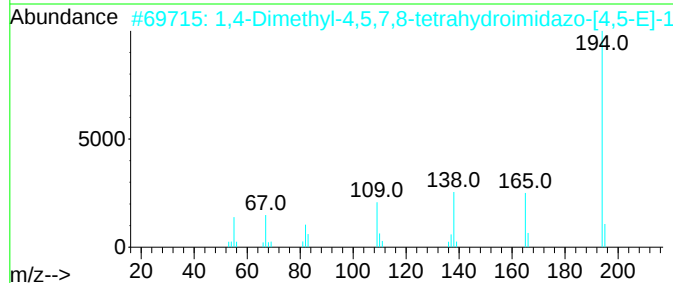
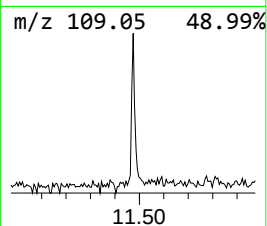
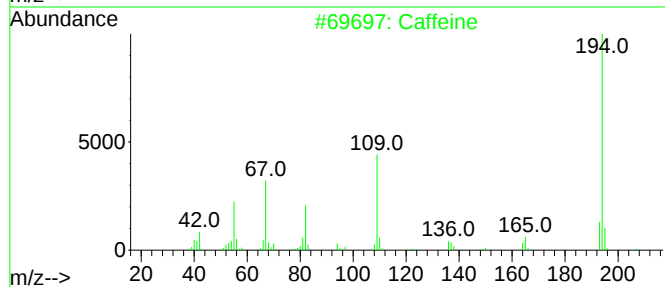
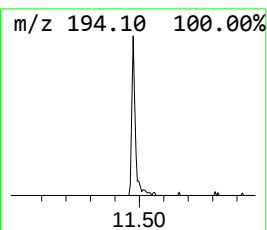
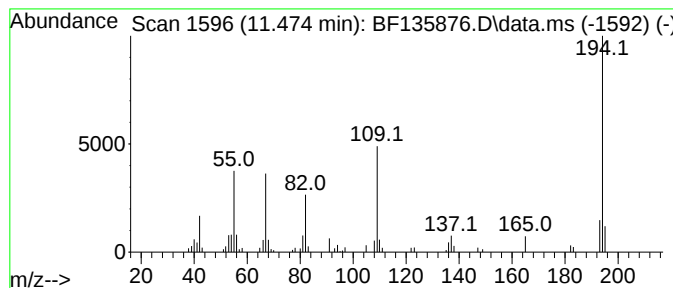
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Caffeine Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.474	2.10 ng	41977	Phenanthrene-d10	11.275

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Caffeine	194	C8H10N4O2	000058-08-2	96
2			1,4-Dimethyl-4,5,7,8-tetrahydroi...	194	C8H10N4O2	130063-15-9	50
3			1H-Benzimidazole, 2-phenyl-	194	C13H10N2	000716-79-0	43
4			2,5-cyclohexadiene-1,4-dione, 2,...	194	C10H14N2O2	1000400-30-4	38
5			6-Amino-1,3-dimethyl-1,6-dihydro...	194	C8H10N4O2	1000301-21-5	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101823\
Data File : BF135876.D
Acq On : 19 Oct 2023 06:00
Operator : CG\JU
Sample : 04740-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

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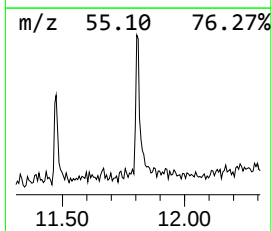
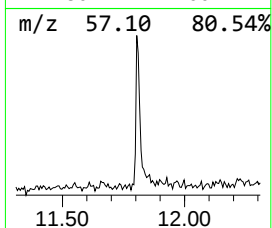
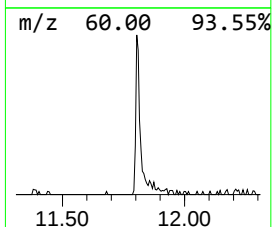
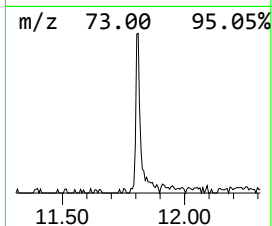
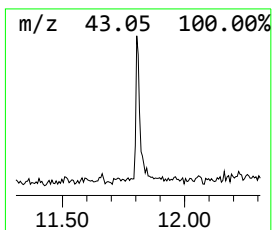
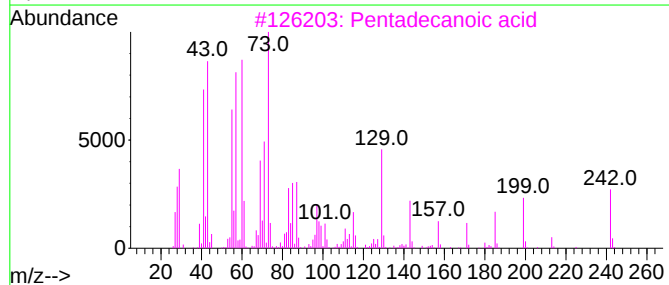
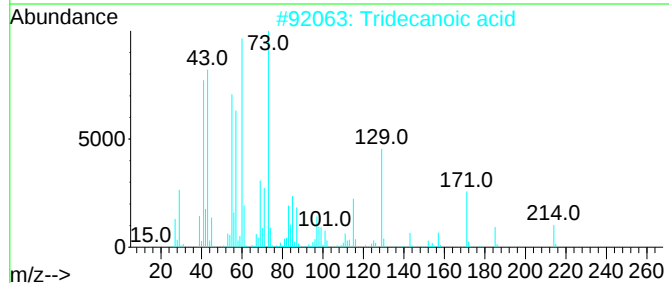
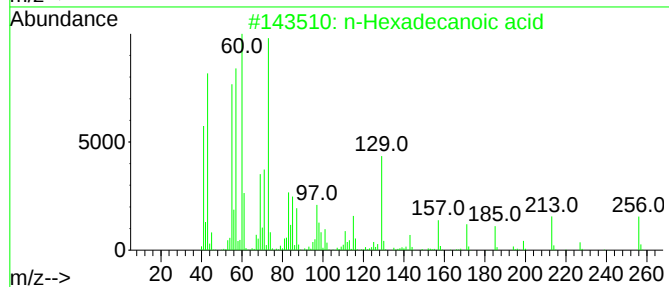
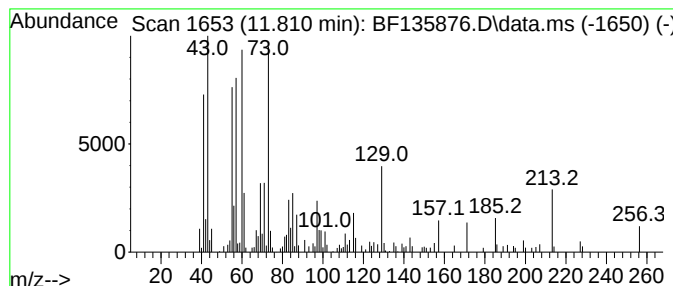
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.810	5.89 ng	117964	Phenanthrene-d10	11.275

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Tridecanoic acid	214	C13H26O2	000638-53-9	93
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	72
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	64
5			Dodecanoic acid	200	C12H24O2	000143-07-7	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101823\
Data File : BF135876.D
Acq On : 19 Oct 2023 06:00
Operator : CG\JU
Sample : 04740-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

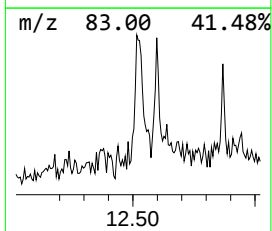
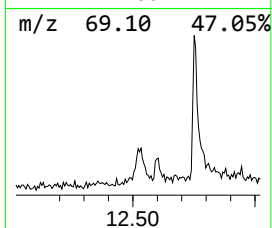
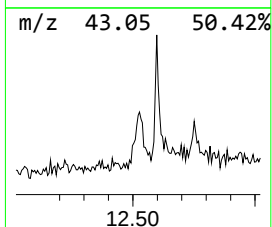
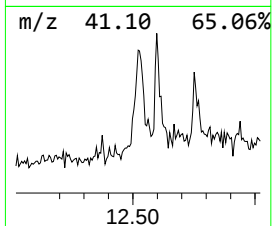
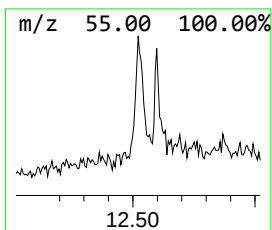
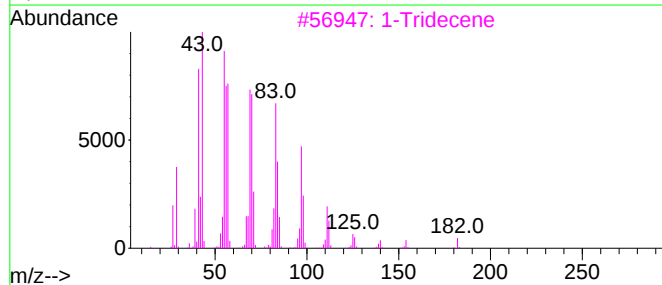
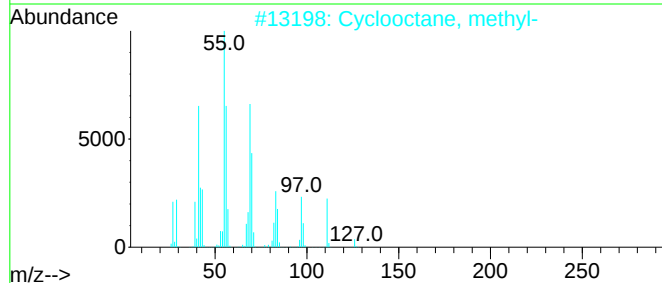
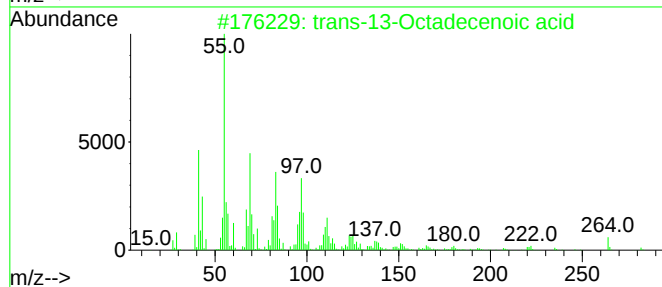
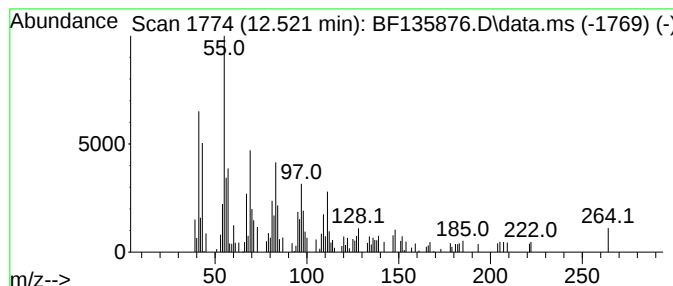
Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 trans-13-Octadecenoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.521	4.63 ng	92671	Phenanthrene-d10	11.275	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	trans-13-Octadecenoic acid	282	C18H34O2	000693-71-0	72
2	Cyclooctane, methyl-	126	C9H18	001502-38-1	64
3	1-Tridecene	182	C13H26	002437-56-1	53
4	Z-8-Methyl-9-tetradecenoic acid	240	C15H28O2	1000130-84-5	53
5	9-Oxabicyclo[6.1.0]nonane	126	C8H14O	000286-62-4	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101823\
Data File : BF135876.D
Acq On : 19 Oct 2023 06:00
Operator : CG\JU
Sample : 04740-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

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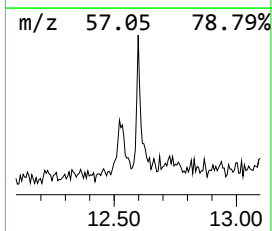
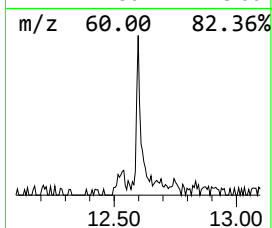
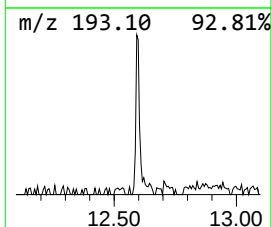
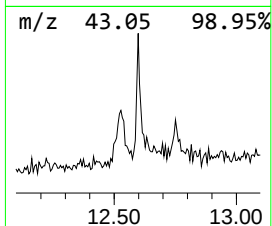
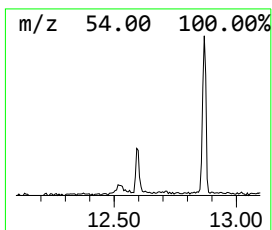
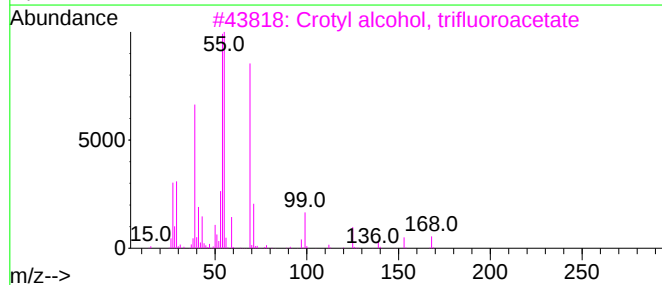
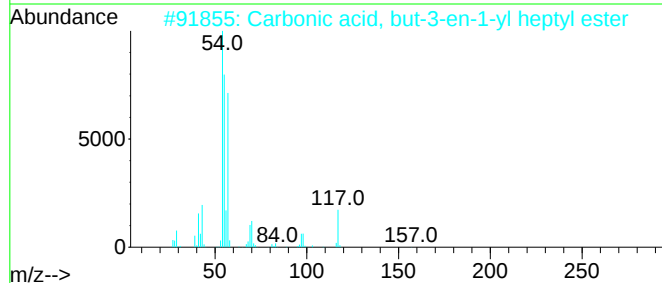
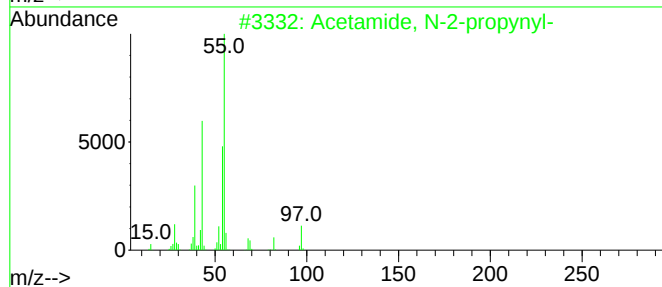
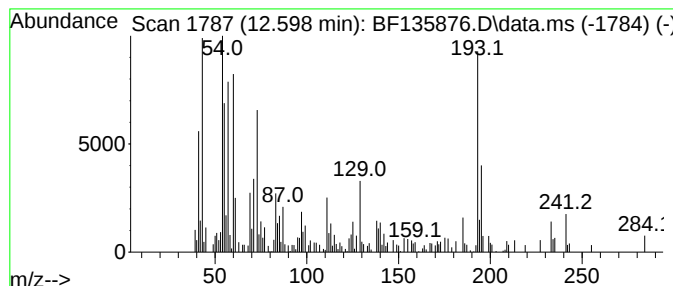
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown12.598 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.598	4.66 ng	93283	Phenanthrene-d10	11.275

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetamide, N-2-propynyl-	97	C5H7NO	065881-41-6	43
2			Carbonic acid, but-3-en-1-yl hep...	214	C12H22O3	1000383-22-8	38
3			Crotyl alcohol, trifluoroacetate	168	C6H7F3O2	1000352-26-6	38
4			Carbonic acid, but-3-en-1-yl und...	270	C16H30O3	1010383-23-2	38
5			Aziridine, 2-methylene-1-(1-meth...	97	C6H11N	055268-35-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101823\
Data File : BF135876.D
Acq On : 19 Oct 2023 06:00
Operator : CG\JU
Sample : 04740-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

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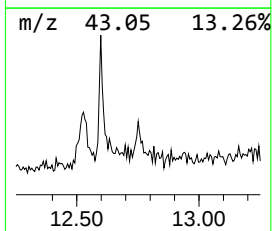
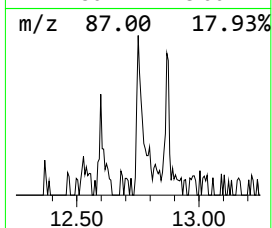
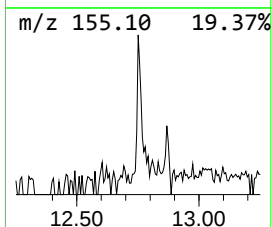
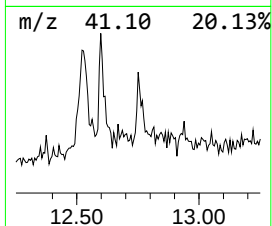
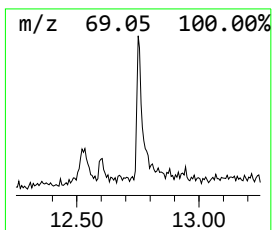
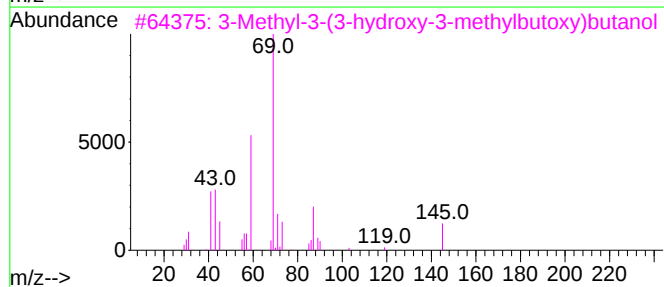
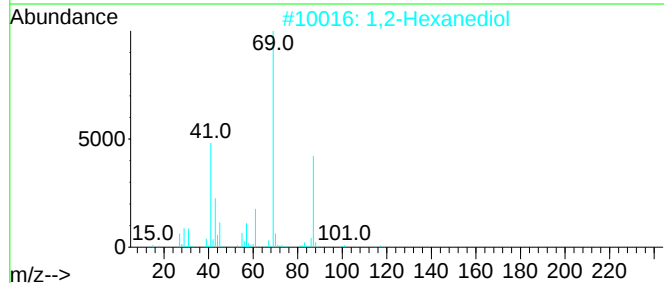
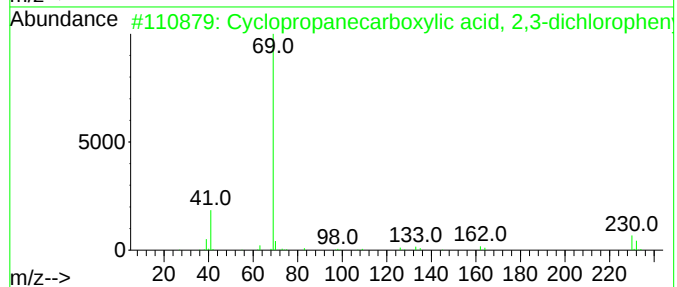
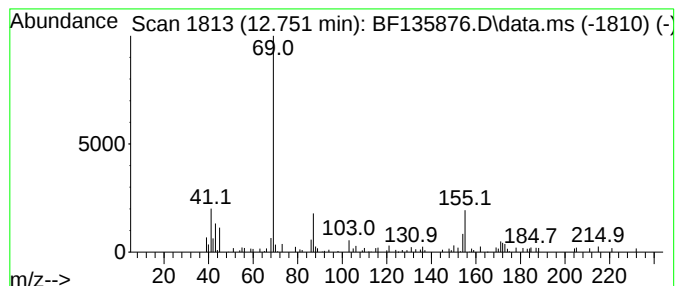
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown12.751 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.751	2.69 ng	48778	Chrysene-d12	13.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropanecarboxylic acid, 2,3...	230	C10H8Cl2O2	959085-02-0	38
2			1,2-Hexanediol	118	C6H14O2	006920-22-5	38
3			3-Methyl-3-(3-hydroxy-3-methylbu...	190	C10H22O3	1000432-16-8	33
4			Geraniol	154	C10H18O	000106-24-1	28
5			Cyclopropanecarboxylic acid, buty...	142	C8H14O2	054947-39-6	28



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101823\
Data File : BF135876.D
Acq On : 19 Oct 2023 06:00
Operator : CG\JU
Sample : 04740-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

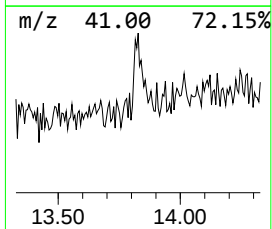
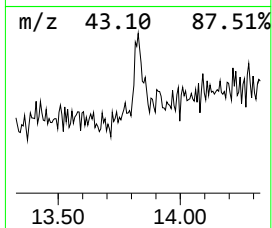
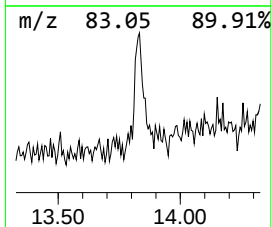
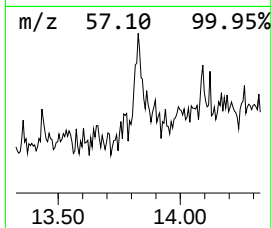
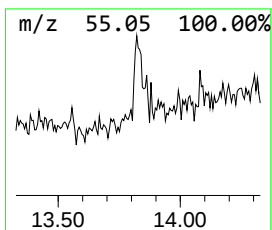
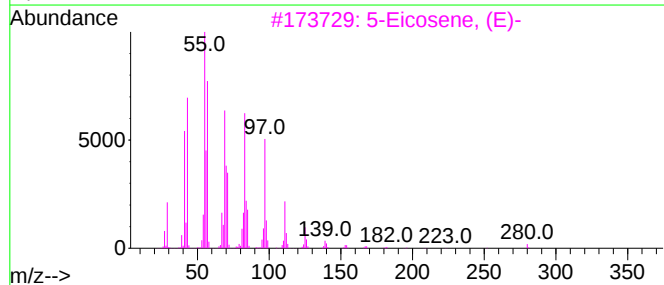
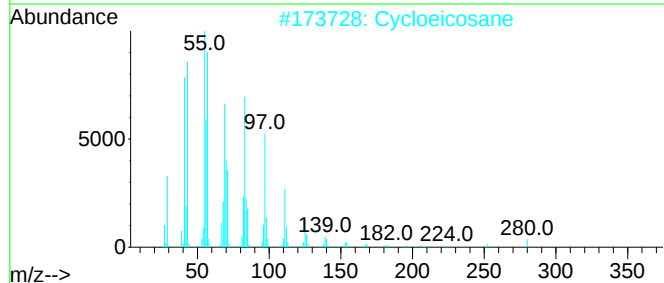
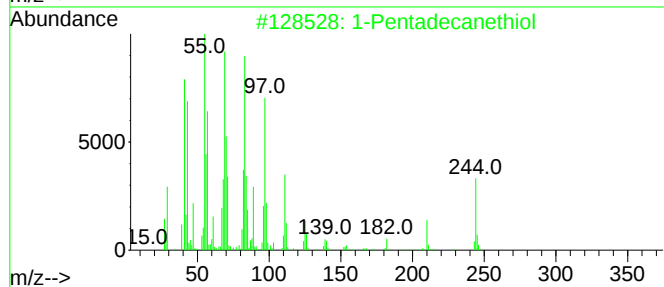
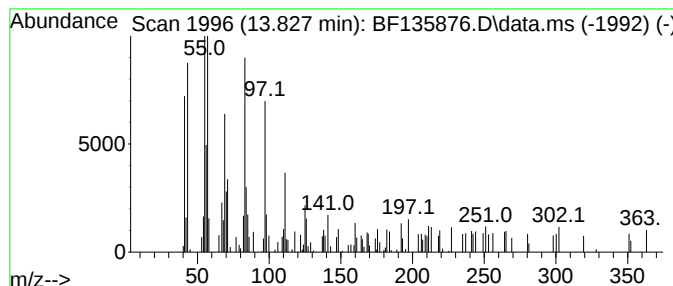
Instrument :
BNA_F
ClientSampleId :
MW4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1-Pentadecanethiol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.827	2.60 ng	47118	Chrysene-d12	13.939	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentadecanethiol	244	C15H32S	025276-70-4	64
2	Cycloeicosane	280	C20H40	000296-56-0	60
3	5-Eicosene, (E)-	280	C20H40	074685-30-6	49
4	1-Eicosanol	298	C20H42O	000629-96-9	49
5	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101823\
Data File : BF135876.D
Acq On : 19 Oct 2023 06:00
Operator : CG\JU
Sample : 04740-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW4

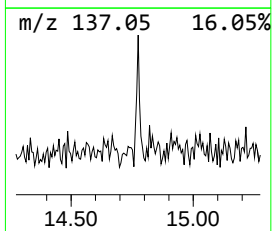
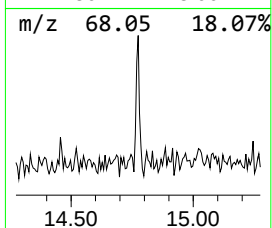
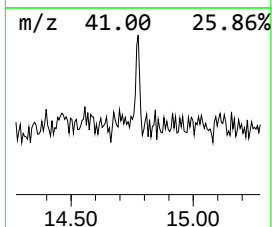
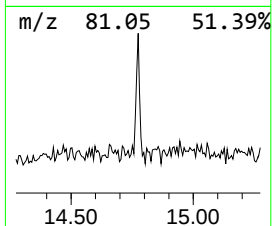
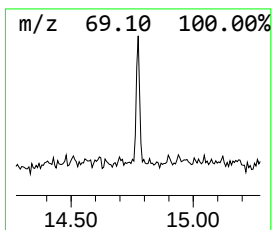
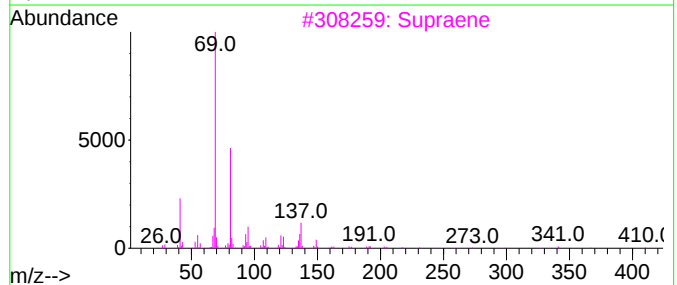
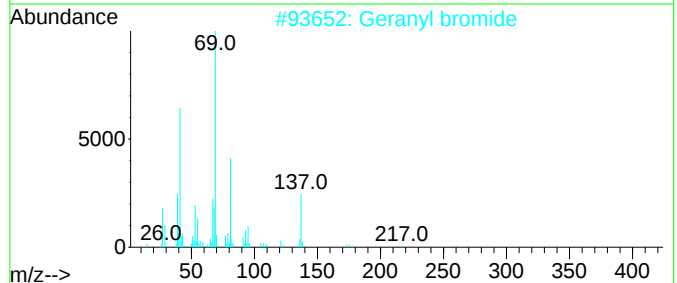
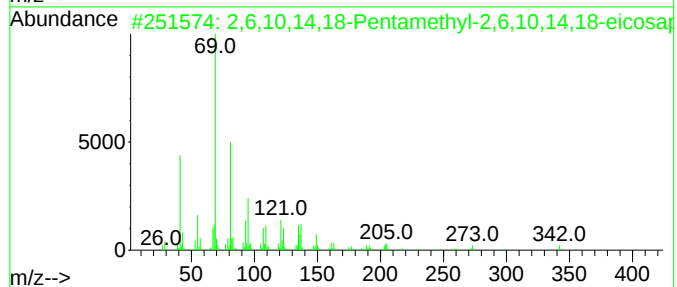
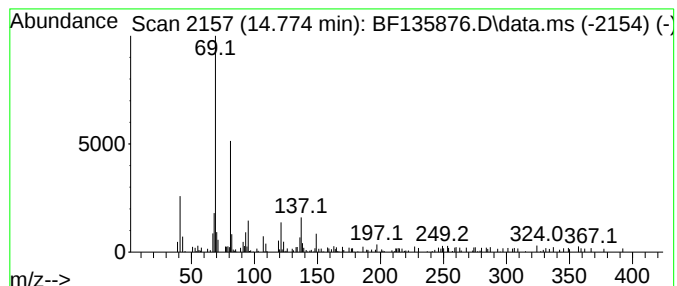
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101323.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 2,6,10,14,18-Pentamethyl-2,... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.774	3.07 ng	46439	Perylene-d12		15.421

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,10,14,18-Pentamethyl-2,6,10,...	342	C25H42		075581-03-2	72
2	Geranyl bromide	216	C10H17Br		006138-90-5	64
3	Supraene	410	C30H50		007683-64-9	64
4	4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O		066408-55-7	64
5	3,7,11-Tridecatrienoic acid, 4,8...	264	C17H28O2		036237-70-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101823\
Data File : BF135876.D
Acq On : 19 Oct 2023 06:00
Operator : CG\JU
Sample : 04740-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101323.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
2-Pentanone, 4-...	4.963	3.1	ng	60968	1	6.739	395224	20.0
Benzenamine, N-...	7.586	7.4	ng	178608	2	8.016	480964	20.0
N,N-Diethylaniline	8.186	25.3	ng	608470	2	8.016	480964	20.0
1H-Isoindole, 3...	9.316	2.2	ng	48125	3	9.775	442290	20.0
N-(1-Methyl-3-o...	10.310	2.1	ng	45661	3	9.775	442290	20.0
2-methoxy Ketamine	10.657	5.2	ng	103813	4	11.275	400731	20.0
Oxazolidin-2-on...	11.010	3.4	ng	68551	4	11.275	400731	20.0
Caffeine	11.474	2.1	ng	41977	4	11.275	400731	20.0
n-Hexadecanoic ...	11.810	5.9	ng	117964	4	11.275	400731	20.0
trans-13-Octade...	12.521	4.6	ng	92671	4	11.275	400731	20.0
unknown12.598	12.598	4.7	ng	93283	4	11.275	400731	20.0
unknown12.751	12.751	2.7	ng	48778	5	13.939	362635	20.0
1-Pentadecanethiol	13.827	2.6	ng	47118	5	13.939	362635	20.0
2,6,10,14,18-Pe...	14.774	3.1	ng	46439	6	15.421	302206	20.0