

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092823\
 Data File : BF135502.D
 Acq On : 28 Sep 2023 19:39
 Operator : CG\JU
 Sample : 04606-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ETGI-327

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-BF091123.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF135502.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.381	556	560	574	rBV	2034516	2297777	97.35%	8.264%
2	6.398	727	733	745	rBV	2168690	2360433	100.00%	8.490%
3	6.746	789	792	800	rBV	955088	789091	33.43%	2.838%
4	6.810	800	803	809	rBV	255995	238759	10.12%	0.859%
5	7.310	885	888	891	rBV	1598984	1288459	54.59%	4.634%
6	7.434	903	909	911	rBV	199555	251134	10.64%	0.903%
7	7.769	963	966	973	rBV5	89581	104663	4.43%	0.376%
8	8.022	1006	1009	1012	rBV	1012343	840010	35.59%	3.021%
9	8.051	1012	1014	1018	rVB2	367019	302657	12.82%	1.089%
10	8.310	1054	1058	1061	rBV5	113741	139693	5.92%	0.502%
11	8.445	1077	1081	1083	rBV	239538	261111	11.06%	0.939%
12	8.557	1097	1100	1102	rBV4	176161	203749	8.63%	0.733%
13	8.710	1123	1126	1129	rVB2	190151	220455	9.34%	0.793%
14	8.739	1129	1131	1133	rBV	311181	227045	9.62%	0.817%
15	8.763	1133	1135	1138	rBV3	171793	156252	6.62%	0.562%
16	8.792	1138	1140	1143	rBV2	125825	168126	7.12%	0.605%
17	8.904	1156	1159	1161	rBV3	312936	315612	13.37%	1.135%
18	9.010	1175	1177	1180	rBV3	359478	384865	16.30%	1.384%
19	9.051	1180	1184	1187	rVB	838724	800207	33.90%	2.878%
20	9.104	1187	1193	1196	rBV	1967306	2069052	87.66%	7.442%
21	9.139	1196	1199	1201	rBV2	268421	277883	11.77%	0.999%
22	9.187	1201	1207	1208	rBV4	881438	1509563	63.95%	5.429%
23	9.487	1256	1258	1260	rBV2	1233312	1031581	43.70%	3.710%
24	9.510	1260	1262	1265	rVB	1368339	1188054	50.33%	4.273%
25	9.539	1265	1267	1271	rBV2	748378	710215	30.09%	2.554%
26	9.604	1275	1278	1279	rBV2	554535	565885	23.97%	2.035%
27	9.651	1283	1286	1291	rBV4	1039050	2071337	87.75%	7.450%
28	9.692	1291	1293	1296	rVB	1912170	1765044	74.78%	6.348%
29	9.734	1297	1300	1302	rBV4	781283	997254	42.25%	3.587%
30	9.763	1302	1305	1307	rBV3	1191501	1424453	60.35%	5.123%
31	14.980	2189	2192	2195	rBV	603912	747798	31.68%	2.690%
32	15.280	2240	2243	2249	rVB	462923	570765	24.18%	2.053%
33	15.339	2250	2253	2259	rVB	511869	615459	26.07%	2.214%
34	15.404	2260	2264	2267	rBV	441590	535355	22.68%	1.925%
35	16.868	2510	2513	2521	rVB	202096	374337	15.86%	1.346%

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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-BF091123.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

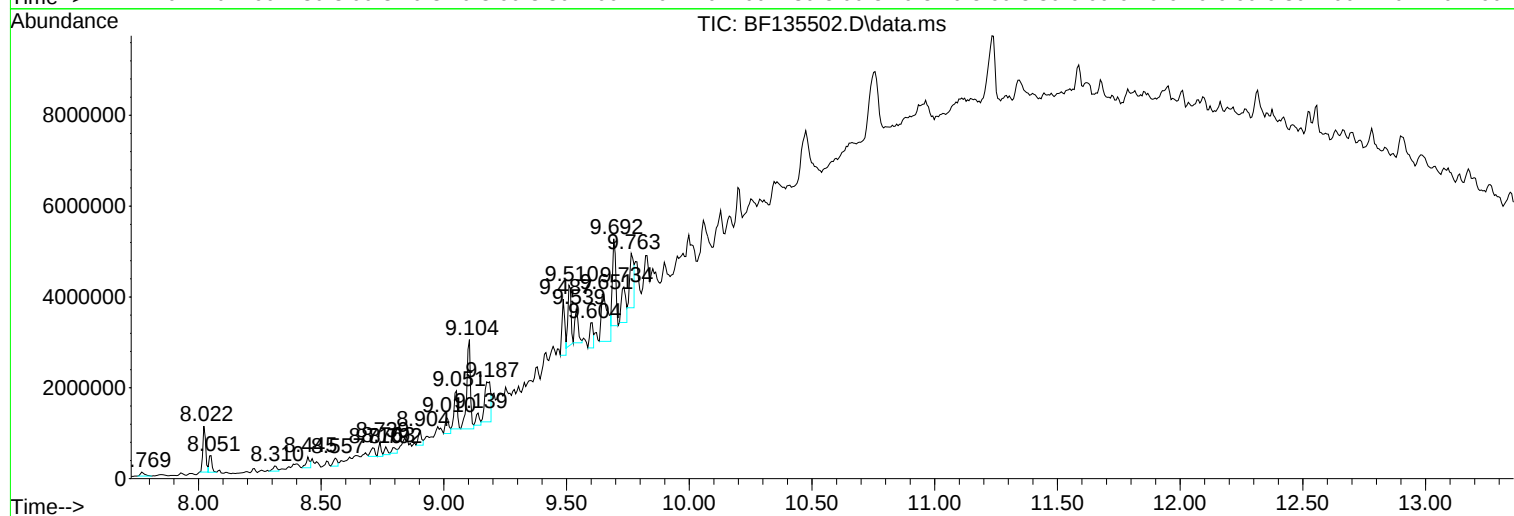
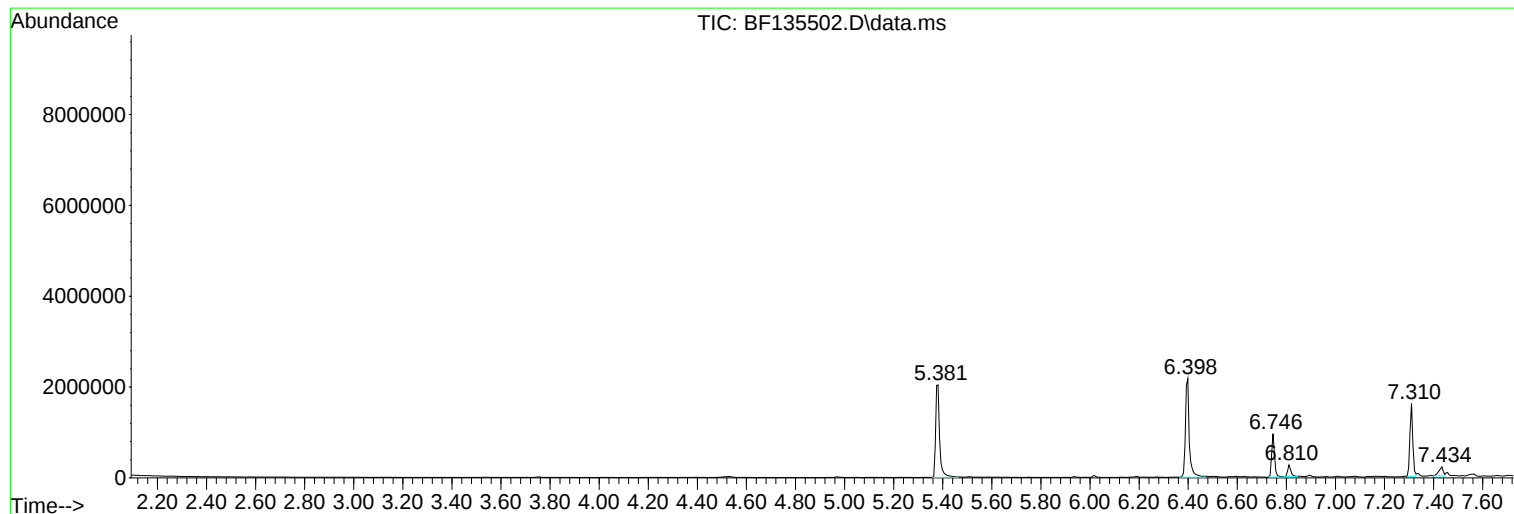
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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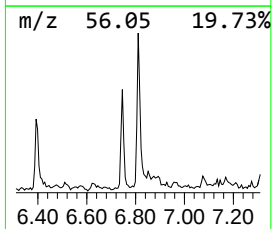
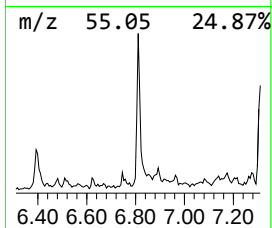
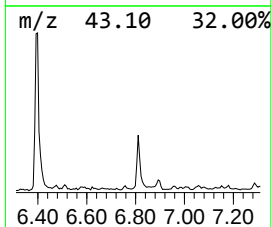
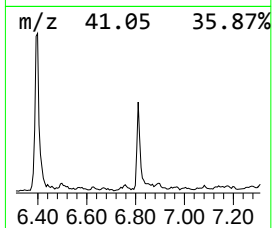
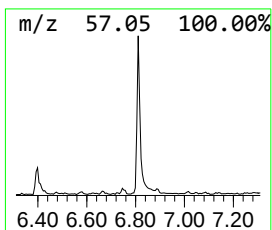
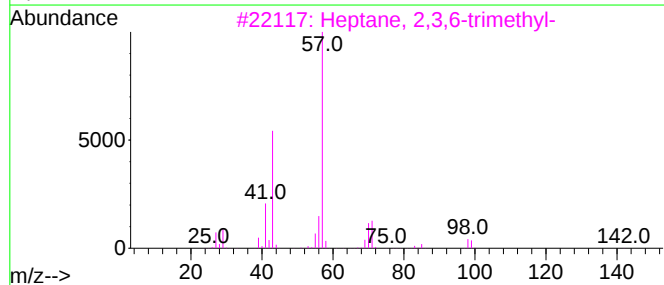
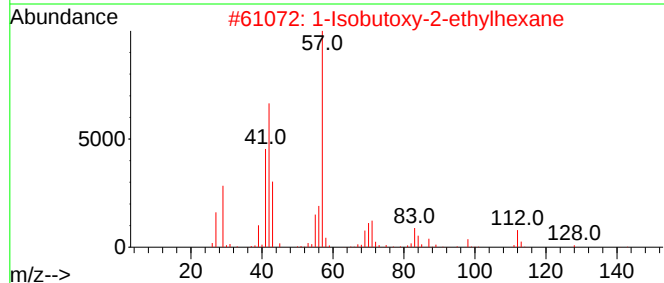
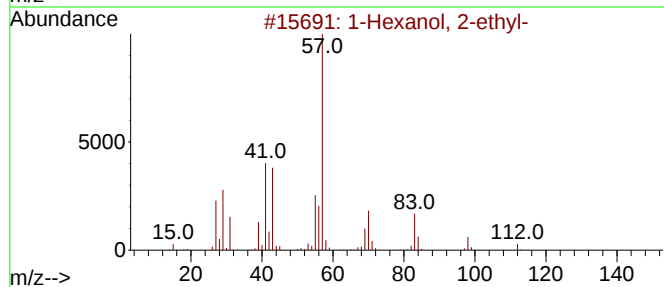
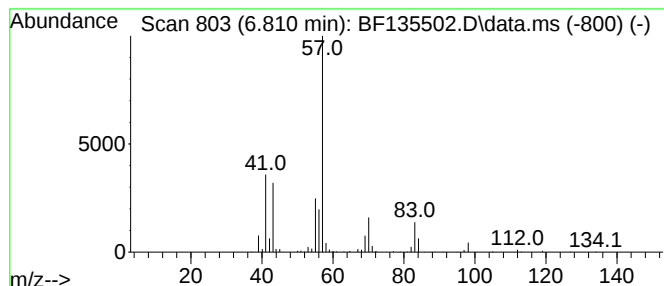
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 1-Hexanol, 2-ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.810	6.05 ng	238759	1,4-Dichlorobenzene-d4	6.745

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2			1-Isobutoxy-2-ethylhexane	186	C12H26O	1000139-90-3	64
3			Heptane, 2,3,6-trimethyl-	142	C10H22	004032-93-3	64
4			1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	56
5			Heptyl isobutyl carbonate	216	C12H24O3	959068-08-7	53



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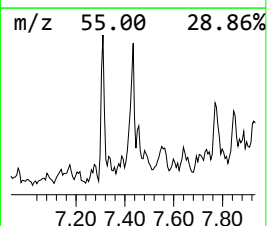
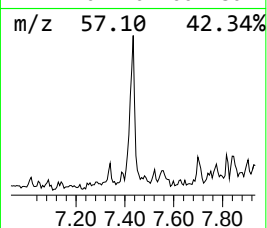
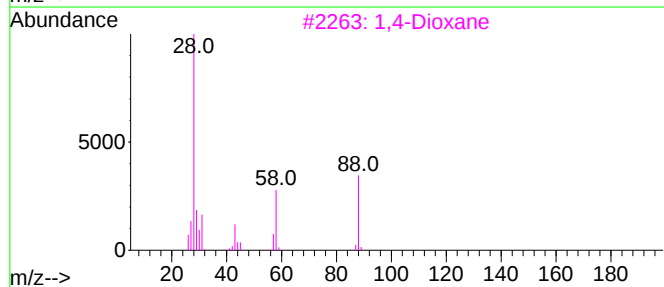
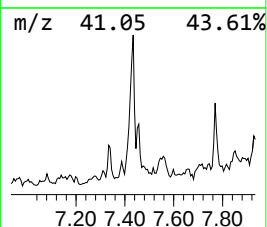
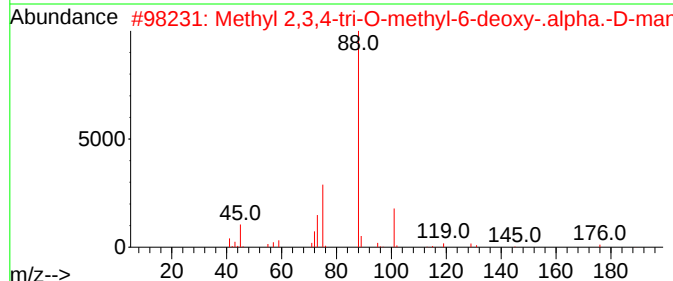
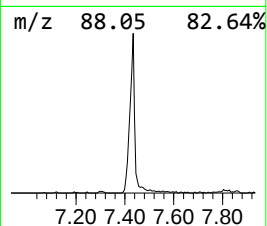
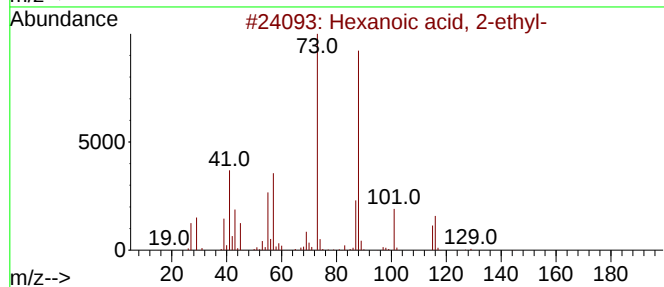
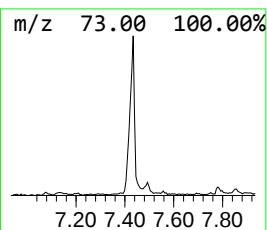
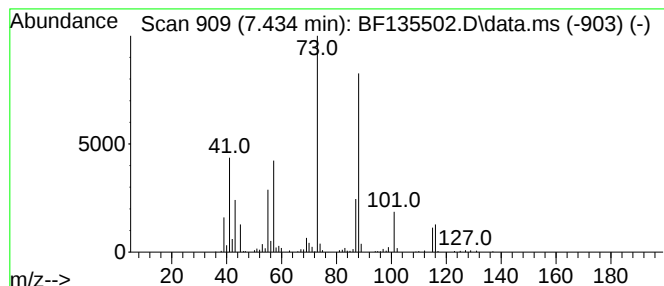
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Hexanoic acid, 2-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.434	5.98 ng	251134	Naphthalene-d8	8.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2			Methyl 2,3,4-tri-O-methyl-6-deox...	220	C10H20O5	072983-16-5	50
3			1,4-Dioxane	88	C4H8O2	000123-91-1	47
4			Acetic acid, diethyl-	116	C6H12O2	000088-09-5	42
5			Heptanoic acid, 2-ethyl-	158	C9H18O2	003274-29-1	42



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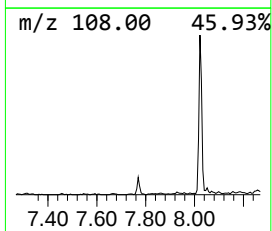
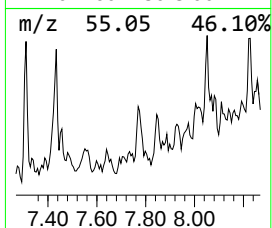
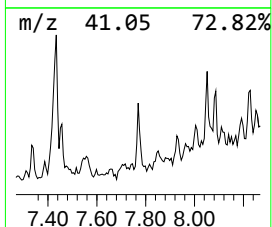
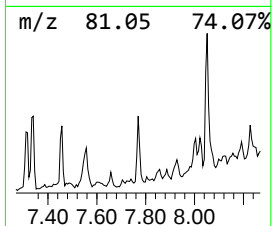
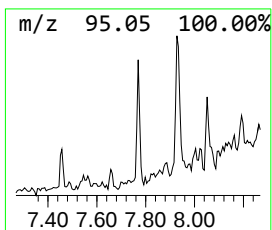
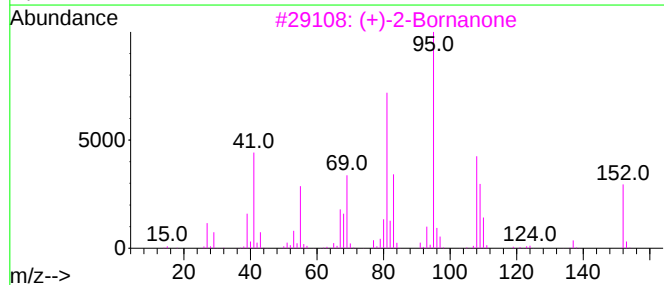
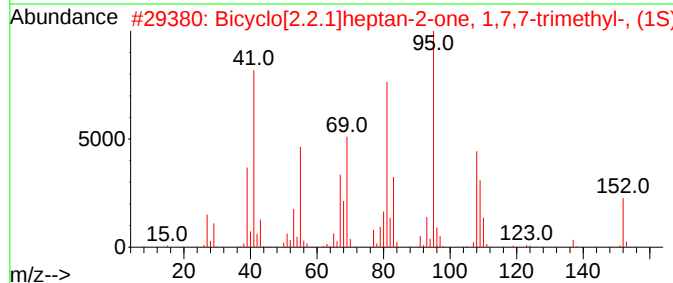
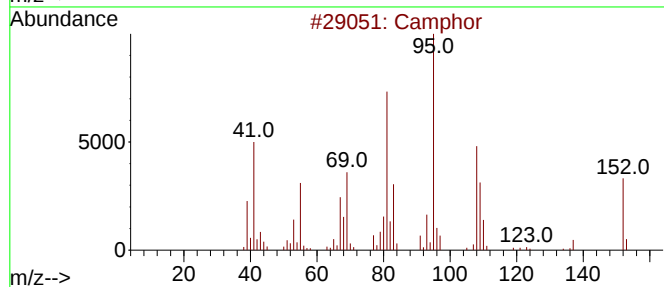
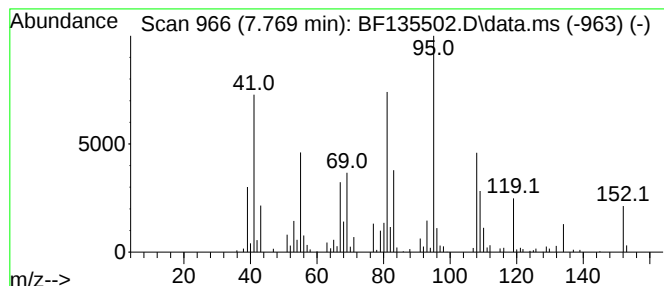
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Camphor Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.769	2.49 ng	104663	Naphthalene-d8	8.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Camphor	152	C10H16O	000076-22-2	93
2			Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	93
3			(+)-2-Bornanone	152	C10H16O	000464-49-3	93
4			2,4-Decadienal, (E,E)-	152	C10H16O	025152-84-5	38
5			Cyclohexanone, 5-methyl-2-(1-met...	152	C10H16O	000529-00-0	30



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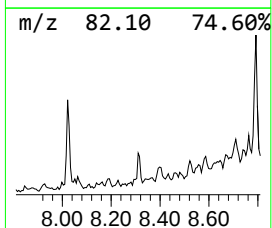
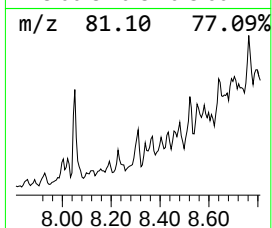
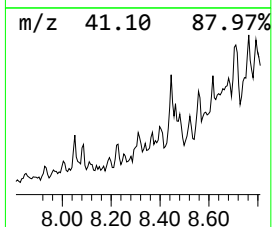
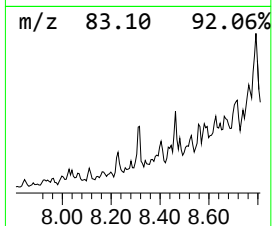
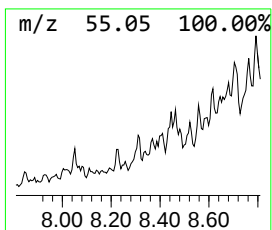
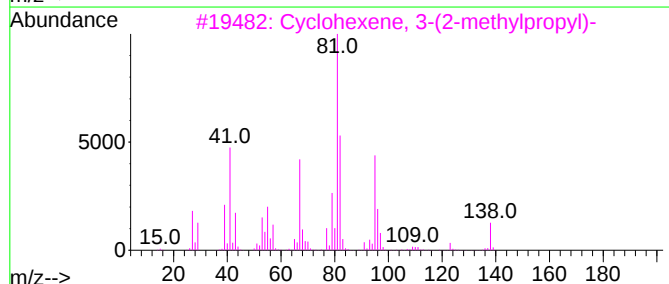
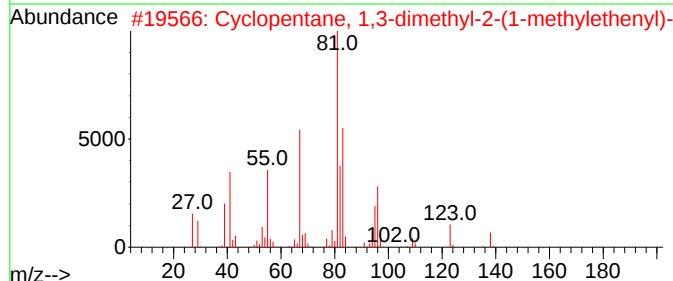
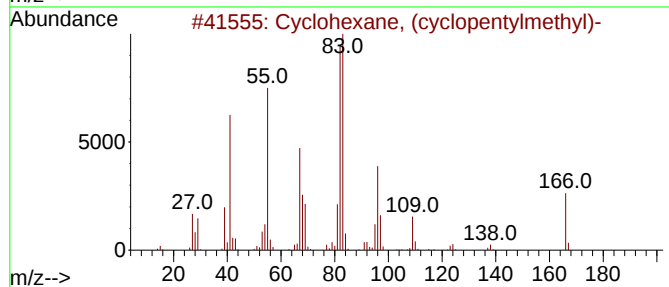
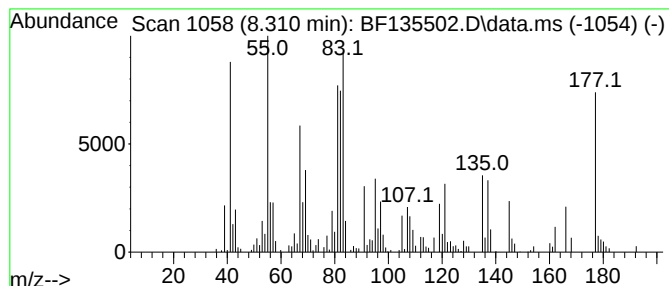
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown8.310 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.310	3.33 ng	139693	Naphthalene-d8	8.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (cyclopentylmethyl)-	166	C12H22	004431-89-4	38
2			Cyclopentane, 1,3-dimethyl-2-(1-...	138	C10H18	061142-31-2	30
3			Cyclohexene, 3-(2-methylpropyl)-	138	C10H18	004104-56-7	25
4			Cyclohexane, hexyl-	168	C12H24	004292-75-5	25
5			Cyclohexane, (4-methylpentyl)-	168	C12H24	061142-20-9	25



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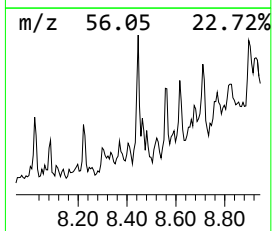
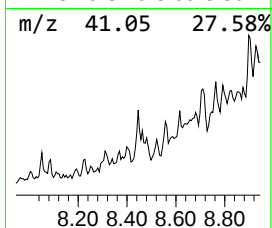
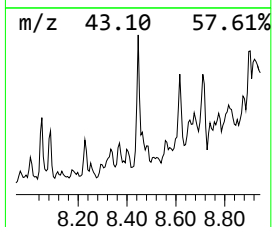
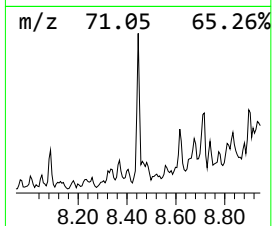
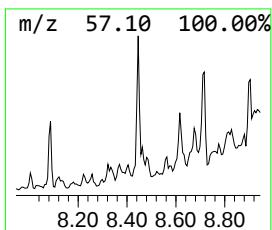
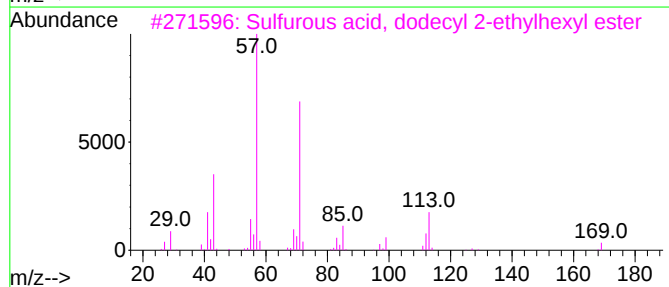
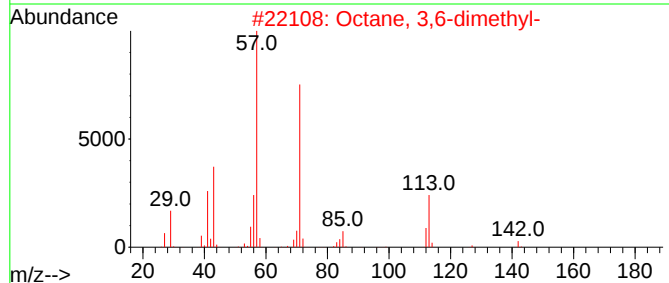
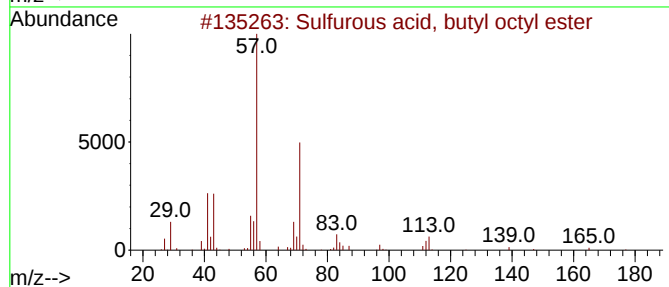
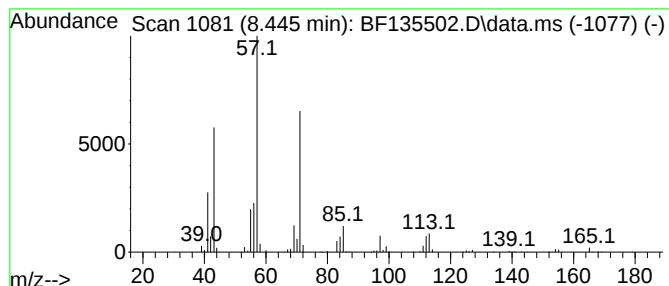
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Sulfurous acid, butyl octyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.445	6.22 ng	261111	Naphthalene-d8	8.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, butyl octyl ester	250	C12H26O3S	1000309-17-5	64
2			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	64
3			Sulfurous acid, dodecyl 2-ethylh...	362	C20H42O3S	1000309-19-5	64
4			Nonane, 3-methyl-	142	C10H22	005911-04-6	64
5			6-Methyl-2-heptanol, 2-methylpro...	200	C12H24O2	1000447-36-7	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092823\
Data File : BF135502.D
Acq On : 28 Sep 2023 19:39
Operator : CG\JU
Sample : 04606-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-327

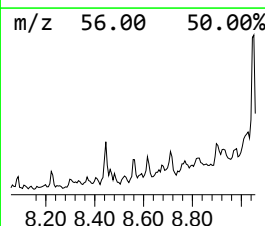
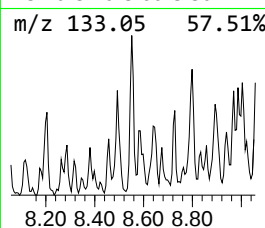
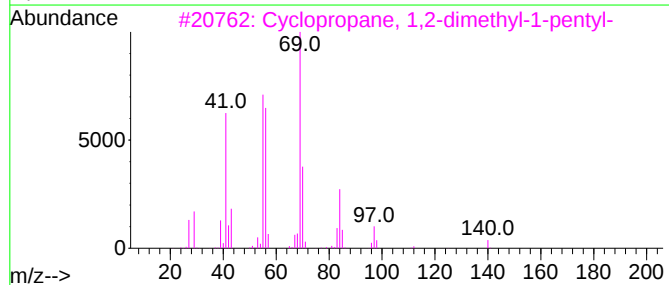
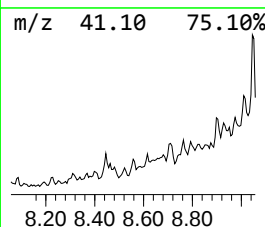
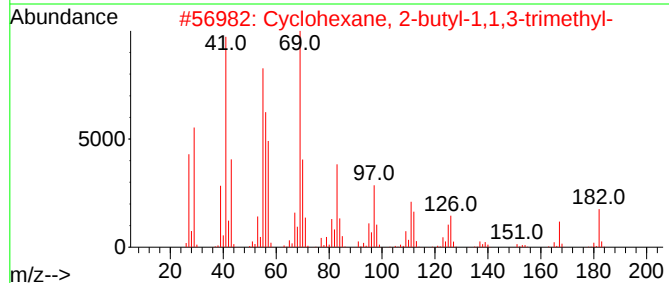
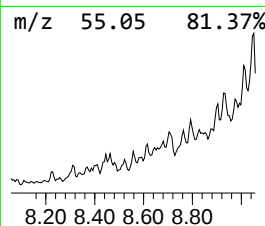
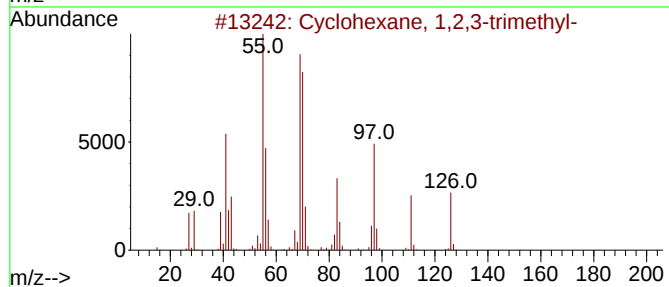
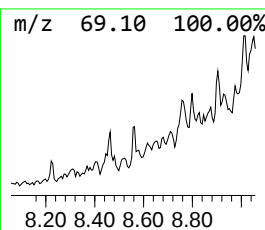
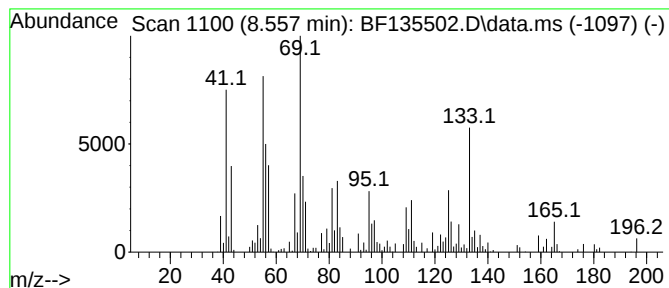
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown8.557 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.557	4.85 ng	203749	Naphthalene-d8	8.022

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 1,2,3-trimethyl-	126	C9H18	001678-97-3	43
2		Cyclohexane, 2-butyl-1,1,3-trime...	182	C13H26	054676-39-0	43
3		Cyclopropane, 1,2-dimethyl-1-pen...	140	C10H20	062238-04-4	41
4		Cyclohexane, 1,2,3-trimethyl-, (...)	126	C9H18	001678-81-5	41
5		trans-2-Methyl-3-octene	126	C9H18	052937-36-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092823\
Data File : BF135502.D
Acq On : 28 Sep 2023 19:39
Operator : CG\JU
Sample : 04606-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

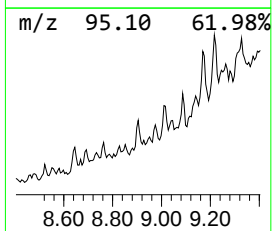
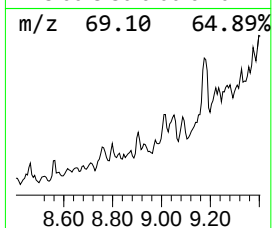
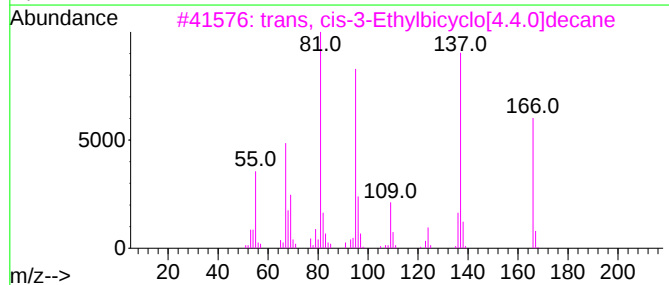
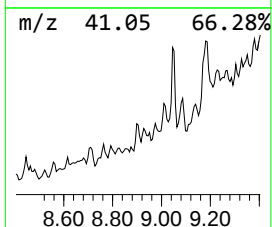
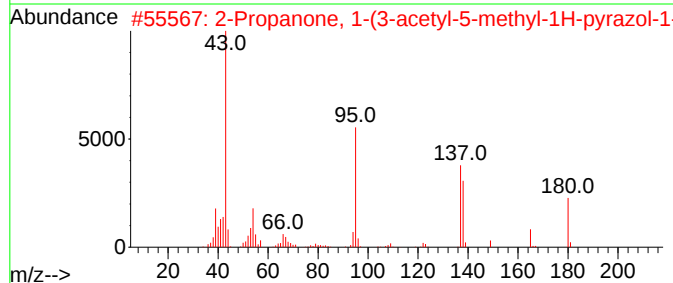
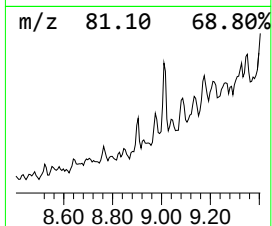
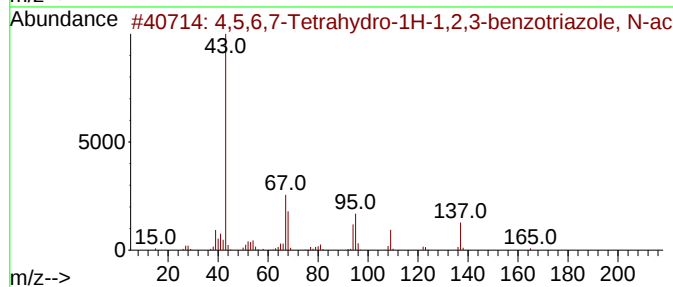
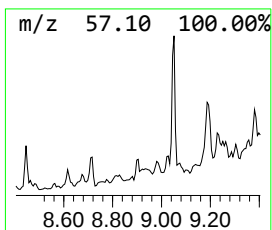
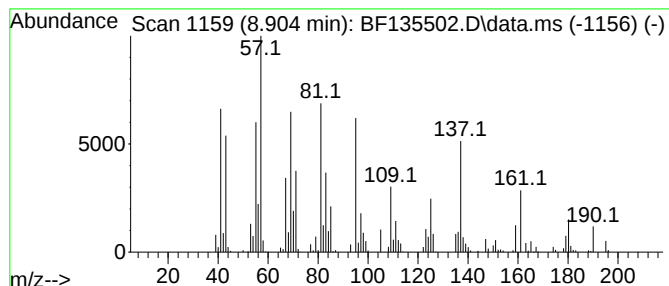
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown8.904 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.904	4.43 ng	315612	Acenaphthene-d10			9.786
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,5,6,7-Tetrahydro-1H-1,2,3-benz...	165	C8H11N3O	074581-99-0	43	
2	2-Propanone, 1-(3-acetyl-5-methy...	180	C9H12N2O2	1000362-64-1	38	
3	trans, cis-3-Ethylbicyclo[4.4.0]...	166	C12H22	066660-43-3	38	
4	Bicyclo[4.1.0]heptane, 2-methyl-...	180	C13H24	055937-92-3	35	
5	Naphthalene, 1,1'-ethylidenebis[...	302	C22H38	054934-70-2	22	



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092823\
Data File : BF135502.D
Acq On : 28 Sep 2023 19:39
Operator : CG\JU
Sample : 04606-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

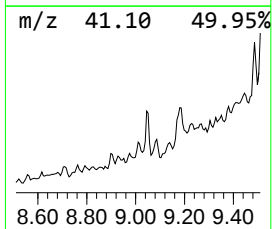
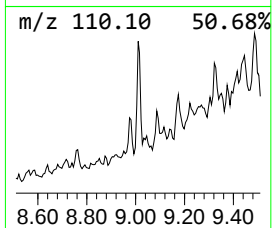
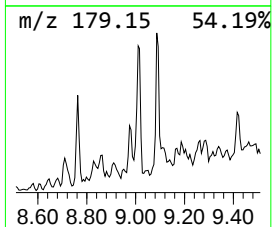
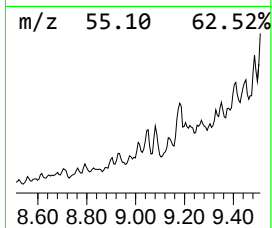
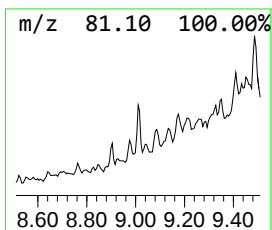
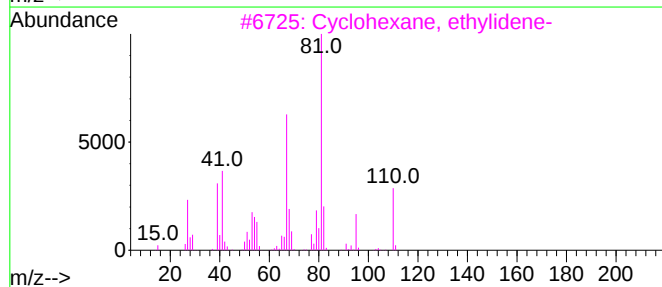
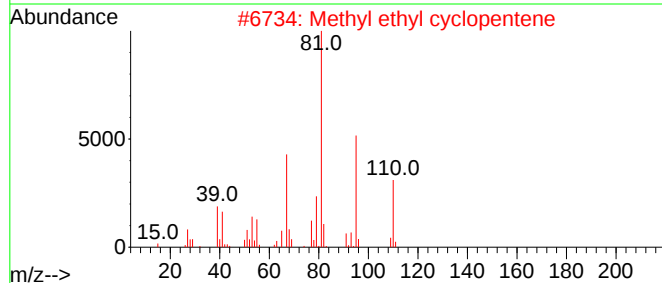
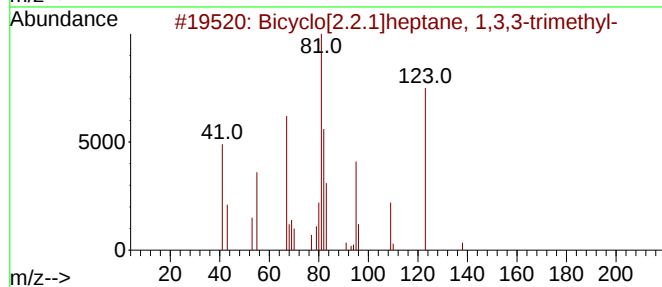
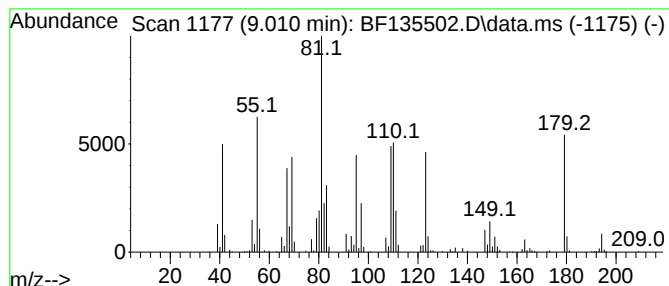
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Bicyclo[2.2.1]heptane, 1,3,... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD			R.T.
9.010	5.40 ng	384865	Acenaphthene-d10			9.786
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptane, 1,3,3-tri...		138	C10H18	006248-88-0	53
2	Methyl ethyl cyclopentene		110	C8H14	019780-56-4	46
3	Cyclohexane, ethylidene-		110	C8H14	001003-64-1	43
4	Cyclohexene, 1-ethyl-		110	C8H14	001453-24-3	38
5	2-Propenyl-3-vinyloxirane		110	C7H10O	070597-49-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092823\
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Operator : CG\JU
Sample : 04606-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

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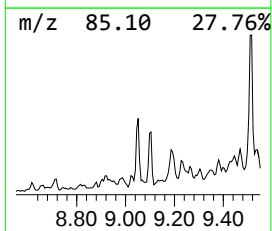
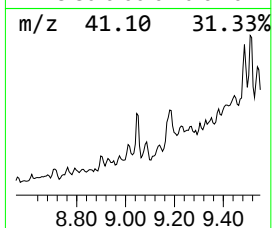
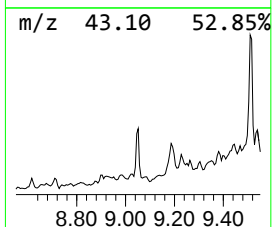
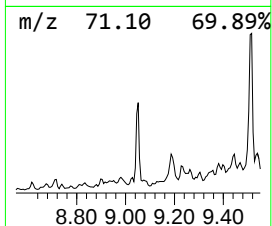
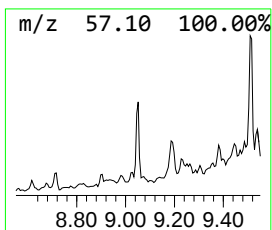
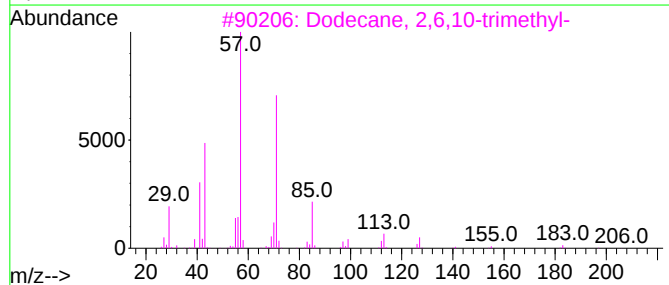
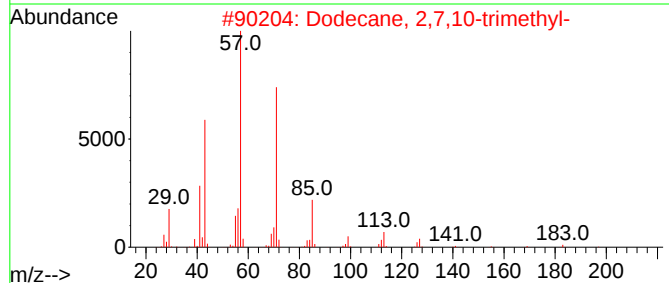
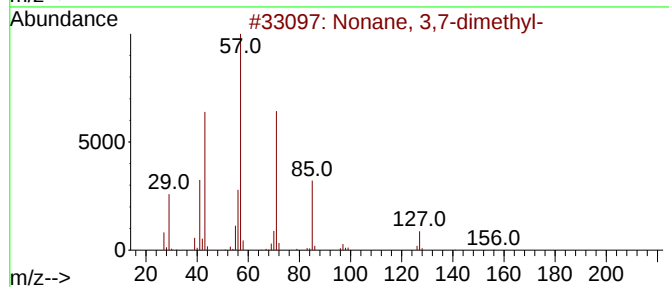
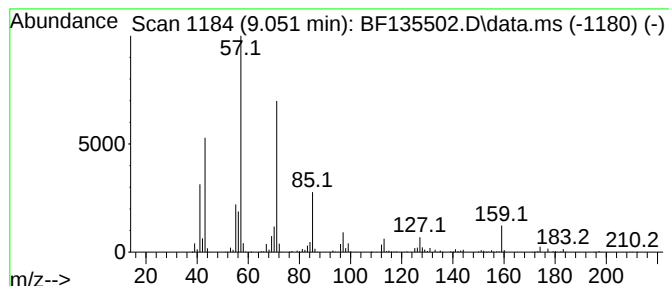
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Nonane, 3,7-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.051	11.24 ng	800207	Acenaphthene-d10	9.786

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	62
2			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	58
3			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	58
4			Hexadecane	226	C16H34	000544-76-3	53
5			Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092823\
Data File : BF135502.D
Acq On : 28 Sep 2023 19:39
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Sample : 04606-01
Misc :
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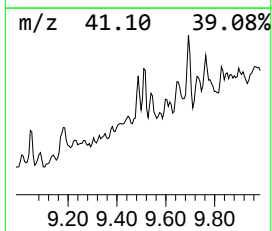
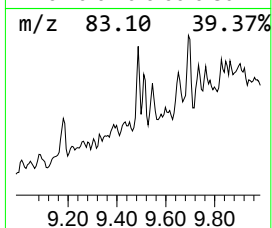
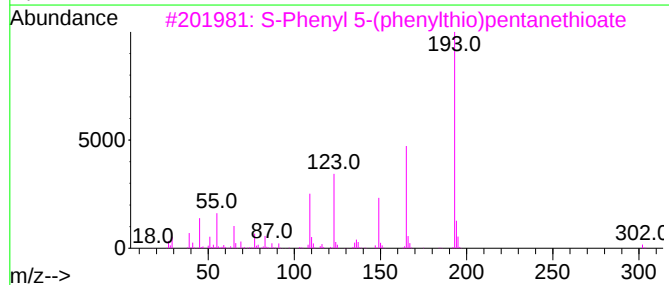
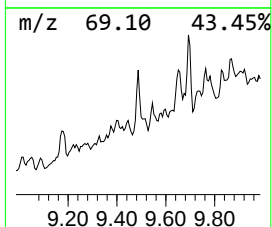
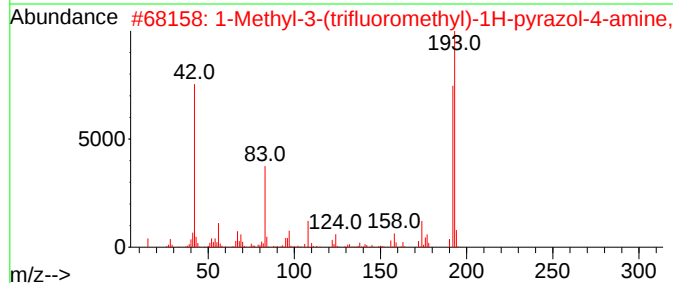
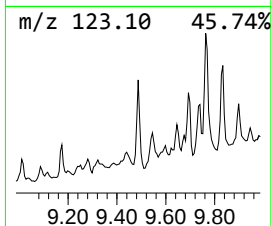
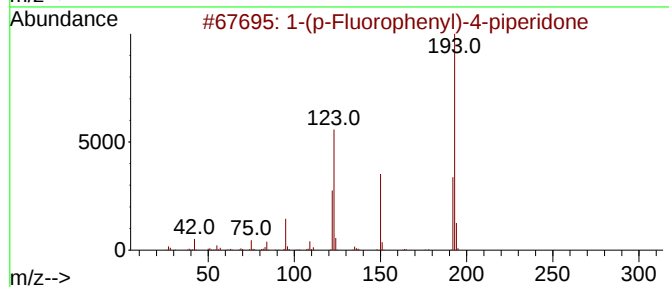
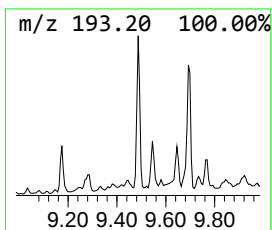
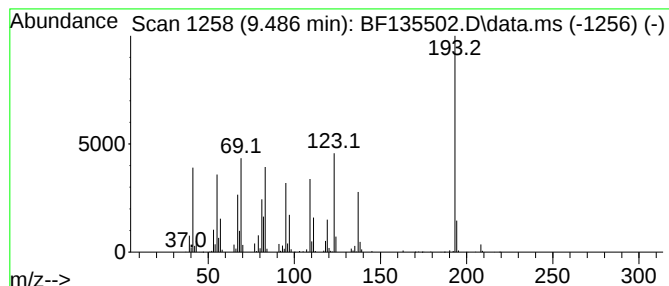
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown9.486 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.486	14.48 ng	1031580	Acenaphthene-d10	9.786

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		S-Phenyl 5-(phenylthio)pentaneth...	302	C17H18OS2	1000234-40-7	38
4		Borinic acid, diethyl-, 3,3,5-tr...	208	C13H25BO	057387-76-5	36
5		diallyl(2,2,4a,7,7-pentamethyl-1...	384	C19H33N2O4P	1000187-42-4	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092823\
Data File : BF135502.D
Acq On : 28 Sep 2023 19:39
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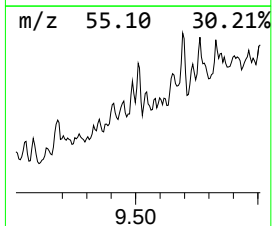
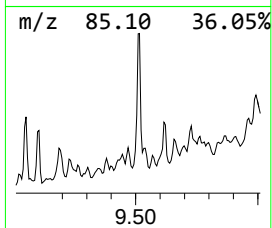
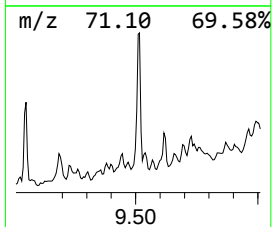
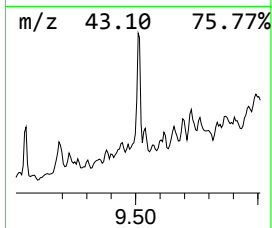
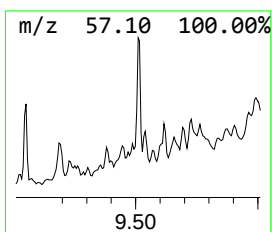
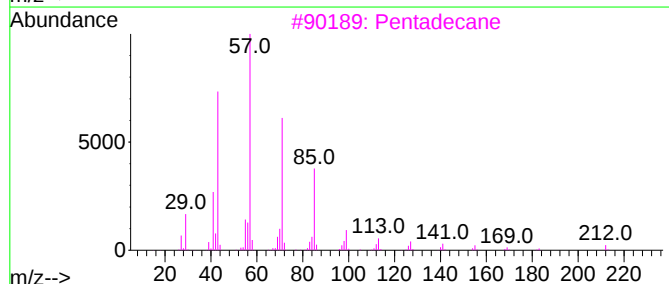
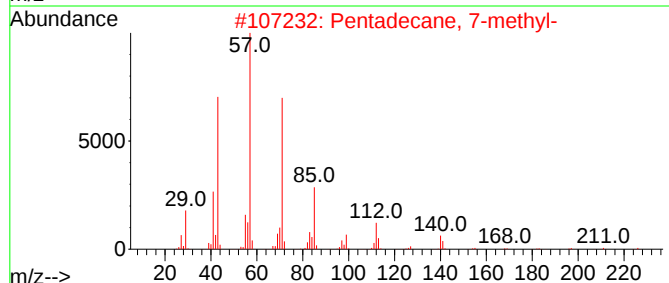
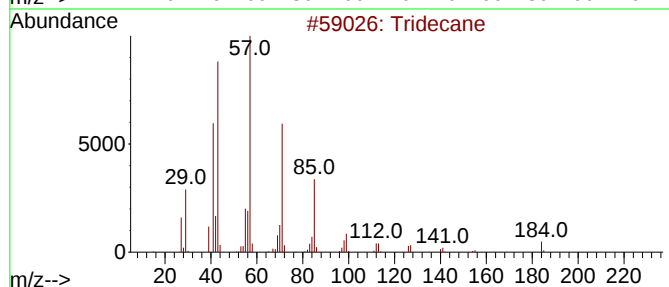
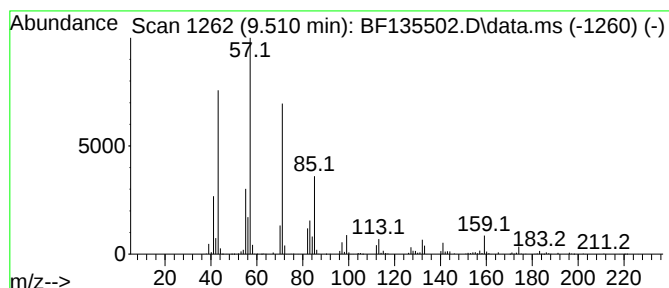
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tridecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.510	16.68 ng	1188050	Acenaphthene-d10	9.786

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane	184	C13H28	000629-50-5	72
2		Pentadecane, 7-methyl-	226	C16H34	006165-40-8	64
3		Pentadecane	212	C15H32	000629-62-9	64
4		Tetradecane	198	C14H30	000629-59-4	64
5		Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092823\
Data File : BF135502.D
Acq On : 28 Sep 2023 19:39
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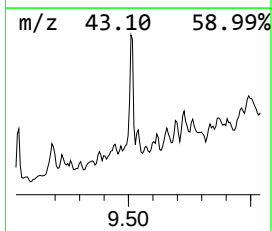
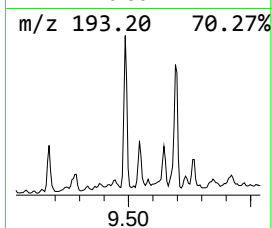
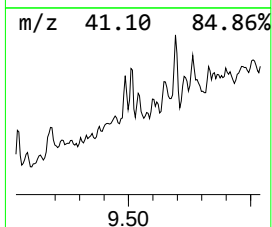
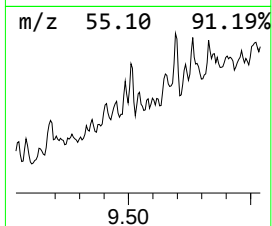
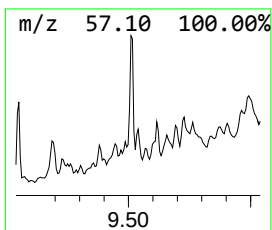
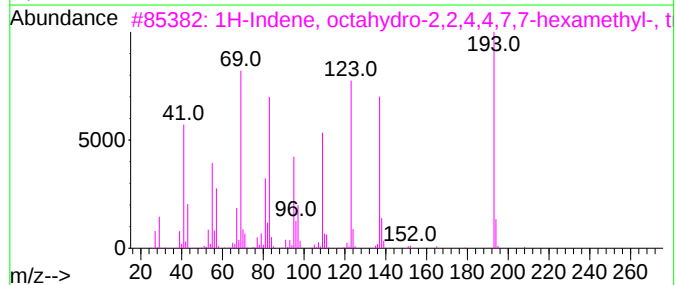
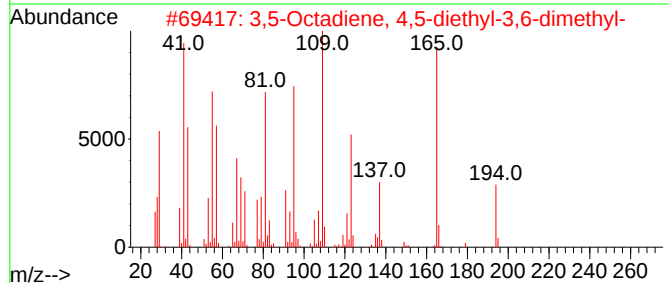
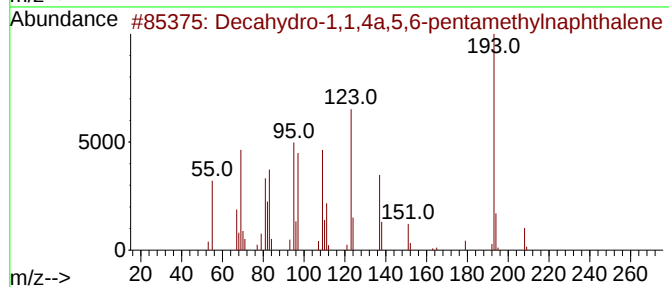
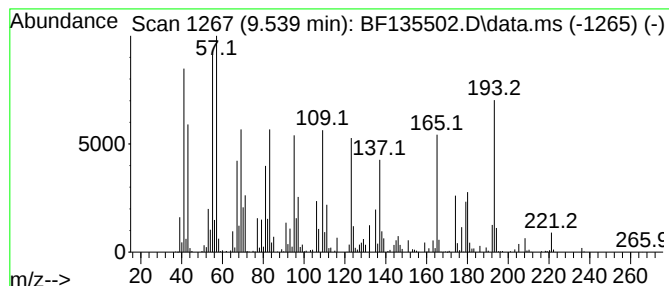
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Decahydro-1,1,4a,5,6-pentam... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.539	9.97 ng	710215	Acenaphthene-d10	9.786

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	52
2			3,5-Octadiene, 4,5-diethyl-3,6-d...	194	C14H26	061233-79-2	50
3			1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	46
4			4,6-Heptadien-2-one, 3,6-dimethy...	180	C12H20O	077822-50-5	30
5			1-(2-[3-(2-Acetyloxiran-2-yl)-1,...	236	C14H20O3	1010192-75-5	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092823\
Data File : BF135502.D
Acq On : 28 Sep 2023 19:39
Operator : CG\JU
Sample : 04606-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-327

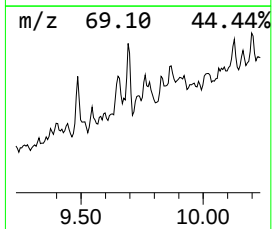
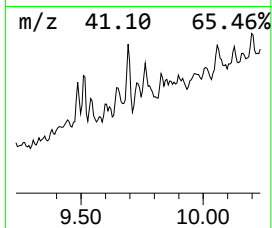
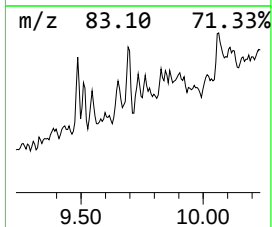
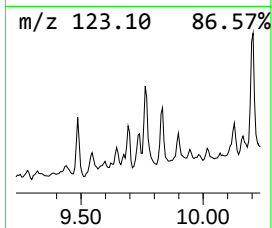
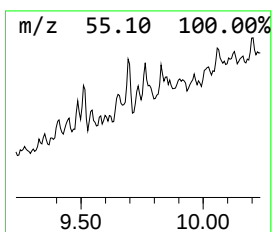
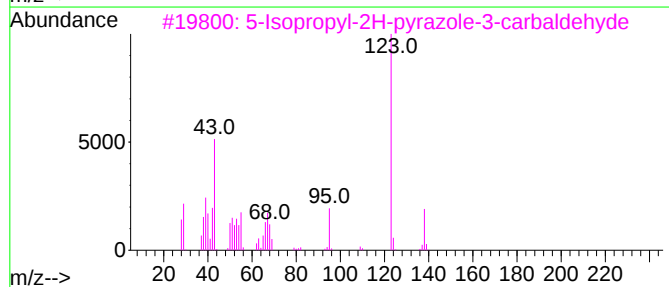
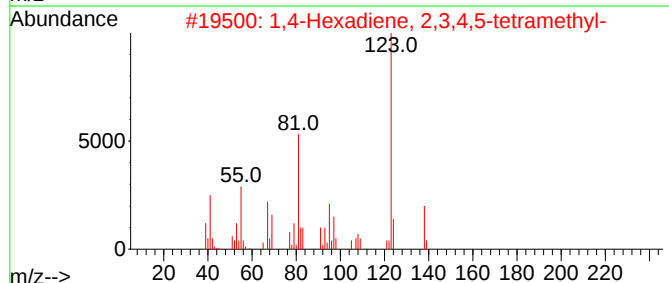
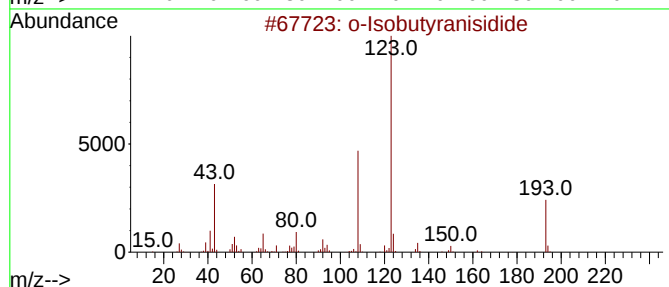
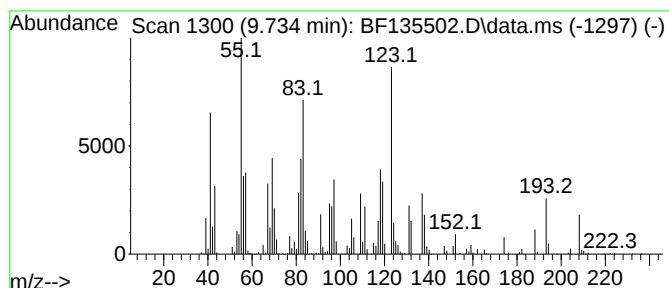
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown9.734 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.734	14.00 ng	997254	Acenaphthene-d10	9.786

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		o-Isobutyranisidide	193	C11H15NO2	071182-38-2	27
2		1,4-Hexadiene, 2,3,4,5-tetramethyl-	138	C10H18	051504-54-2	18
3		5-Isopropyl-2H-pyrazole-3-carbal...	138	C7H10N2O	1000410-48-4	18
4		o-Fluoroacetophenone	138	C8H7FO	000445-27-2	18
5		2(1H)-Pyridinone, 1,3-dimethyl-	123	C7H9NO	006456-92-4	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092823\
Data File : BF135502.D
Acq On : 28 Sep 2023 19:39
Operator : CG\JU
Sample : 04606-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-327

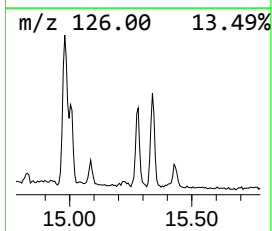
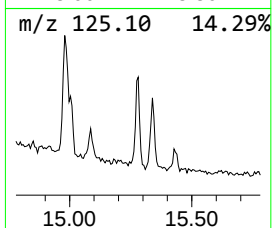
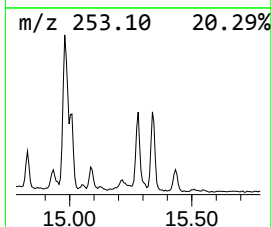
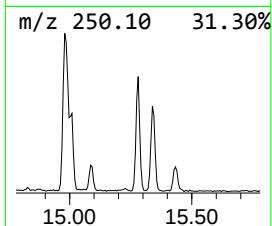
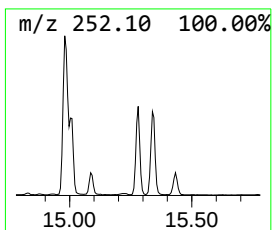
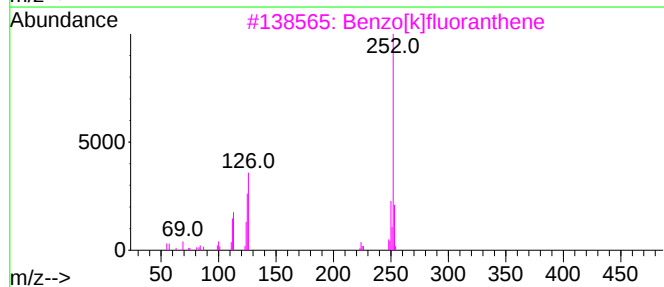
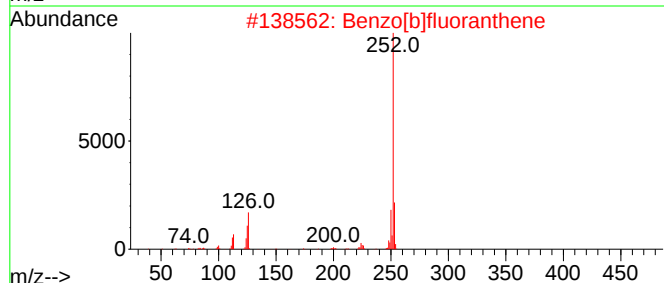
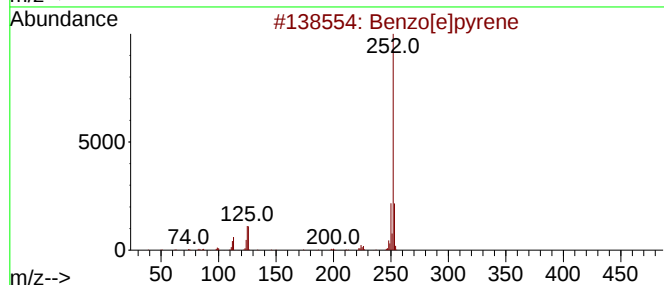
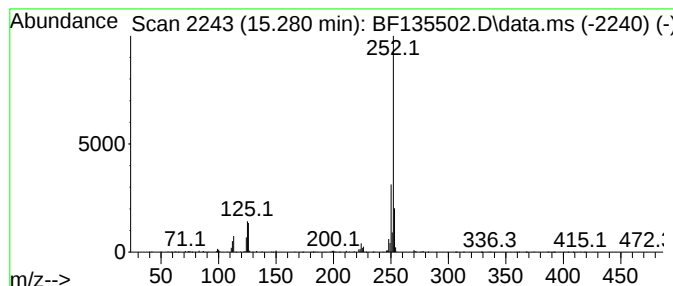
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.280	21.32 ng	570765	Perylene-d12	15.404

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[e]pyrene	252	C20H12	000192-97-2	98
2		Benzo[b]fluoranthene	252	C20H12	000205-99-2	97
3		Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
4		Benzo[a]pyrene	252	C20H12	000050-32-8	93
5		Perylene	252	C20H12	000198-55-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092823\
Data File : BF135502.D
Acq On : 28 Sep 2023 19:39
Operator : CG\JU
Sample : 04606-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-327

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.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
1-Hexanol, 2-et...	6.810	6.0	ng	238759	1	6.745	789091	20.0
Hexanoic acid, ...	7.434	6.0	ng	251134	2	8.022	840010	20.0
Camphor	7.769	2.5	ng	104663	2	8.022	840010	20.0
unknown8.310	8.310	3.3	ng	139693	2	8.022	840010	20.0
Sulfurous acid,...	8.445	6.2	ng	261111	2	8.022	840010	20.0
unknown8.557	8.557	4.8	ng	203749	2	8.022	840010	20.0
unknown8.904	8.904	4.4	ng	315612	3	9.786	1424450	20.0
Bicyclo[2.2.1]h...	9.010	5.4	ng	384865	3	9.786	1424450	20.0
Nonane, 3,7-dim...	9.051	11.2	ng	800207	3	9.786	1424450	20.0
unknown9.486	9.486	14.5	ng	1031580	3	9.786	1424450	20.0
Tridecane	9.510	16.7	ng	1188050	3	9.786	1424450	20.0
Decahydro-1,1,4...	9.539	10.0	ng	710215	3	9.786	1424450	20.0
unknown9.734	9.734	14.0	ng	997254	3	9.786	1424450	20.0
Benzo[e]pyrene	15.280	21.3	ng	570765	6	15.404	535355	20.0