

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061323\
 Data File : BF133731.D
 Acq On : 13 Jun 2023 15:16
 Operator : CG\JU
 Sample : 03094-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-1

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-BF060923.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF133731.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.210	16	21	28	rBV	74827	96582	3.05%	0.497%
2	5.451	564	572	575	rBV	1287584	1305046	41.20%	6.710%
3	6.269	705	711	714	rBV	131034	130055	4.11%	0.669%
4	6.463	740	744	747	rVB	1340566	1164544	36.76%	5.988%
5	6.834	803	807	810	rBV	641701	508579	16.05%	2.615%
6	7.392	899	902	909	rVB	864585	758048	23.93%	3.898%
7	8.116	1022	1025	1028	rVB	674291	503378	15.89%	2.588%
8	8.998	1172	1175	1178	rBV4	82056	98265	3.10%	0.505%
9	9.139	1196	1199	1203	rVB	216525	186027	5.87%	0.957%
10	9.192	1203	1208	1211	rBV	1610714	1356519	42.82%	6.975%
11	9.275	1218	1222	1224	rBV2	279670	374174	11.81%	1.924%
12	9.422	1245	1247	1249	rBV2	128304	110680	3.49%	0.569%
13	9.445	1249	1251	1253	rVV2	169510	152374	4.81%	0.783%
14	9.586	1272	1275	1276	rBV	550827	570362	18.00%	2.933%
15	9.639	1281	1284	1288	rVV3	299019	410215	12.95%	2.109%
16	9.745	1299	1302	1305	rBV2	607006	831905	26.26%	4.278%
17	9.786	1307	1309	1312	rVB	996560	914091	28.86%	4.700%
18	9.827	1314	1316	1318	rBV2	335312	304915	9.63%	1.568%
19	9.927	1329	1333	1335	rBV4	426602	732067	23.11%	3.764%
20	10.292	1392	1395	1400	rBV	1211903	1210881	38.22%	6.226%
21	10.545	1435	1438	1442	rVB	1236480	1225397	38.68%	6.301%
22	10.816	1481	1484	1488	rVB	2645184	2935791	92.68%	15.096%
23	11.292	1562	1565	1568	rVB	2750680	3167815	100.00%	16.289%
24	15.492	2276	2279	2289	rVB	293006	400340	12.64%	2.059%

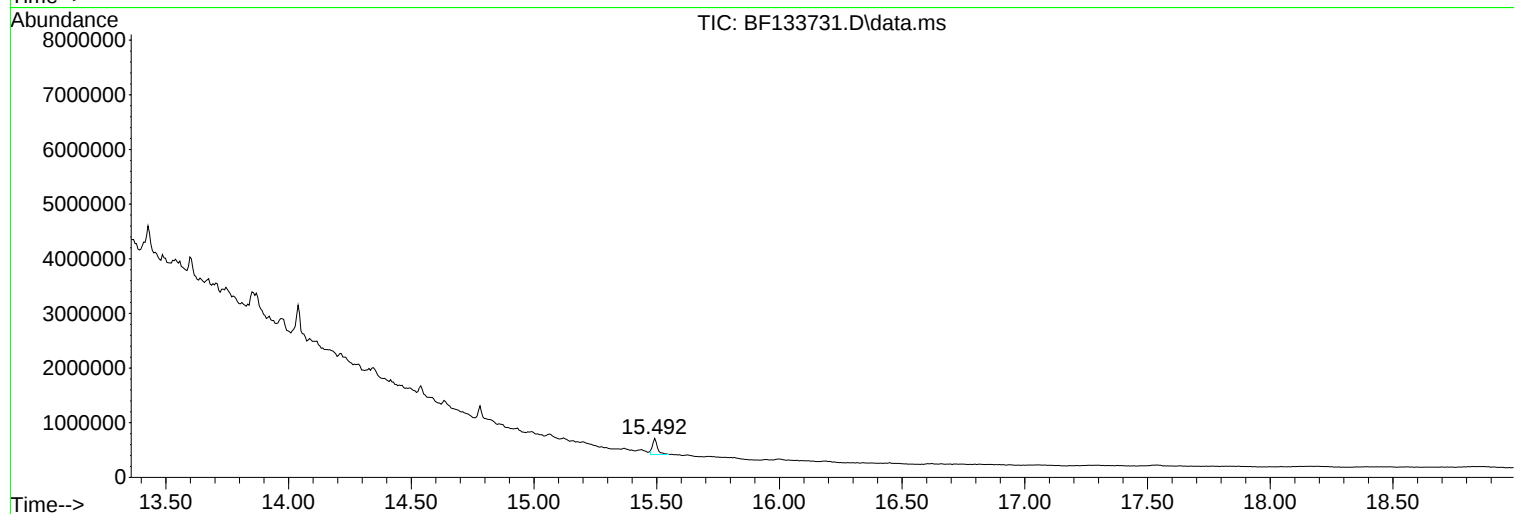
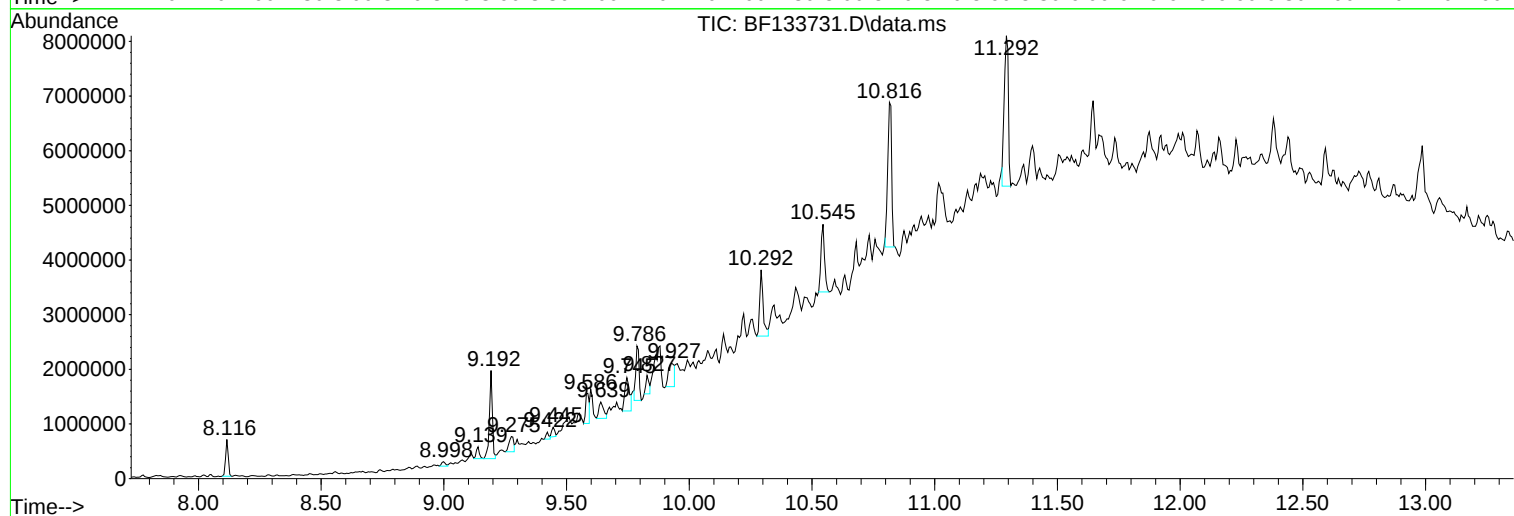
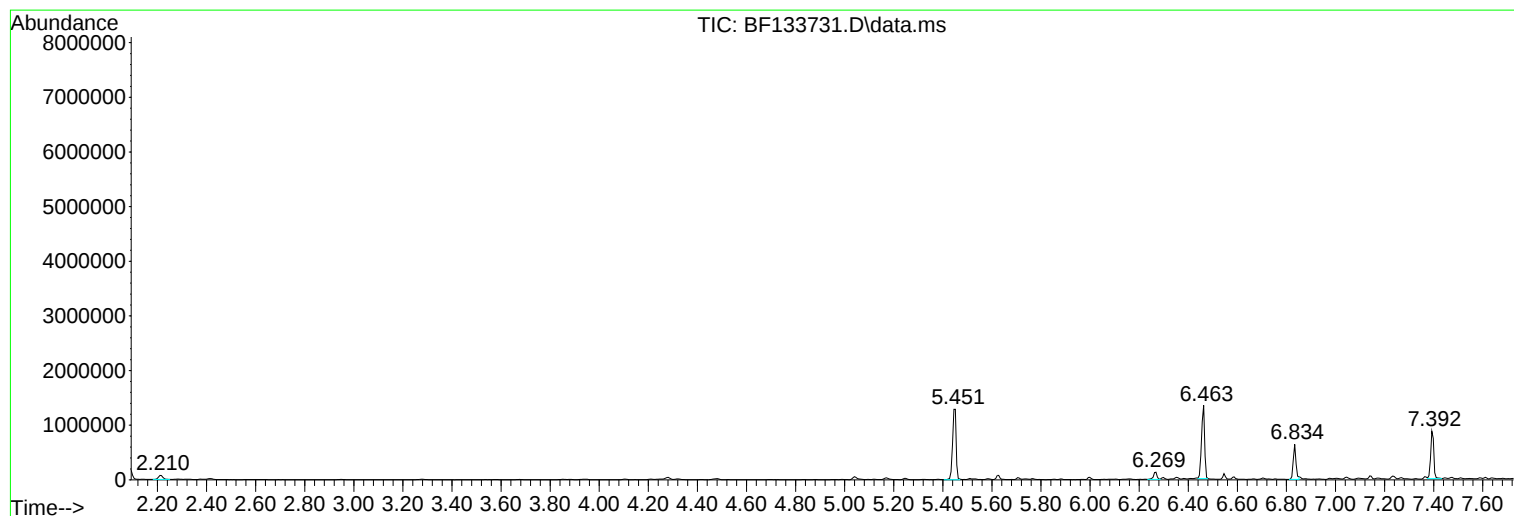
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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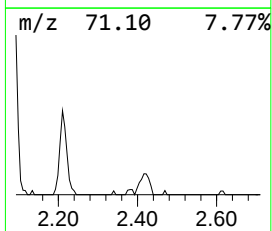
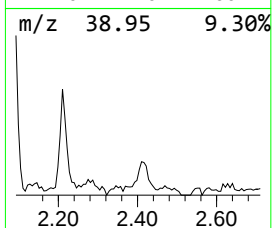
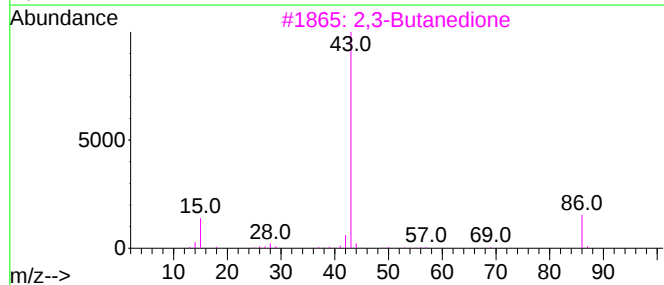
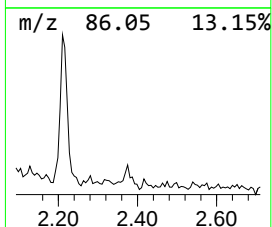
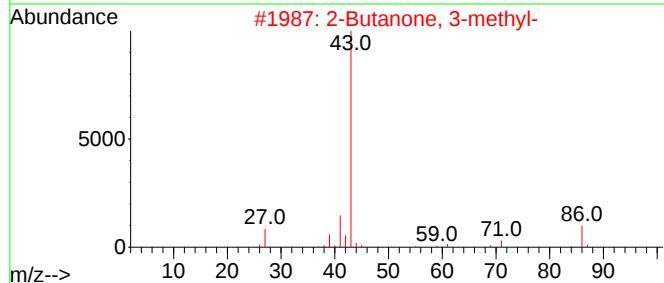
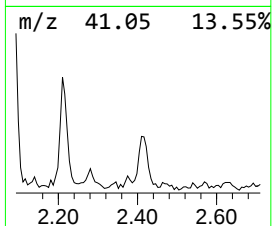
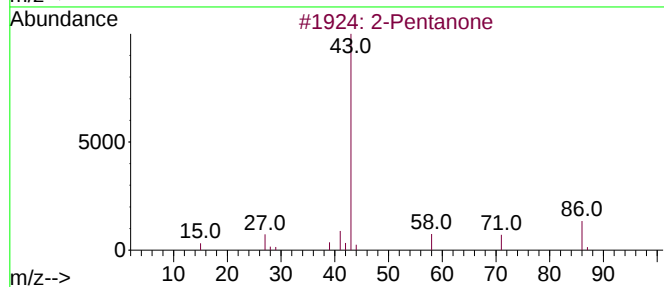
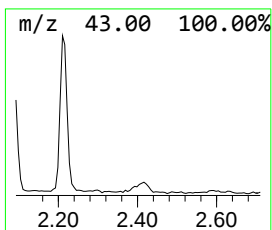
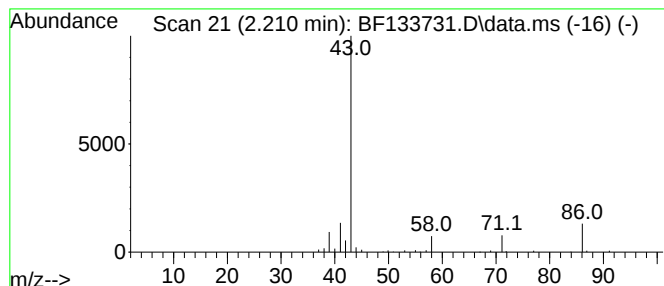
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.210	3.80 ng	96582	1,4-Dichlorobenzene-d4	6.834

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone		86	C5H10O	000107-87-9	64
2	2-Butanone, 3-methyl-		86	C5H10O	000563-80-4	53
3	2,3-Butanedione		86	C4H6O2	000431-03-8	40
4	3-Aminopyrrolidine		86	C4H10N2	079286-79-6	36
5	Ethanone, 1-oxiranyl-		86	C4H6O2	004401-11-0	9



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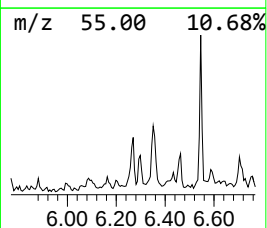
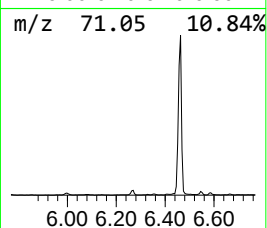
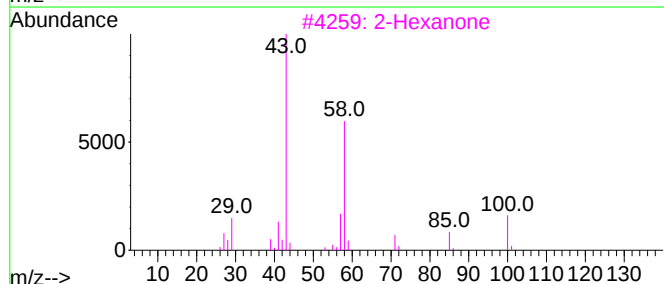
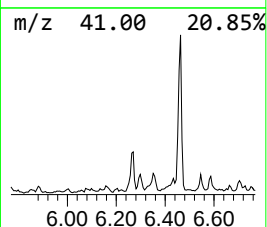
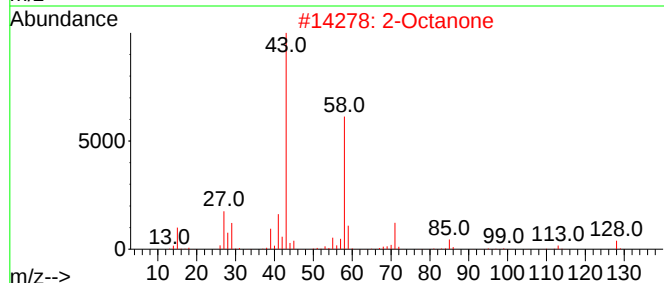
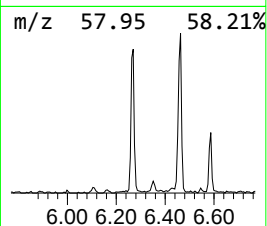
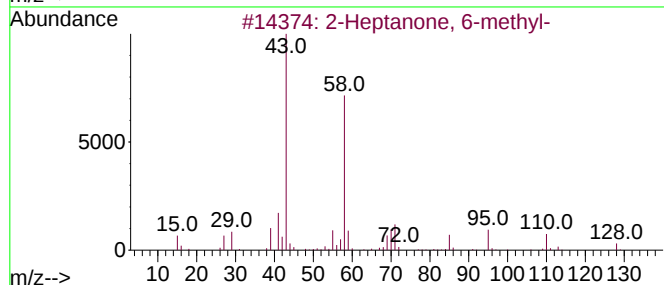
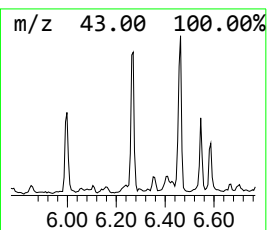
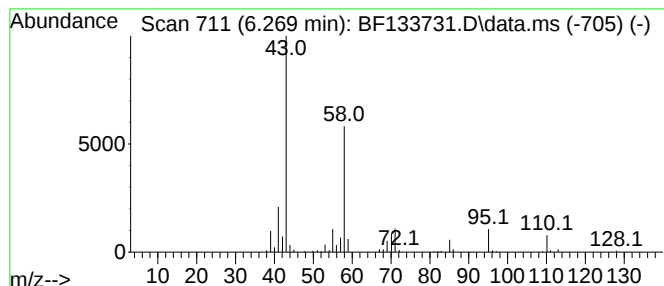
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TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Heptanone, 6-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.269	5.11 ng	130055	1,4-Dichlorobenzene-d4	6.834

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Heptanone, 6-methyl-	128	C8H16O	000928-68-7	86
2		2-Octanone	128	C8H16O	000111-13-7	53
3		2-Hexanone	100	C6H12O	000591-78-6	45
4		2-Hexanone, 4-methyl-	114	C7H14O	000105-42-0	42
5		(Dimethylamino)ethyl methacrylate	157	C8H15NO2	002867-47-2	38



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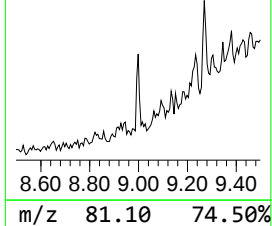
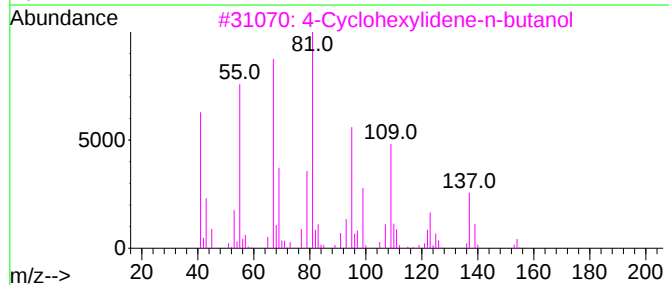
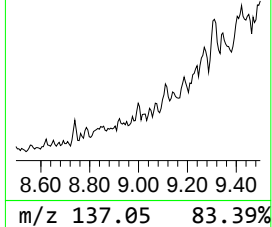
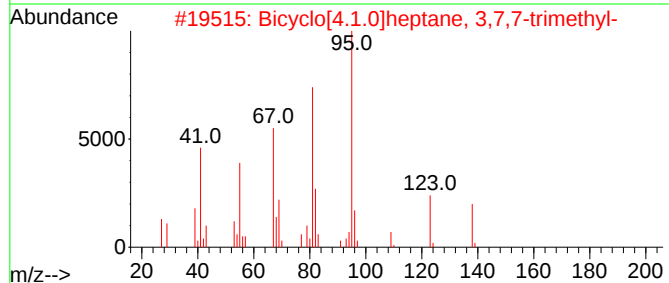
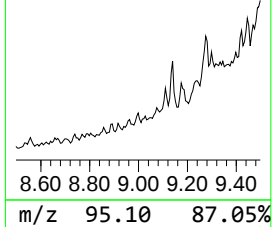
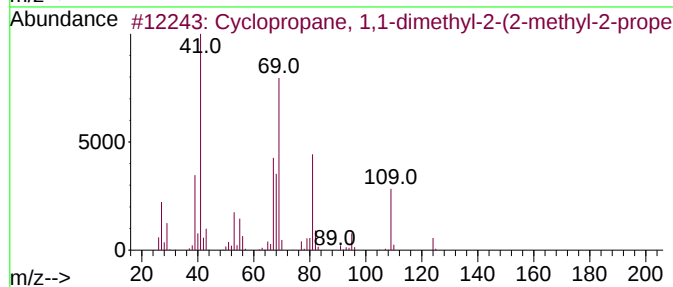
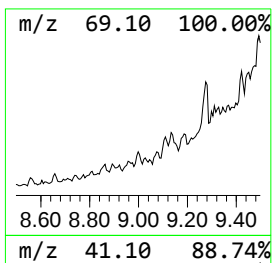
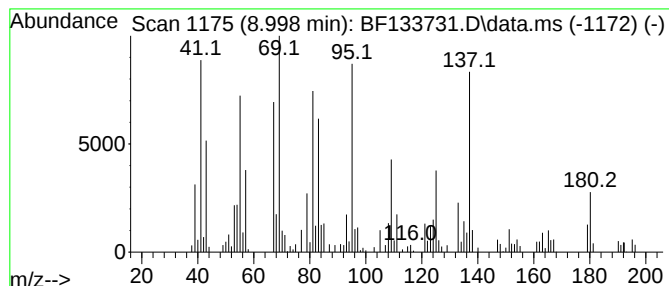
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 unknown8.998 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.998	3.90 ng	98265	Naphthalene-d8	8.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, 1,1-dimethyl-2-(2-...	124	C9H16	069147-03-1	43
2			Bicyclo[4.1.0]heptane, 3,7,7-tri...	138	C10H18	000554-59-6	42
3			4-Cyclohexylidene-n-butanol	154	C10H18O	004441-58-1	42
4			3,5-Dodecadiene, 2-methyl-	180	C13H24	055638-51-2	38
5			2,5-Octadiene, 3,4,5,6-tetramethyl-	166	C12H22	247797-85-9	35



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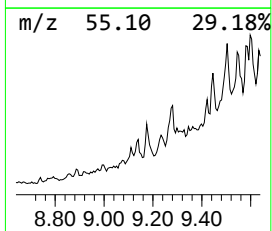
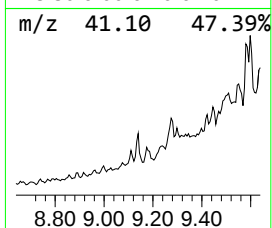
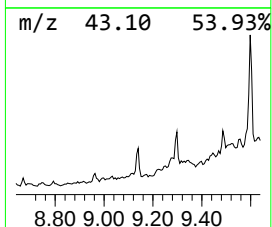
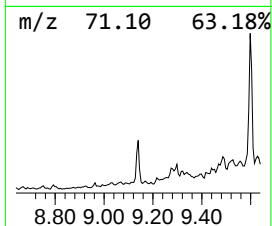
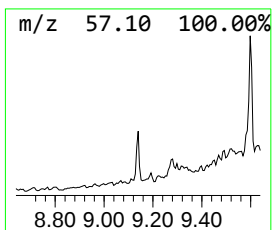
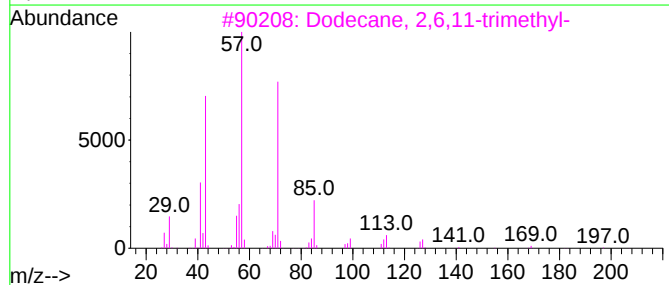
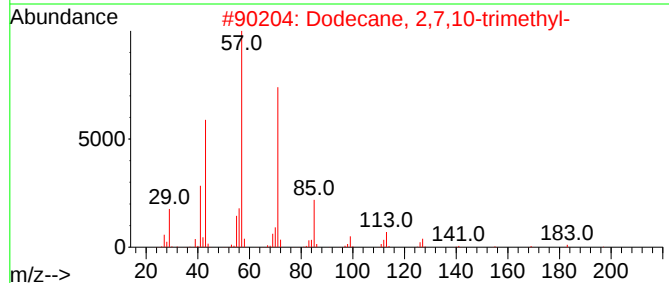
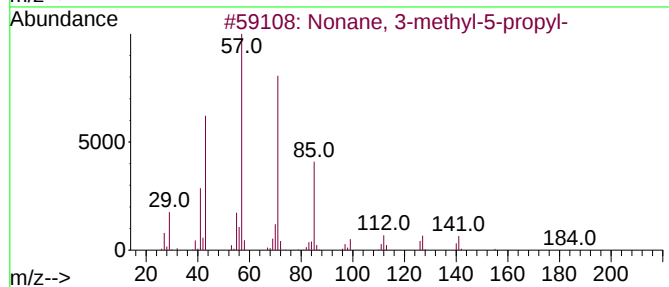
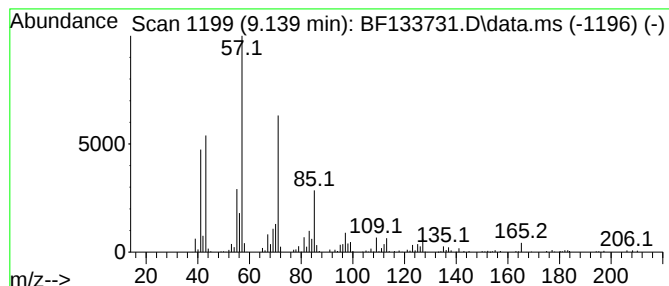
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Peak Number 4 Nonane, 3-methyl-5-propyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.139	5.08 ng	186027	Acenaphthene-d10	9.880

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3-methyl-5-propyl-	184	C13H28	031081-18-2	72
2			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	64
3			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	59
4			Hentriacontane	437	C31H64	000630-04-6	59
5			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	58



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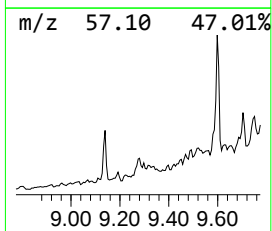
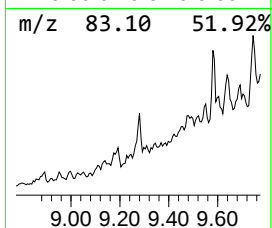
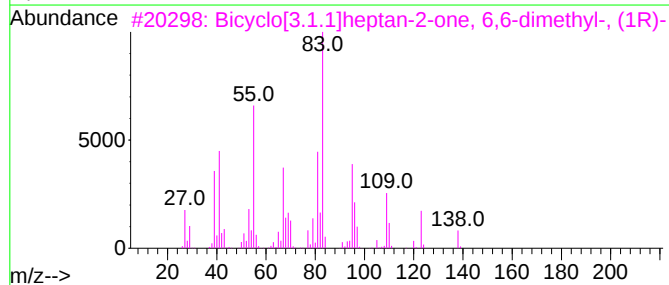
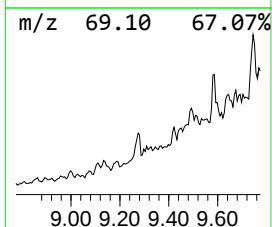
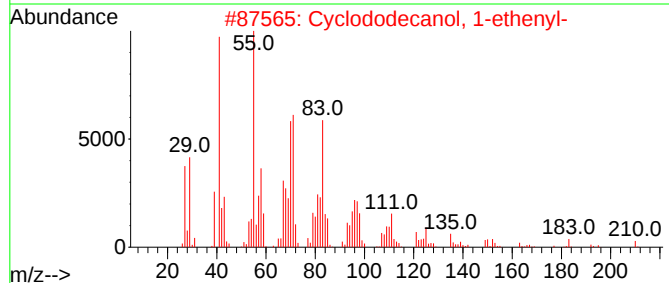
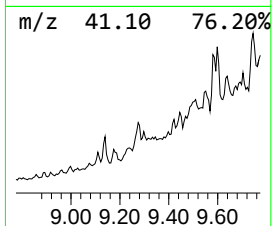
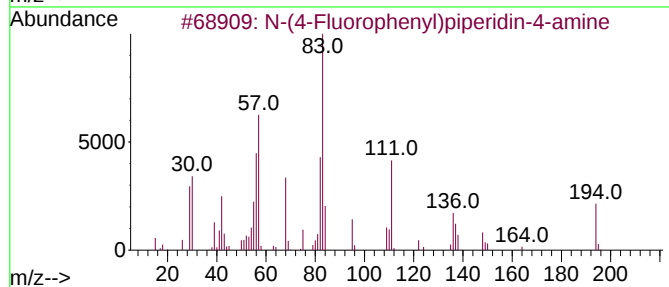
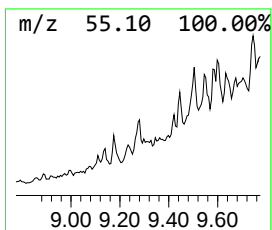
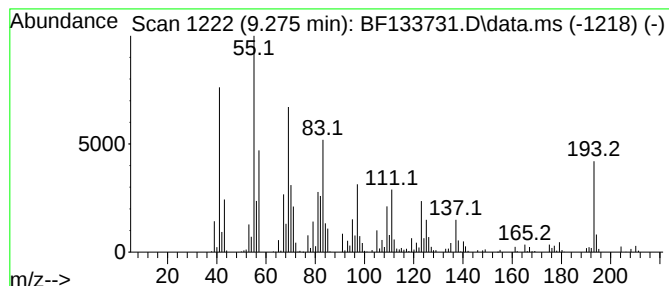
Instrument :
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown9.275 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.275	10.22 ng	374174	Acenaphthene-d10	9.880	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N-(4-Fluorophenyl)piperidin-4-amine	194	C11H15FN2	038043-08-2	46
2	Cyclododecanol, 1-ethenyl-	210	C14H26O	006244-49-1	38
3	Bicyclo[3.1.1]heptan-2-one, 6,6-...	138	C9H14O	038651-65-9	22
4	1-Heneicosanol	312	C21H44O	015594-90-8	18
5	1-Docosene	308	C22H44	001599-67-3	18



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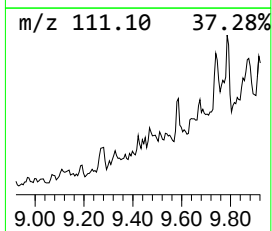
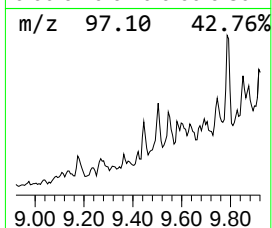
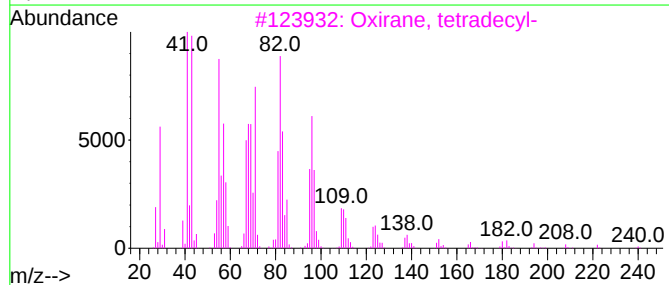
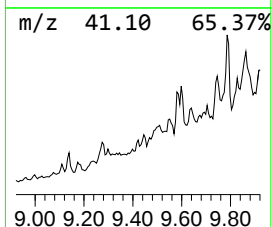
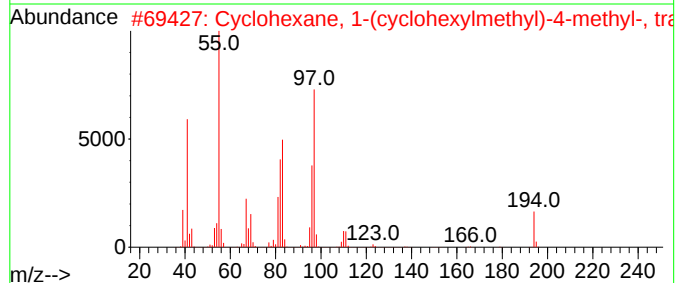
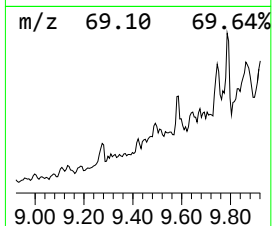
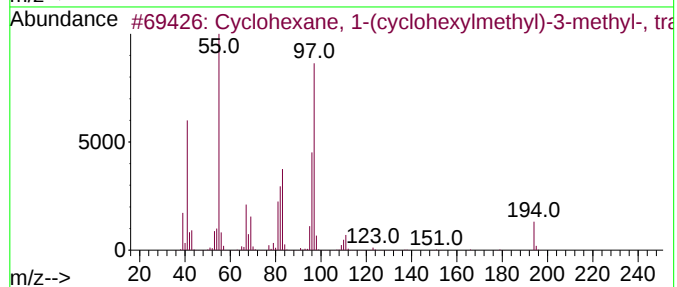
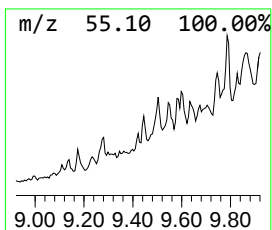
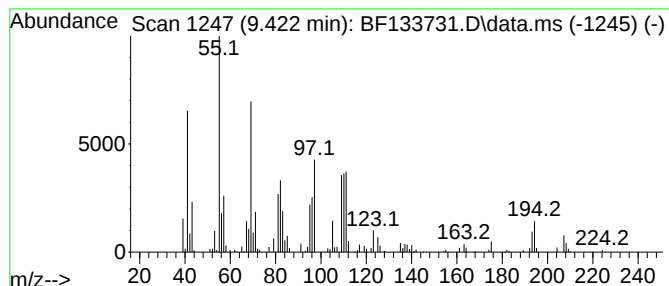
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown9.422 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.422	3.02 ng	110680	Acenaphthene-d10	9.880

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-(cyclohexylmethyl)...	194	C14H26	054823-95-9	35
2			Cyclohexane, 1-(cyclohexylmethyl)...	194	C14H26	054823-98-2	35
3			Oxirane, tetradecyl-	240	C16H32O	007320-37-8	27
4			(2,4-Dimethylcyclohexyl)methanol...	142	C9H18O	1000499-02-8	27
5			4-Isopropyl-5,10-dimethyl-decali...	236	C15H24O2	291278-94-9	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061323\
Data File : BF133731.D
Acq On : 13 Jun 2023 15:16
Operator : CG\JU
Sample : 03094-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

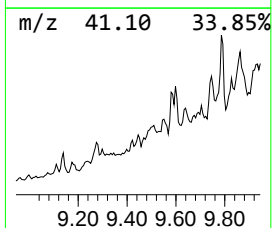
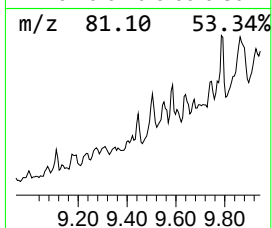
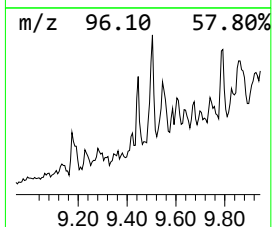
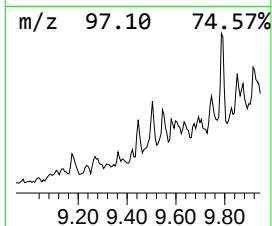
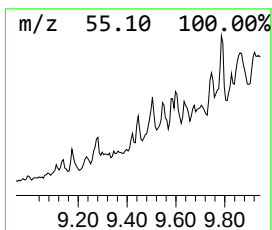
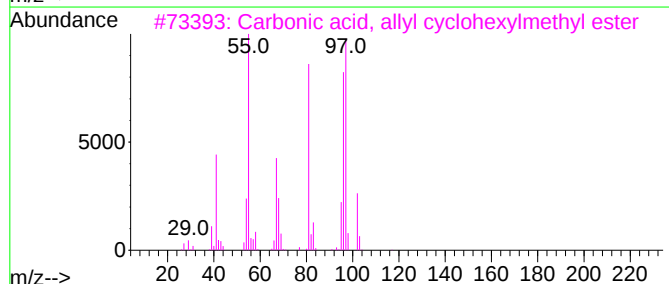
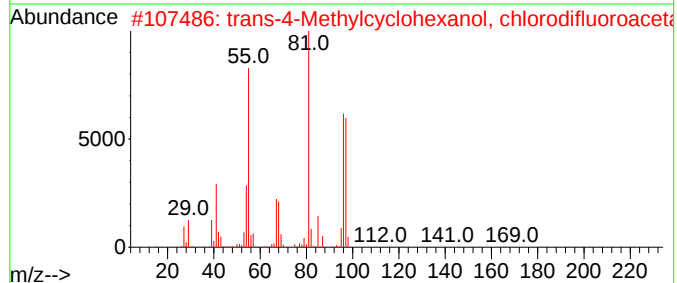
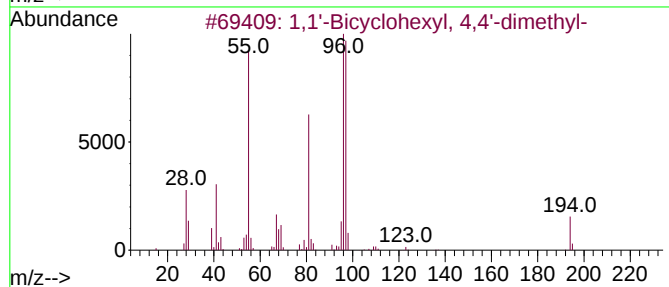
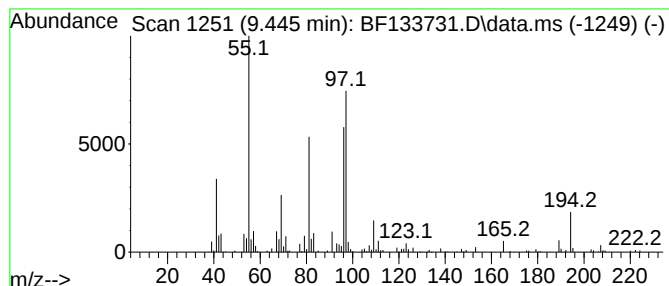
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1,1'-Bicyclohexyl, 4,4'-dim... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.445	4.16 ng	152374	Acenaphthene-d10	9.880	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Bicyclohexyl, 4,4'-dimethyl-	194	C14H26	054823-99-3	76
2	trans-4-Methylcyclohexanol, chlo...	226	C9H13ClF2O2	1000376-26-1	59
3	Carbonic acid, allyl cyclohexylm...	198	C11H18O3	1000357-80-7	59
4	Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	001678-82-6	59
5	Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	006069-98-3	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061323\
Data File : BF133731.D
Acq On : 13 Jun 2023 15:16
Operator : CG\JU
Sample : 03094-01
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ALS Vial : 12 Sample Multiplier: 1

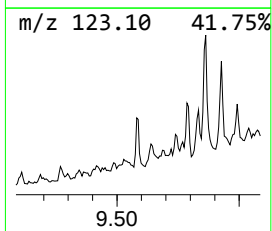
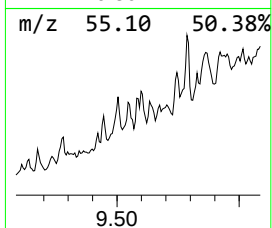
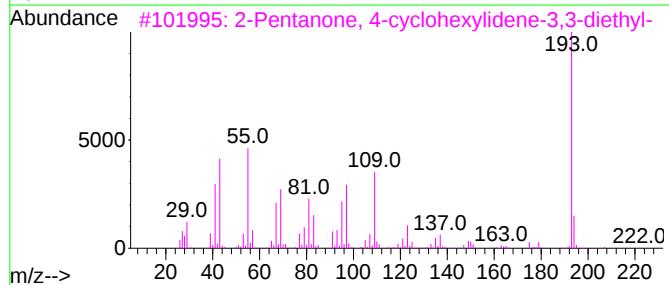
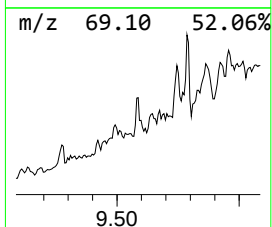
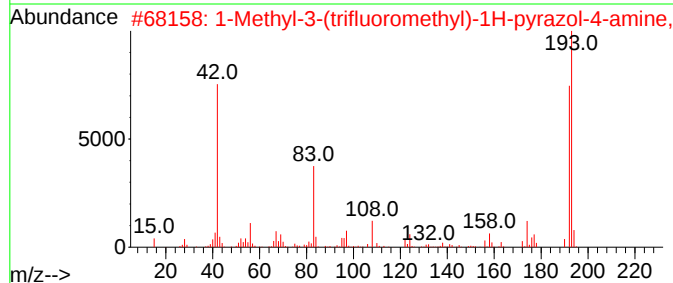
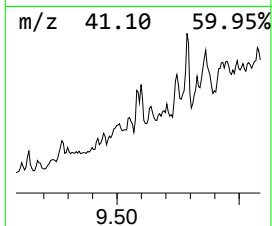
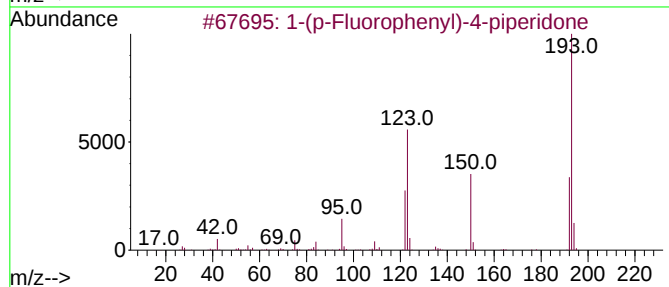
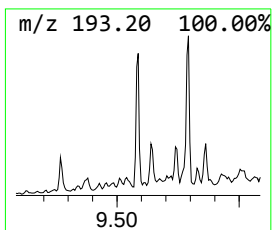
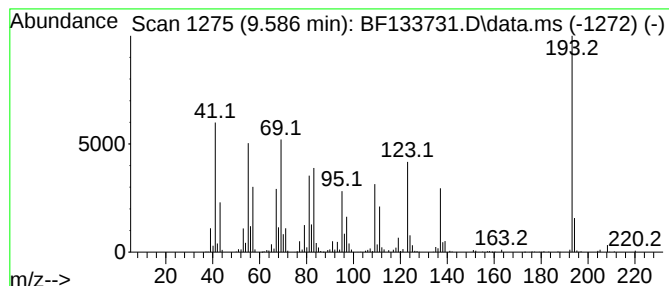
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ClientSampleId :
WC-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown9.586 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.586	15.58 ng	570362	Acenaphthene-d10	9.880	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	43
2	1-Methyl-3-(trifluoromethyl)-1H...	193	C7H10F3N3	1000511-73-1	38
3	2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	37
4	Butan-2-one, 1-[3-(3,3-dimethyl-...	278	C16H26N2O2	1000303-34-2	27
5	2,6-Dodecadiene, 2,6-dimethyl-	194	C14H26	1000164-67-6	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061323\
Data File : BF133731.D
Acq On : 13 Jun 2023 15:16
Operator : CG\JU
Sample : 03094-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

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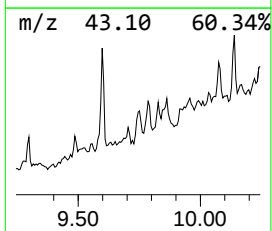
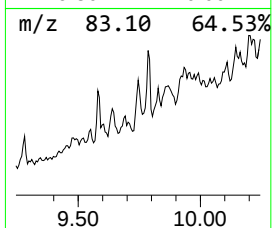
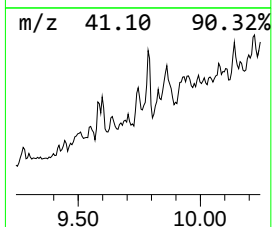
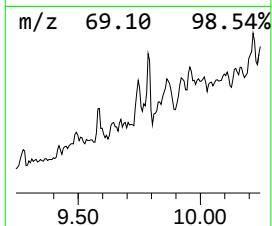
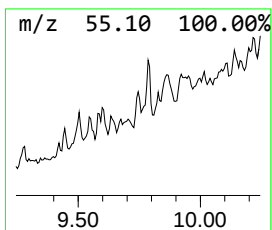
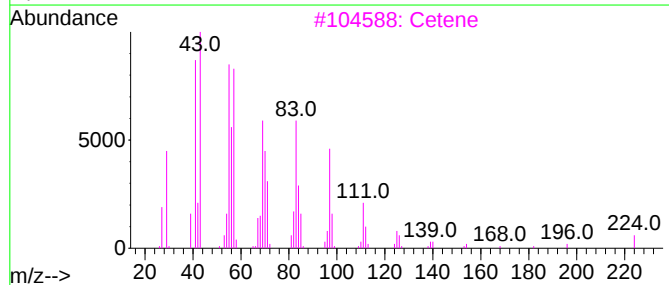
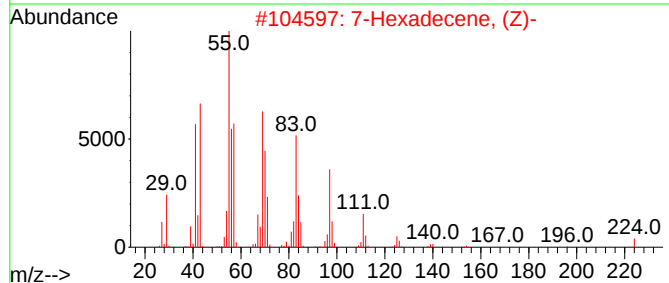
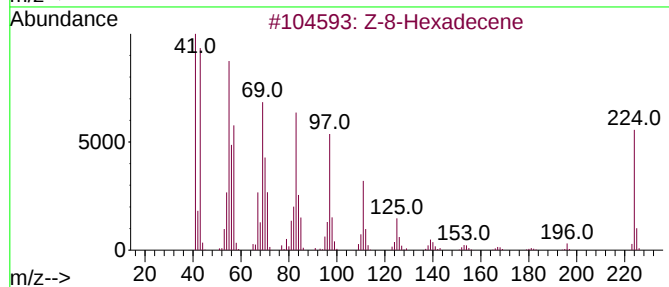
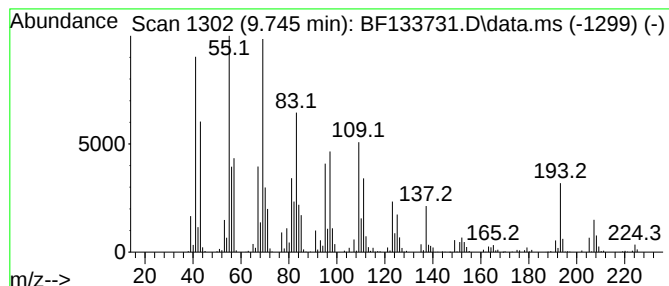
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Z-8-Hexadecene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.745	22.73 ng	831905	Acenaphthene-d10	9.880

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Z-8-Hexadecene	224	C16H32	1000130-87-5	51
2		7-Hexadecene, (Z)-	224	C16H32	035507-09-6	47
3		Cetene	224	C16H32	000629-73-2	46
4		Crotyl methacrylate	140	C8H12O2	007376-45-6	42
5		6-Tridecene, 7-methyl-	196	C14H28	024949-42-6	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061323\
Data File : BF133731.D
Acq On : 13 Jun 2023 15:16
Operator : CG\JU
Sample : 03094-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-1

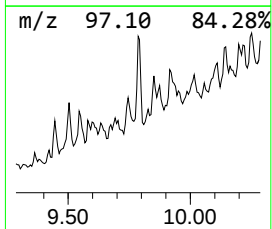
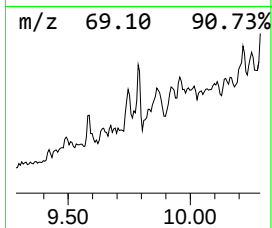
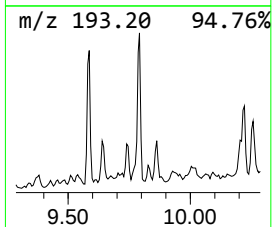
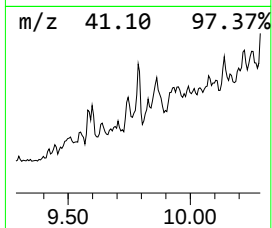
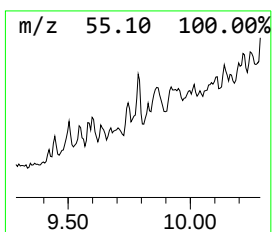
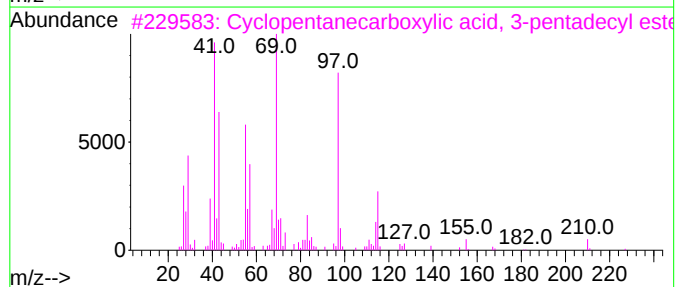
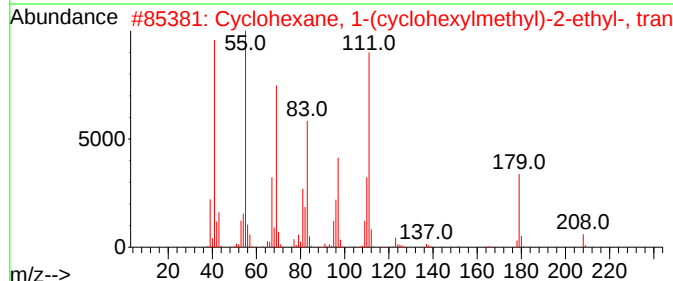
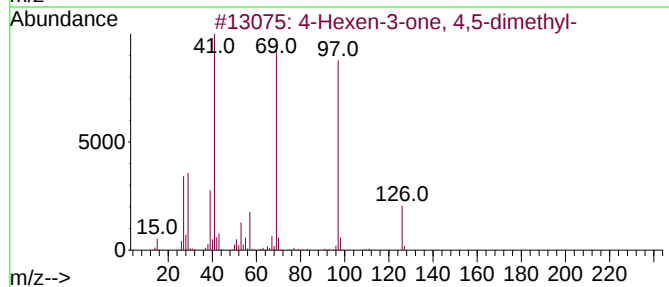
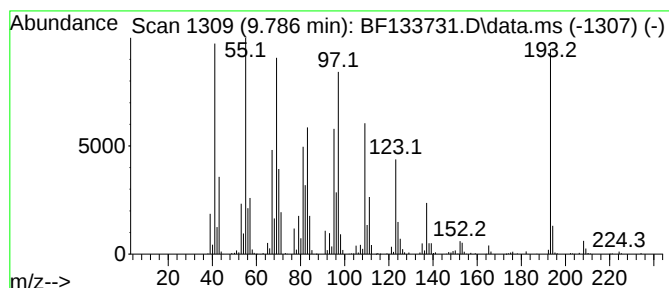
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown9.786 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.786	24.97 ng	914091	Acenaphthene-d10	9.880

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		4-Hexen-3-one, 4,5-dimethyl-	126	C8H14O	017325-90-5	25
2		Cyclohexane, 1-(cyclohexylmethyl)-	208	C15H28	054934-92-8	25
3		Cyclopentanecarboxylic acid, 3-pentadecyl ester	324	C21H40O2	1000280-59-1	22
4		Cyclohexane, 1-(cyclohexylmethyl)-	194	C14H26	054823-96-0	18
5		3-(But-3-enyl)-cyclohexanone	152	C10H16O	003636-03-1	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061323\
Data File : BF133731.D
Acq On : 13 Jun 2023 15:16
Operator : CG\JU
Sample : 03094-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

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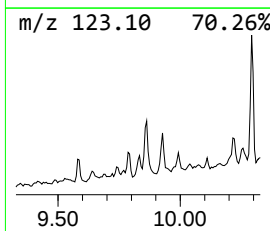
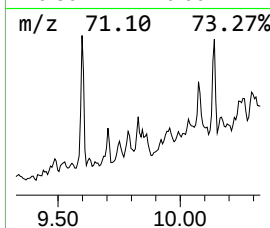
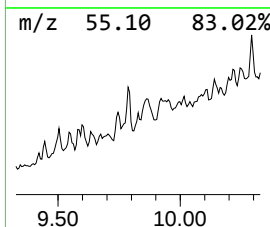
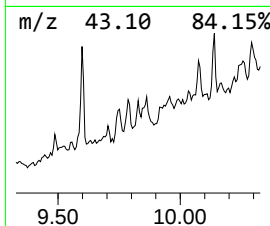
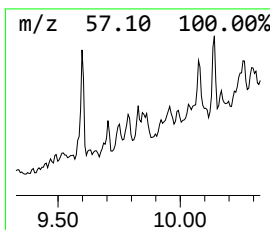
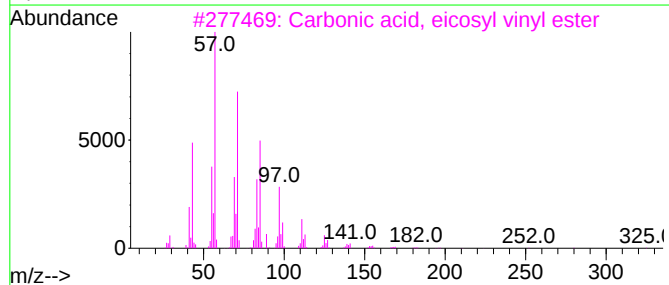
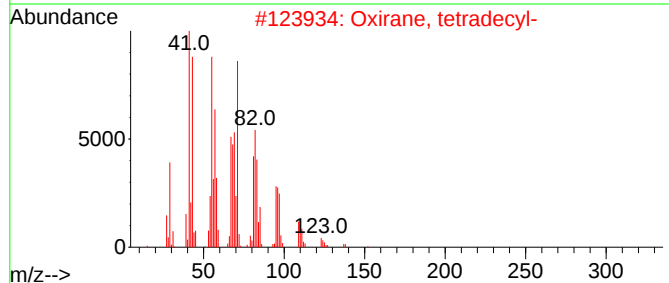
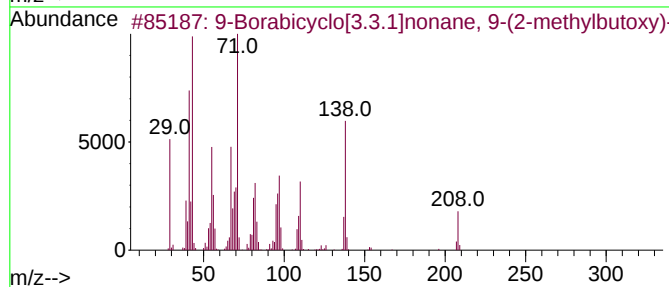
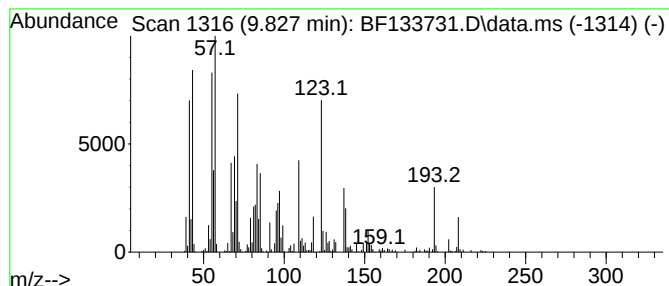
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 9-Borabicyclo[3.3.1]nonane,... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.827	8.33 ng	304915	Acenaphthene-d10	9.880

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Borabicyclo[3.3.1]nonane, 9-(2...	208	C13H25BO	094537-53-8	70
2		Oxirane, tetradecyl-	240	C16H32O	007320-37-8	38
3		Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	25
4		1-(3,5-Dimethyl-1H-pyrazol-4-yl)...	138	C7H10N2O	001123-48-4	25
5		Benzene, 1,4-dimethoxy-	138	C8H10O2	000150-78-7	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061323\
Data File : BF133731.D
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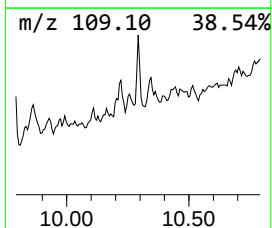
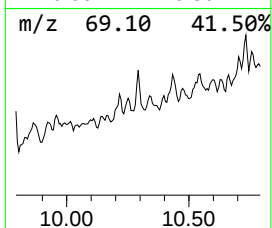
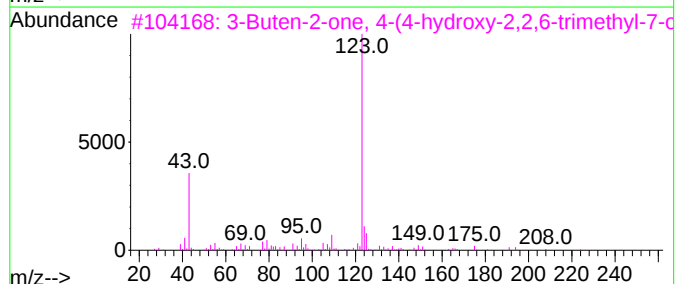
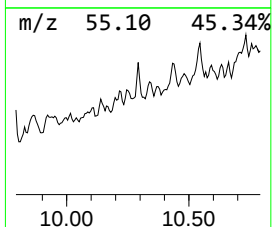
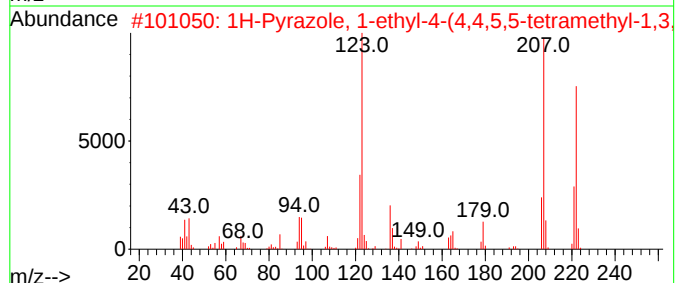
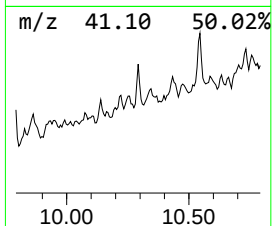
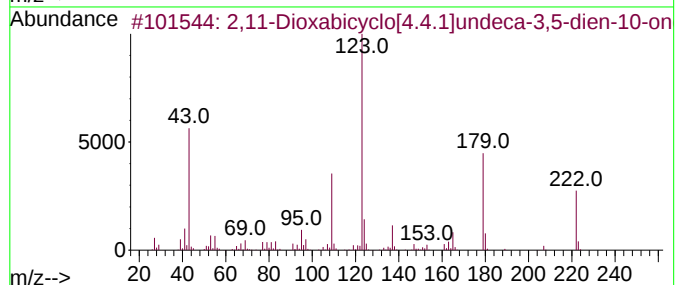
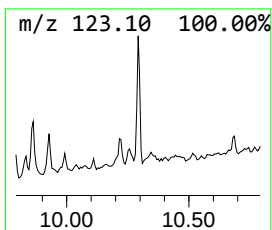
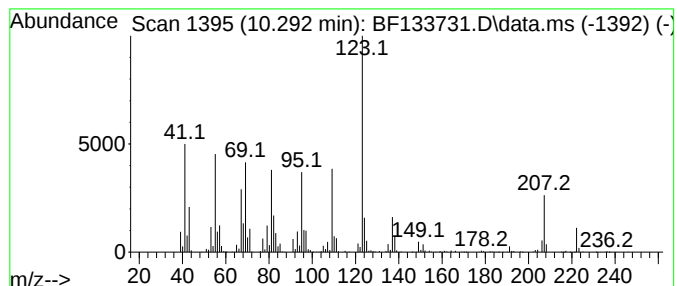
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 2,11-Dioxabicyclo[4.4.1]und... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.292	33.08 ng	1210880	Acenaphthene-d10	9.880	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,11-Dioxabicyclo[4.4.1]undeca-3...	222	C13H18O3	070412-52-1	59
2	1H-Pyrazole, 1-ethyl-4-(4,4,5,5-...	222	C11H19BN2O2	1010319-69-6	50
3	3-Buten-2-one, 4-(4-hydroxy-2,2,...	224	C13H20O3	038274-01-0	47
4	3-Fluorobenzoic acid, 3-methylbu...	208	C12H13FO2	154559-38-3	46
5	(Phenylthio)acetic acid, cyclobu...	222	C12H14O2S	1010299-95-5	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061323\
Data File : BF133731.D
Acq On : 13 Jun 2023 15:16
Operator : CG\JU
Sample : 03094-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-1

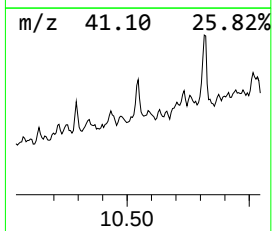
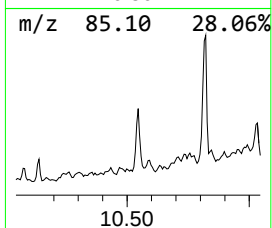
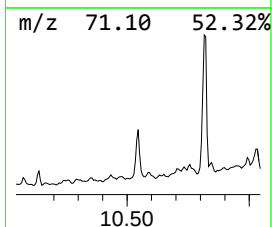
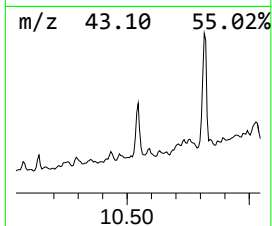
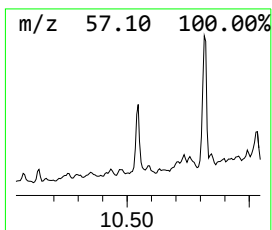
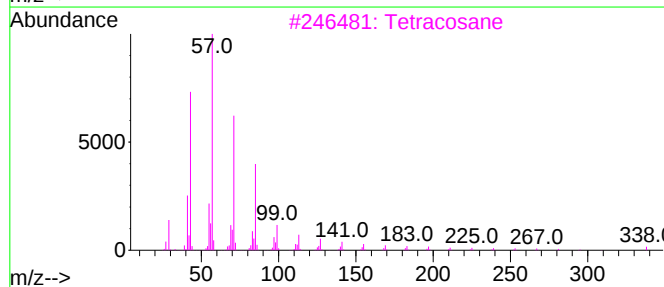
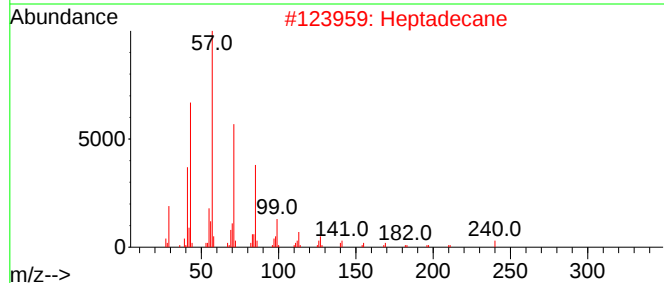
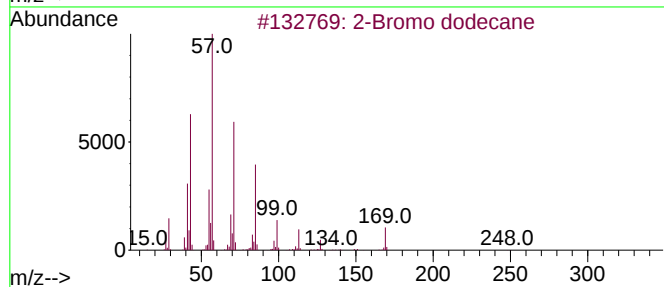
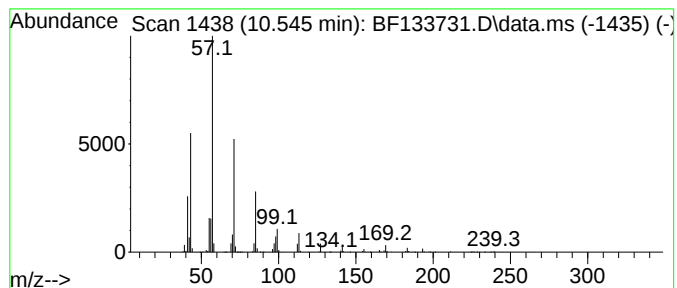
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 2-Bromo dodecane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.545	33.48 ng	1225400	Acenaphthene-d10	9.880

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Bromo dodecane	248	C12H25Br	013187-99-0	90
2		Heptadecane	240	C17H36	000629-78-7	90
3		Tetracosane	338	C24H50	000646-31-1	86
4		Octacosane	394	C28H58	000630-02-4	86
5		Heptacosane	380	C27H56	000593-49-7	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061323\
Data File : BF133731.D
Acq On : 13 Jun 2023 15:16
Operator : CG\JU
Sample : 03094-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-1

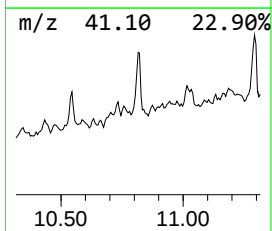
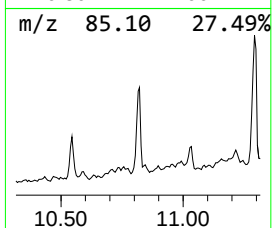
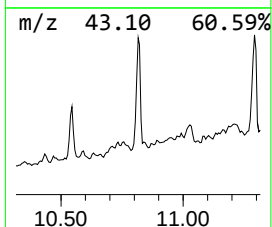
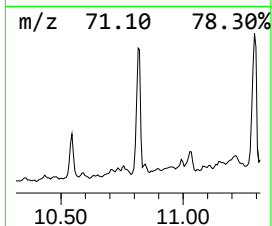
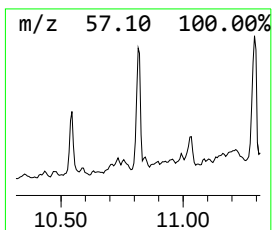
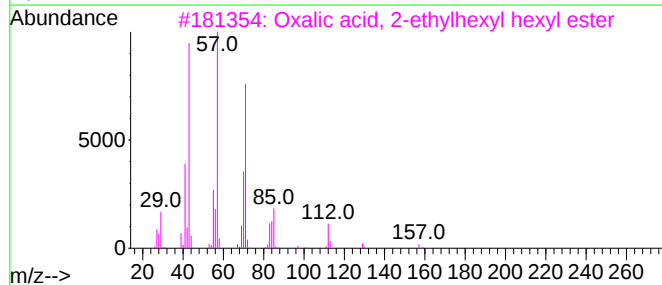
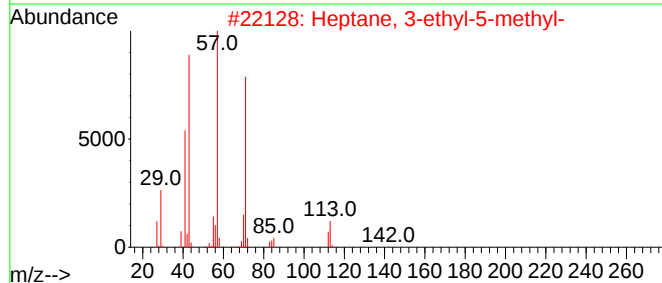
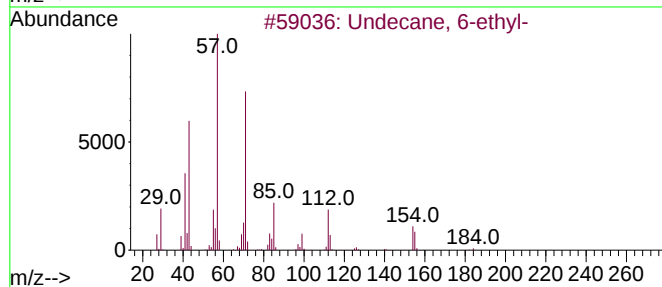
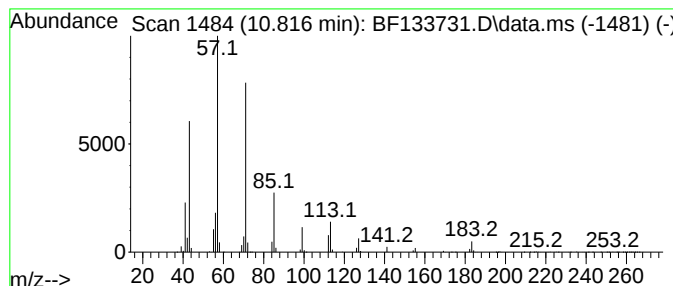
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Undecane, 6-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.816	18.54 ng	2935790	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 6-ethyl-	184	C13H28	017312-60-6	80
2			Heptane, 3-ethyl-5-methyl-	142	C10H22	052896-90-9	64
3			Oxalic acid, 2-ethylhexyl hexyl ...	286	C16H30O4	1000309-38-9	53
4			3-Hexanone, 4-ethyl-	128	C8H16O	006137-12-8	53
5			Hexane, 2,4,4-trimethyl-	128	C9H20	016747-30-1	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061323\
Data File : BF133731.D
Acq On : 13 Jun 2023 15:16
Operator : CG\JU
Sample : 03094-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

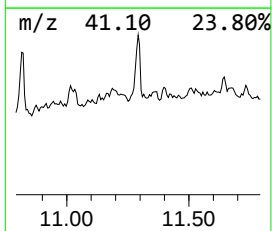
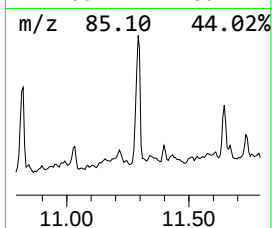
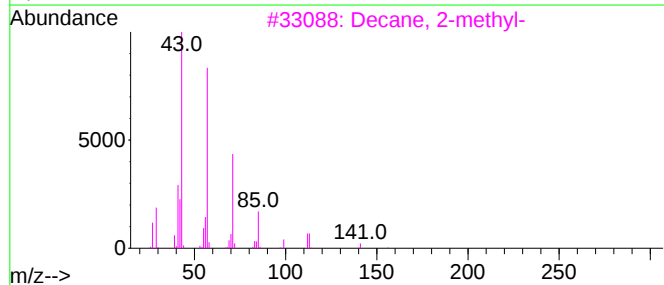
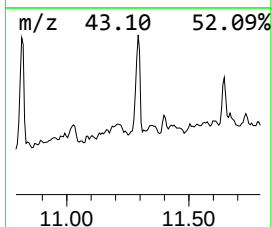
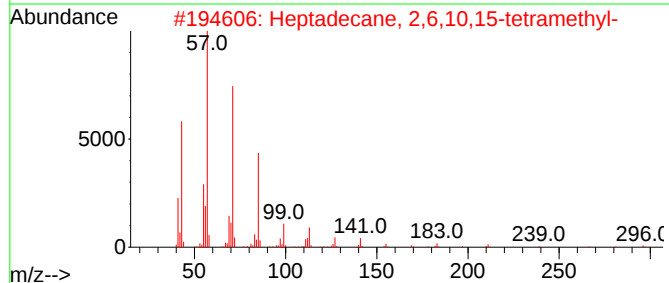
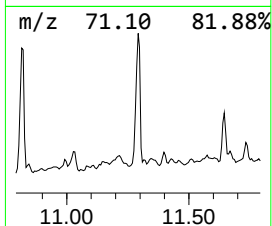
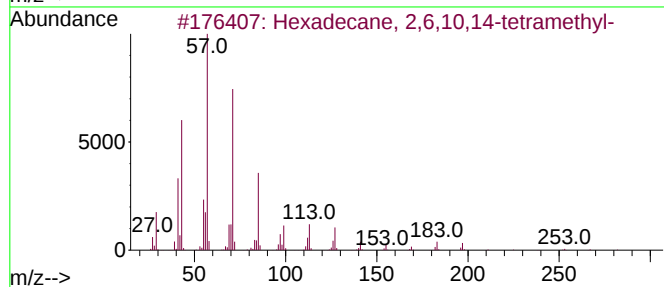
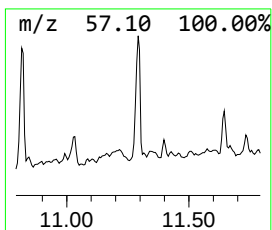
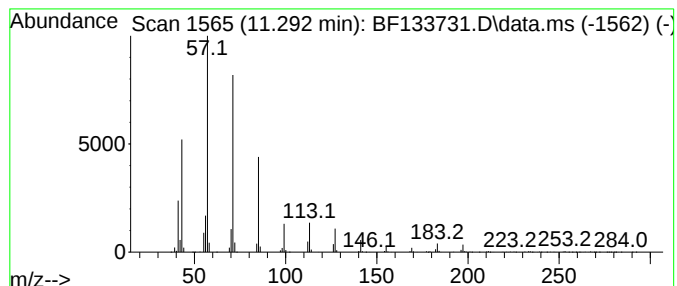
Instrument :
BNA_F
ClientSampleId :
WC-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Hexadecane, 2,6,10,14-tetra... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.292	20.00 ng	3167820	Phenanthrene-d10	11.386	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	91
2	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
3	Decane, 2-methyl-	156	C11H24	006975-98-0	83
4	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	81
5	Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061323\
Data File : BF133731.D
Acq On : 13 Jun 2023 15:16
Operator : CG\JU
Sample : 03094-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060923.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	2.210	3.8	ng	96582	1	6.834	508579	20.0
2-Heptanone, 6-...	6.269	5.1	ng	130055	1	6.834	508579	20.0
unknown8.998	8.998	3.9	ng	98265	2	8.116	503378	20.0
Nonane, 3-methy...	9.139	5.1	ng	186027	3	9.880	732067	20.0
unknown9.275	9.275	10.2	ng	374174	3	9.880	732067	20.0
unknown9.422	9.422	3.0	ng	110680	3	9.880	732067	20.0
1,1'-Bicyclohex...	9.445	4.2	ng	152374	3	9.880	732067	20.0
unknown9.586	9.586	15.6	ng	570362	3	9.880	732067	20.0
unknown9.639	9.639	11.2	ng	410215	3	9.880	732067	20.0
Z-8-Hexadecene	9.745	22.7	ng	831905	3	9.880	732067	20.0
unknown9.786	9.786	25.0	ng	914091	3	9.880	732067	20.0
9-Borabicyclo[3...	9.827	8.3	ng	304915	3	9.880	732067	20.0
2,11-Dioxabicyc...	10.292	33.1	ng	1210880	3	9.880	732067	20.0
2-Bromo dodecane	10.545	33.5	ng	1225400	3	9.880	732067	20.0
Undecane, 6-ethyl-	10.816	18.5	ng	2935790	4	11.386	3167820	20.0
Hexadecane, 2,6...	11.292	20.0	ng	3167820	4	11.386	3167820	20.0