

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110223\
 Data File : BF136102.D
 Acq On : 02 Nov 2023 18:30
 Operator : CG\JU
 Sample : 05049-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B1

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-BF103023.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF136102.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.040	498	502	510	rVB	59011	66613	1.86%	0.353%
2	5.440	566	570	574	rBV	1407535	1325878	37.05%	7.024%
3	6.445	737	741	749	rVB	897611	839922	23.47%	4.449%
4	6.816	801	804	808	rBV	782032	626764	17.51%	3.320%
5	6.887	812	816	822	rBV2	46266	81558	2.28%	0.432%
6	7.216	868	872	881	rBV	40638	48018	1.34%	0.254%
7	7.381	895	900	903	rBV	1969442	2012043	56.23%	10.658%
8	7.481	910	917	921	rBV	88461	122024	3.41%	0.646%
9	7.845	973	979	985	rVB3	116733	163920	4.58%	0.868%
10	8.010	1004	1007	1010	rBV	45894	49952	1.40%	0.265%
11	8.039	1010	1012	1018	rVB	88786	69180	1.93%	0.366%
12	8.098	1018	1022	1025	rBV	1108616	828898	23.16%	4.391%
13	8.122	1025	1026	1031	rVB	60580	48183	1.35%	0.255%
14	8.375	1066	1069	1072	rVB	93592	68424	1.91%	0.362%
15	8.451	1076	1082	1087	rBV	196201	212048	5.93%	1.123%
16	8.498	1087	1090	1097	rVB3	44011	69769	1.95%	0.370%
17	9.010	1173	1177	1180	rBV	47550	50280	1.41%	0.266%
18	9.086	1184	1190	1195	rBV	50707	71334	1.99%	0.378%
19	9.181	1202	1206	1209	rBV	4154519	3578517	100.00%	18.956%
20	9.857	1317	1321	1324	rBV	1136632	878632	24.55%	4.654%
21	10.275	1388	1392	1396	rBV2	68494	92938	2.60%	0.492%
22	10.651	1449	1456	1459	rBV	2200284	2014887	56.31%	10.673%
23	10.863	1489	1492	1496	rVB5	31847	40601	1.13%	0.215%
24	11.075	1525	1528	1532	rVB2	64365	60138	1.68%	0.319%
25	11.222	1550	1553	1558	rVB5	33644	53187	1.49%	0.282%
26	11.351	1571	1575	1578	rVB	855388	740862	20.70%	3.925%
27	11.469	1592	1595	1601	rBV3	64129	95790	2.68%	0.507%
28	11.527	1602	1605	1612	rVB2	59494	87190	2.44%	0.462%
29	11.892	1663	1667	1675	rBV2	222451	233049	6.51%	1.235%
30	12.680	1798	1801	1810	rVB	101660	109366	3.06%	0.579%
31	12.769	1813	1816	1820	rBV	281444	232483	6.50%	1.232%
32	12.951	1843	1847	1850	rBV	2856380	2517864	70.36%	13.338%
33	12.980	1850	1852	1858	rVB	143690	119724	3.35%	0.634%
34	13.745	1979	1982	1985	rBV	168329	128974	3.60%	0.683%
35	13.804	1989	1992	1997	rBV	55680	58370	1.63%	0.309%
36	13.898	2004	2008	2013	rBV3	39432	64950	1.81%	0.344%

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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-BF103023.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	14.004	2022	2026	2030	rBV	616481	537225	15.01%	2.846%
38	15.492	2274	2279	2287	rBV	407879	478073	13.36%	2.532%

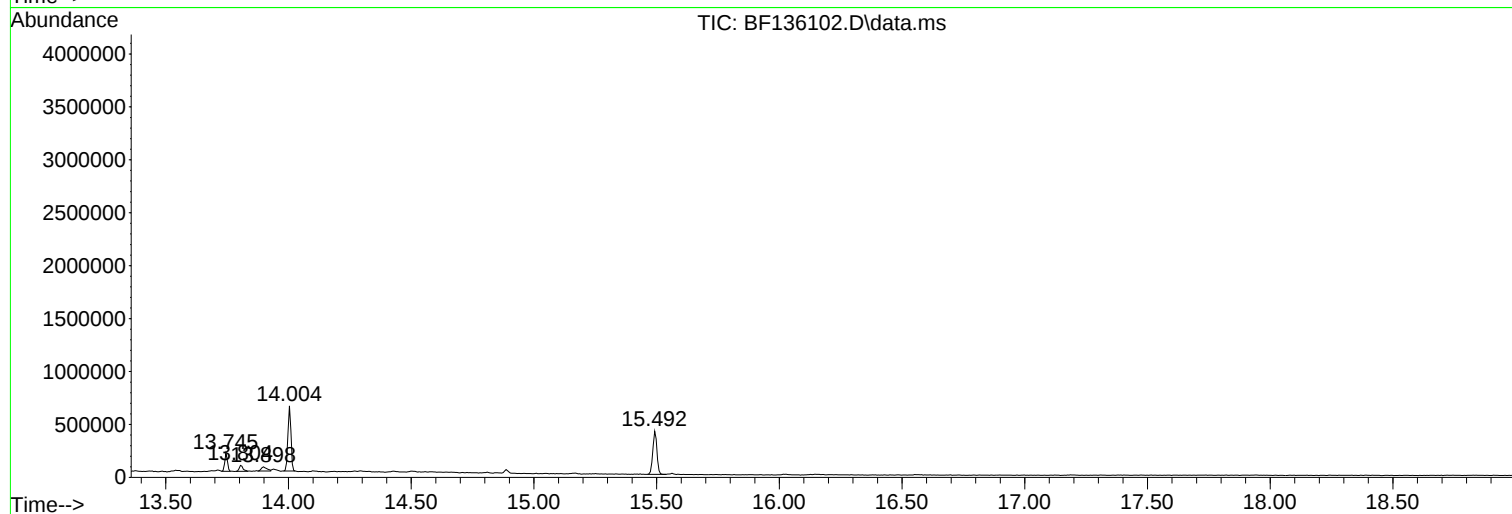
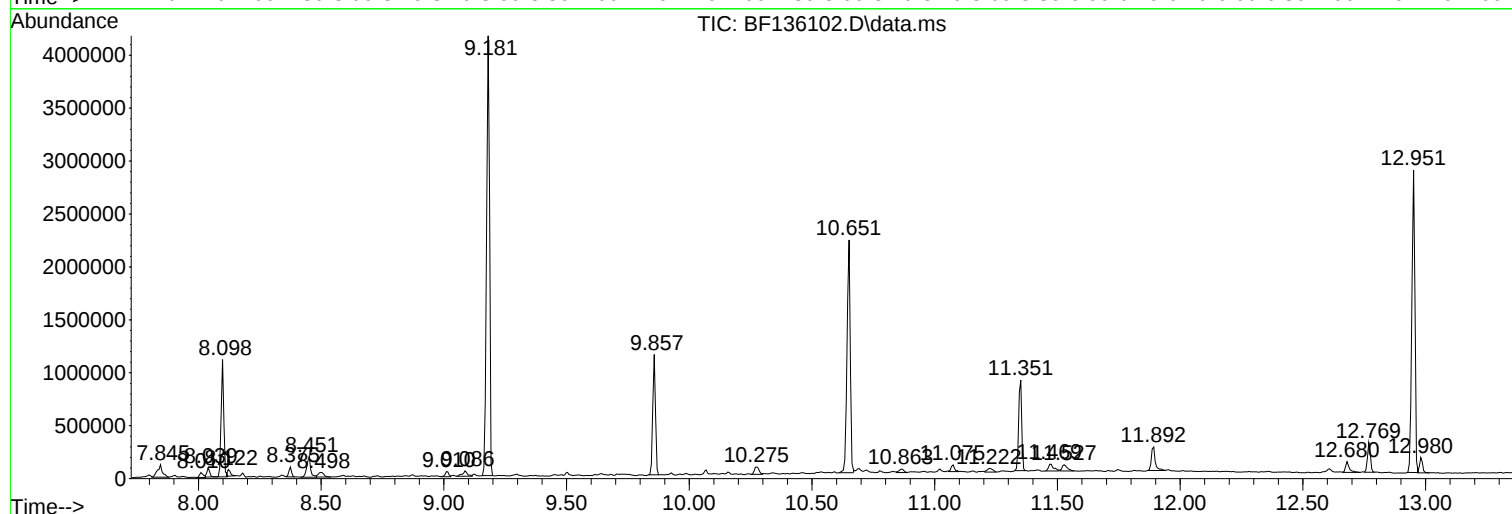
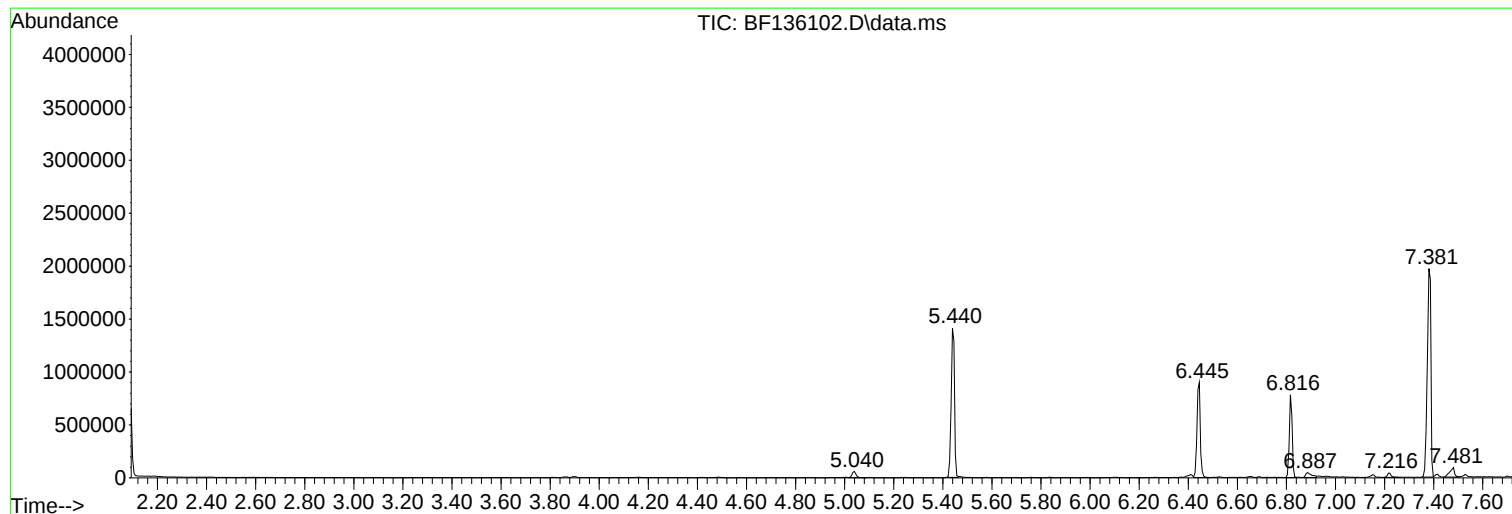
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF103023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Operator : CG\JU
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Misc :
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Instrument :
BNA_F
ClientSampleId :
B1

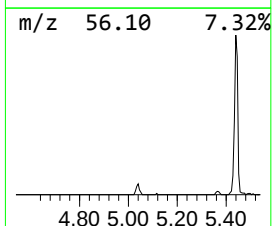
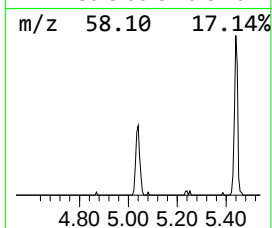
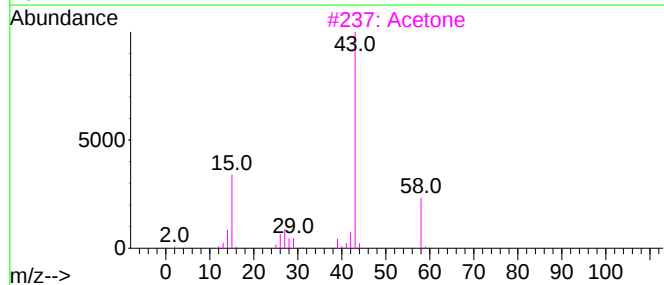
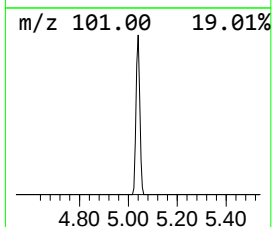
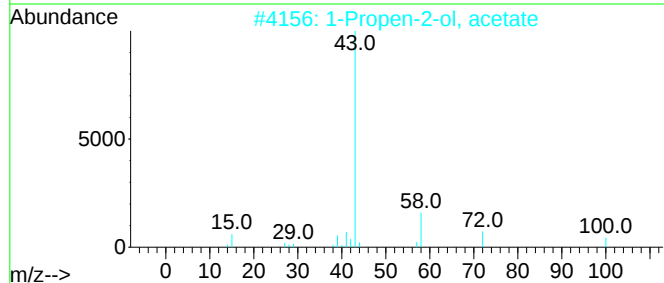
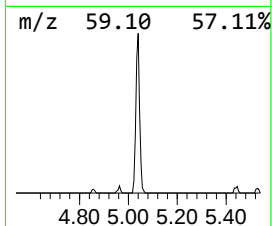
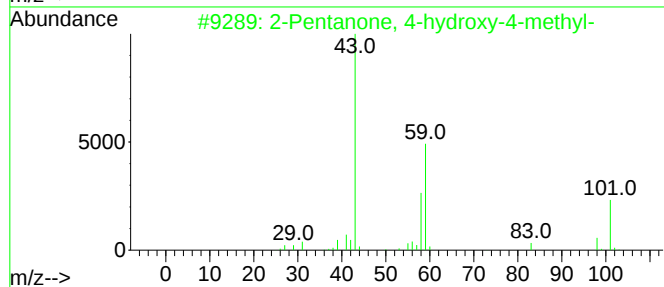
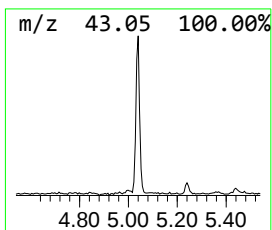
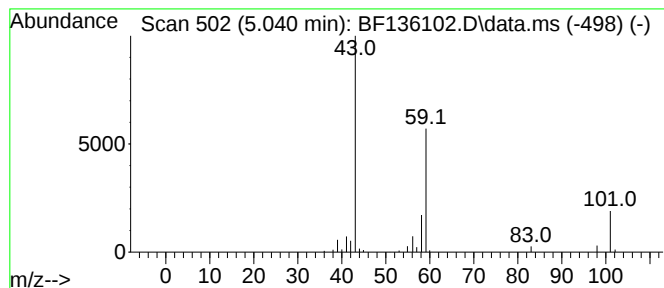
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF103023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.040	2.13 ng	66613	1,4-Dichlorobenzene-d4	6.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64		
2	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12		
3	Acetone	58	C3H6O	000067-64-1	12		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9		
5	Acetamide, N-(cyanomethyl)-	98	C4H6N2O	004814-80-6	9		



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Instrument :
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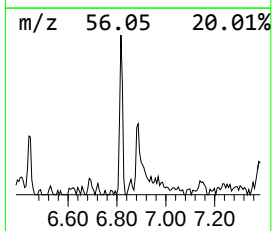
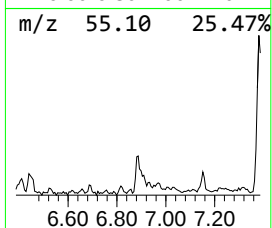
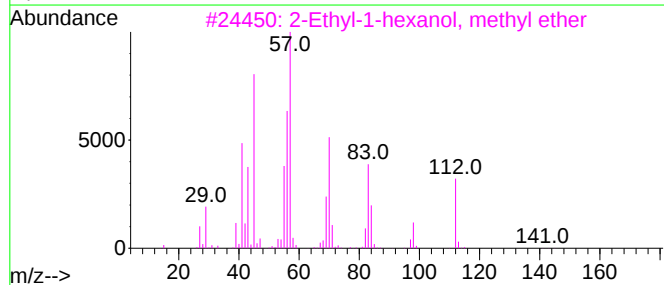
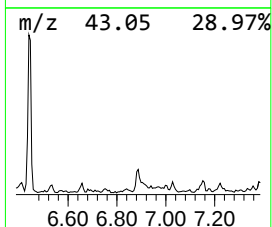
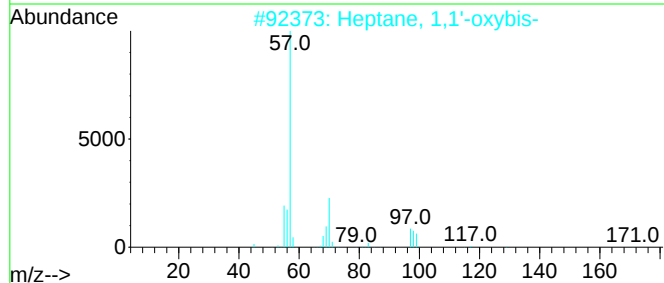
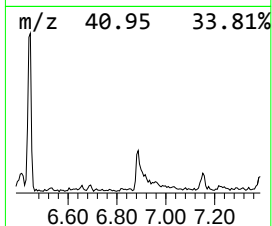
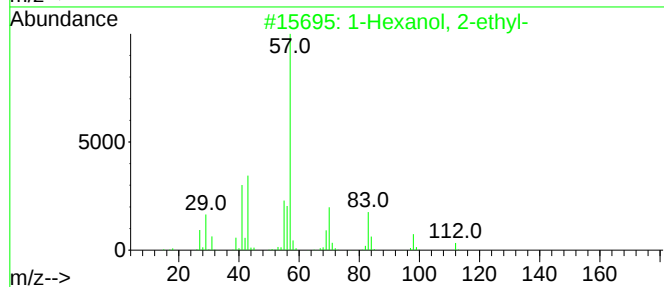
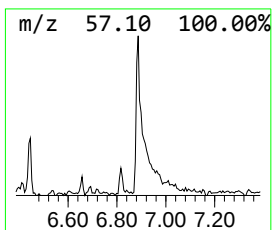
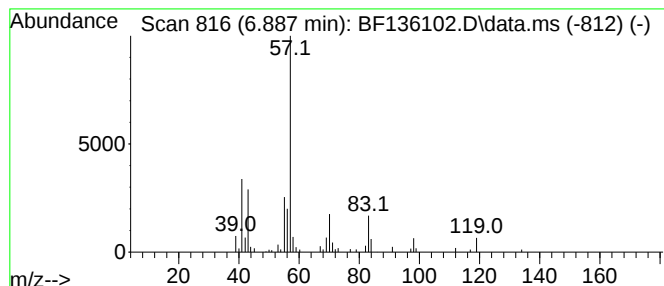
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF103023.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.887	2.60 ng	81558	1,4-Dichlorobenzene-d4	6.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	72
2			Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	53
3			2-Ethyl-1-hexanol, methyl ether	144	C9H20O	1000365-10-2	50
4			Oxalic acid, isobutyl heptyl ester	244	C13H24O4	1000309-37-2	47
5			1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	47



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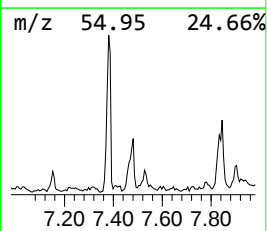
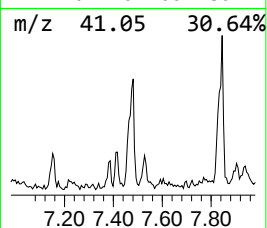
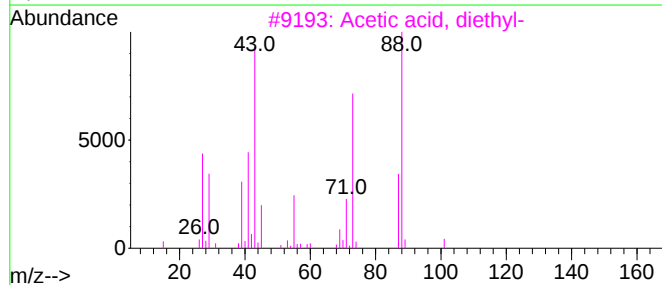
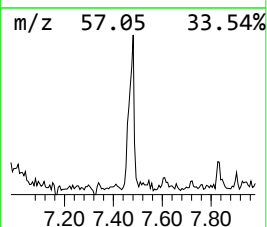
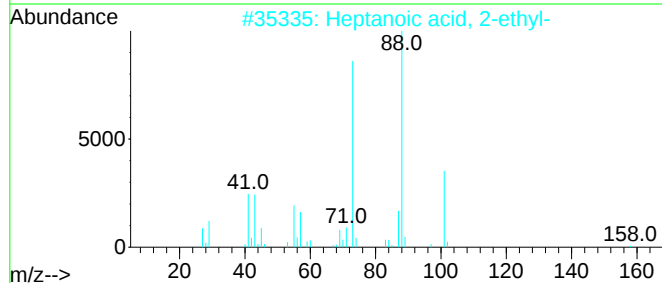
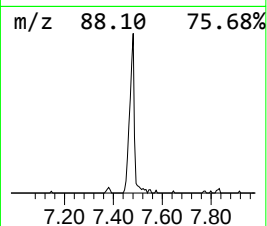
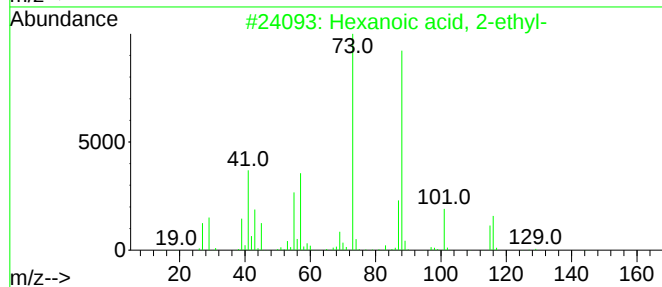
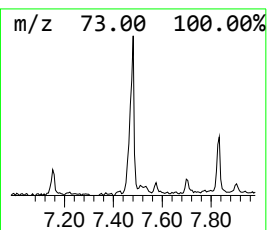
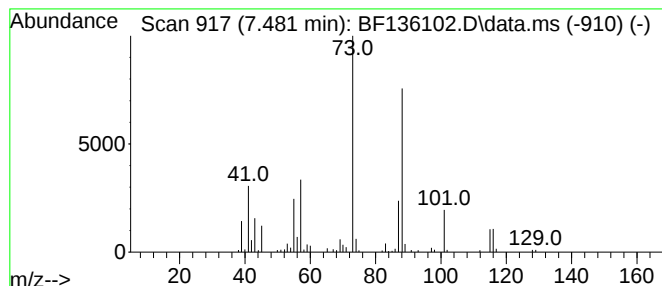
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF103023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Hexanoic acid, 2-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.481	2.94 ng	122024	Naphthalene-d8	8.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	80
2			Heptanoic acid, 2-ethyl-	158	C9H18O2	003274-29-1	53
3			Acetic acid, diethyl-	116	C6H12O2	000088-09-5	53
4			Butanoic acid, 2-methyl-, methyl...	116	C6H12O2	000868-57-5	47
5			2-Ethyl-4-methylpentanoic acid	144	C8H16O2	000108-81-6	45



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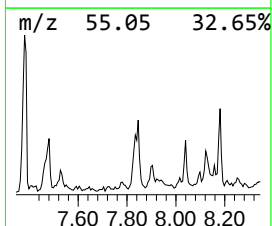
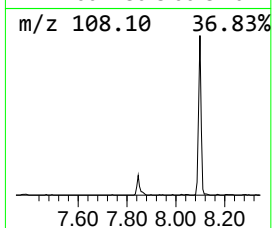
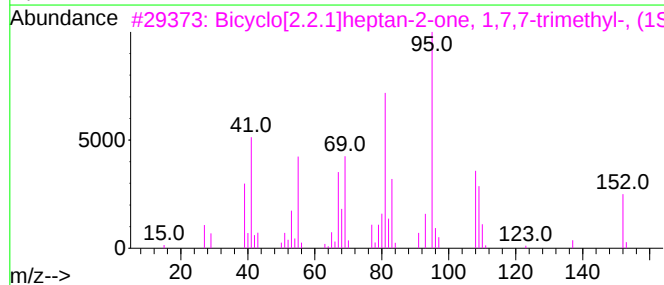
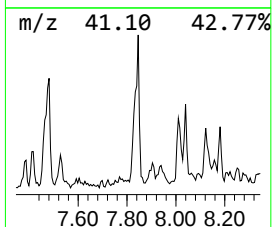
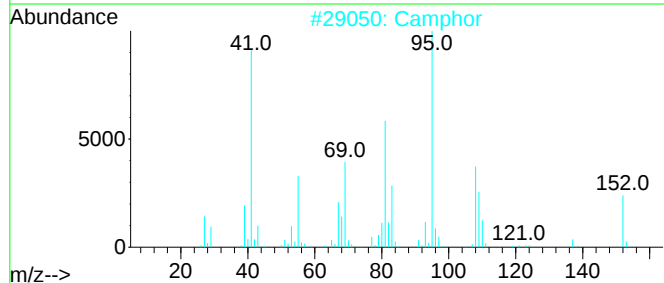
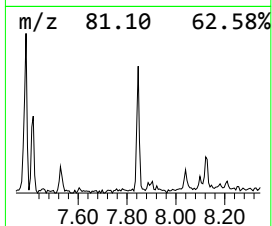
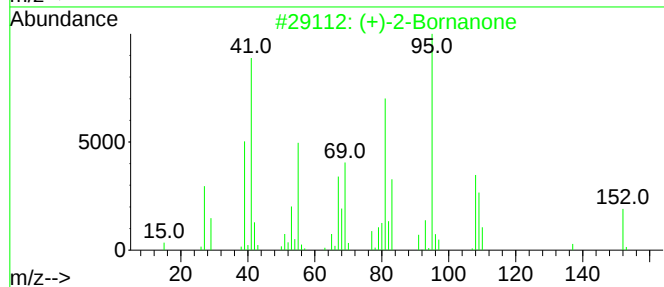
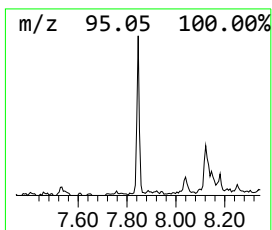
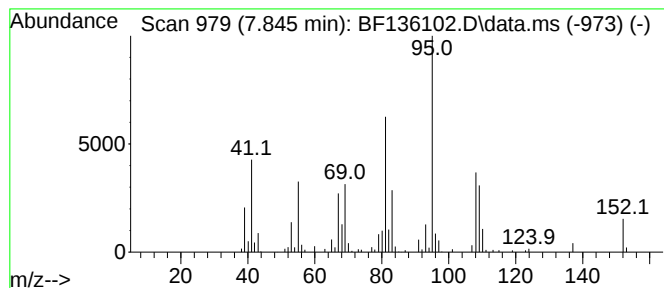
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TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (+)-2-Bornanone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.845	3.96 ng	163920	Naphthalene-d8	8.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(+)-2-Bornanone	152	C10H16O	000464-49-3	96
2			Camphor	152	C10H16O	000076-22-2	95
3			Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	94
4			1,3-Dimethyl-1-cyclohexene	110	C8H14	002808-76-6	46
5			Bicyclo[2.2.1]heptane-2-carboxal...	124	C8H12O	003574-55-8	43



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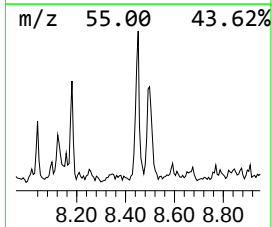
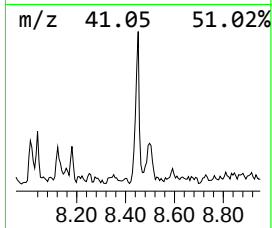
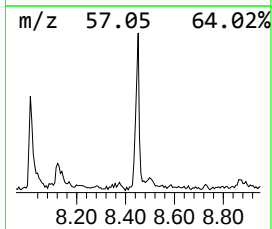
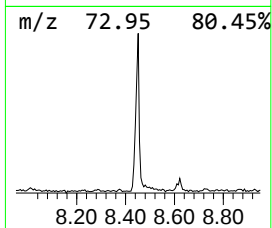
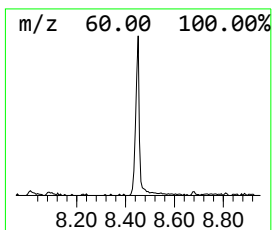
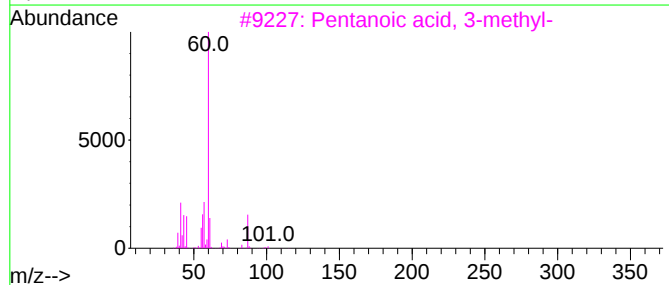
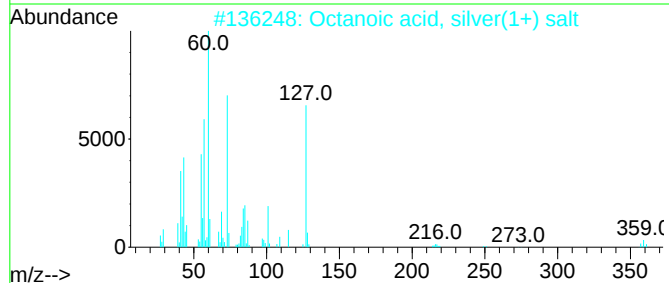
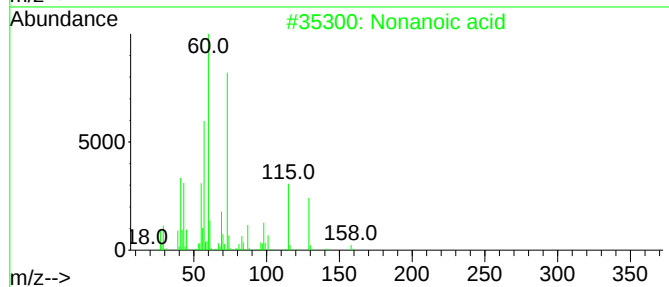
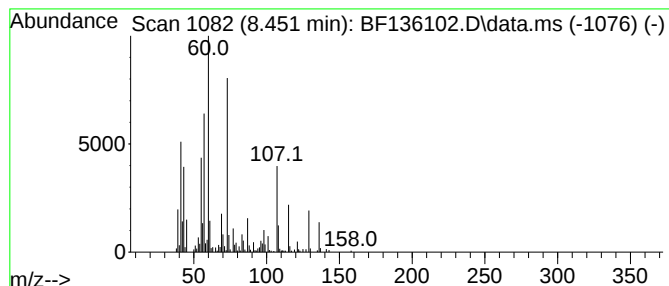
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Nonanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.451	5.12 ng	212048	Naphthalene-d8	8.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonanoic acid	158	C9H18O2	000112-05-0	62
2			Octanoic acid, silver(1+) salt	250	C8H15AgO2	024927-67-1	38
3			Pentanoic acid, 3-methyl-	116	C6H12O2	000105-43-1	35
4			Hexanoic acid	116	C6H12O2	000142-62-1	35
5			Butanoic acid	88	C4H8O2	000107-92-6	35



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Data File : BF136102.D
Acq On : 02 Nov 2023 18:30
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Sample : 05049-01
Misc :
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Instrument :
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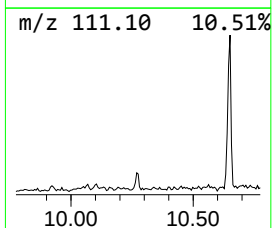
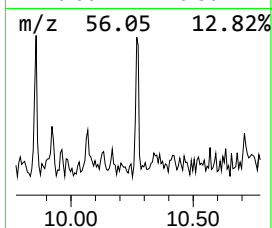
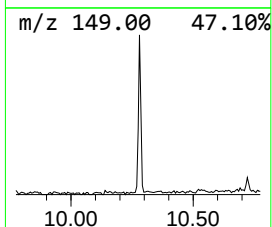
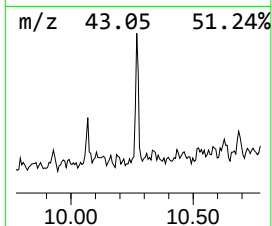
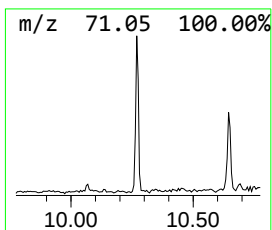
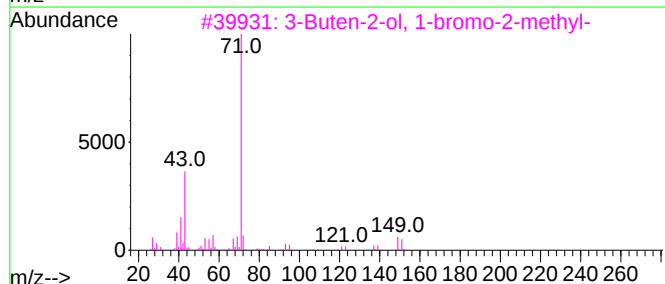
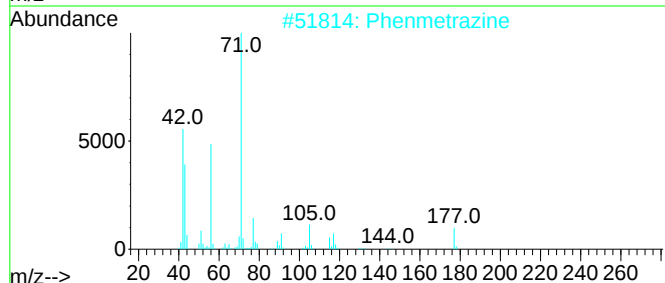
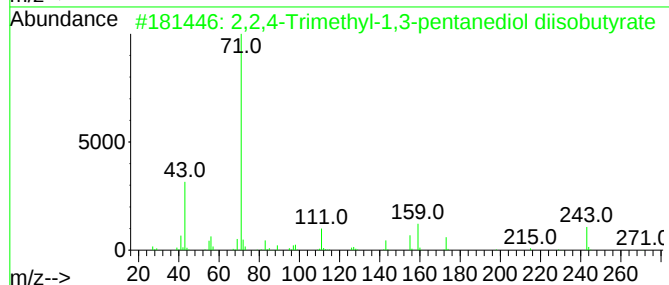
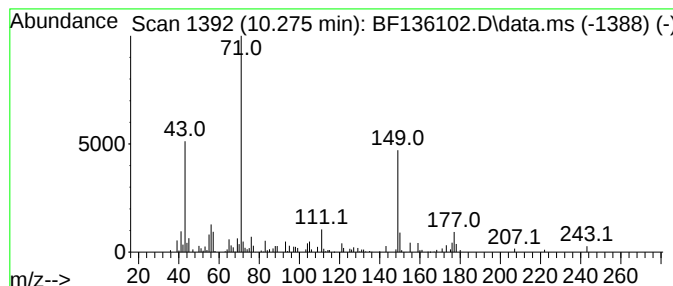
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown10.275 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.275	2.12 ng	92938	Acenaphthene-d10	9.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2,4-Trimethyl-1,3-pentanediol ...	286	C16H30O4	006846-50-0	47		
2	Phenmetrazine	177	C11H15NO	000134-49-6	38		
3	3-Buten-2-ol, 1-bromo-2-methyl-	164	C5H9BrO	036219-40-6	38		
4	Pentane, 1-bromo-	150	C5H11Br	000110-53-2	38		
5	Cyclohexanol, 2,2-dichloro-1-met...	182	C7H12Cl2O	1000115-65-2	38		



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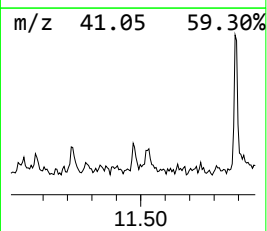
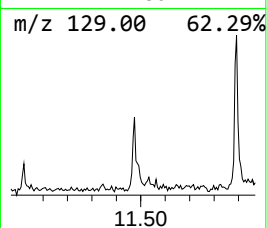
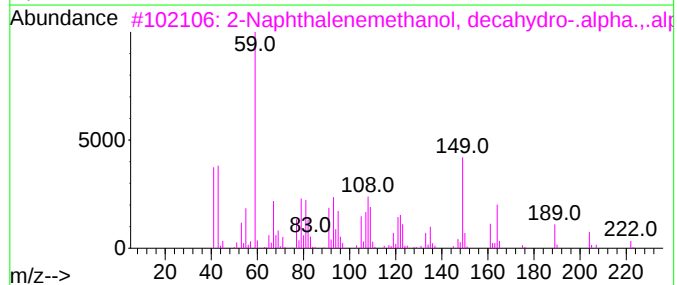
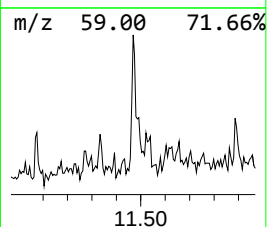
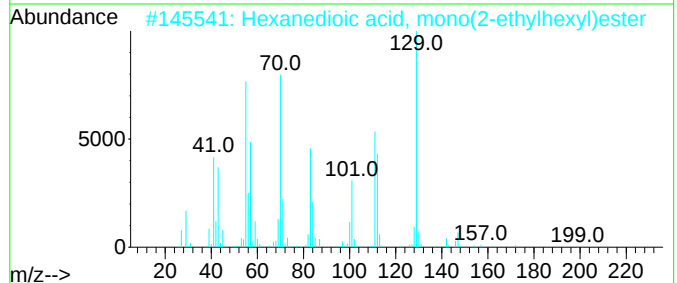
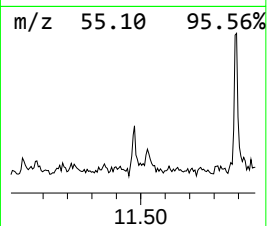
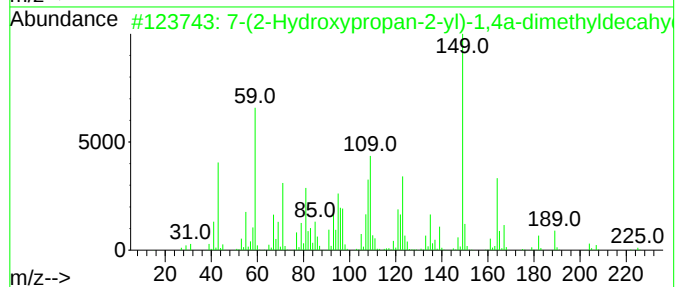
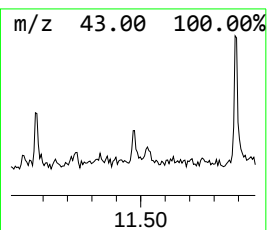
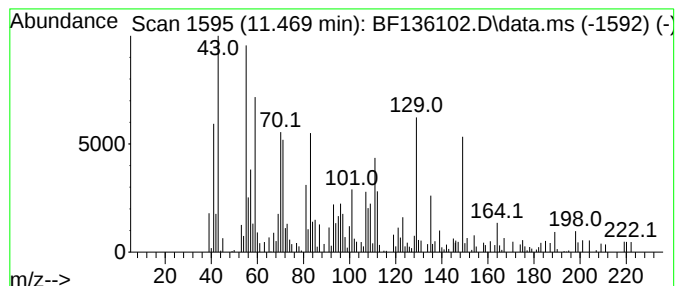
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown11.469 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.469	2.59 ng	95790	Phenanthrene-d10	11.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-(2-Hydroxypropan-2-yl)-1,4a-di...	240	C15H28O2	092857-25-5	27		
2	Hexanedioic acid, mono(2-ethylhe...	258	C14H26O4	004337-65-9	25		
3	2-Naphthalenemethanol, decahydro...	222	C15H26O	000473-15-4	25		
4	2-Naphthalenemethanol, 1,2,3,4,4...	222	C15H26O	000473-16-5	22		
5	Cyclohexanecarboxylic acid, cycl...	196	C12H20O2	005420-91-7	14		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110223\
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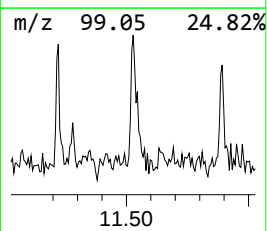
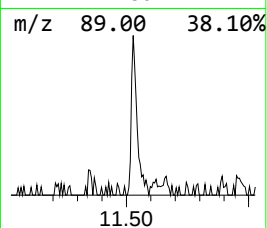
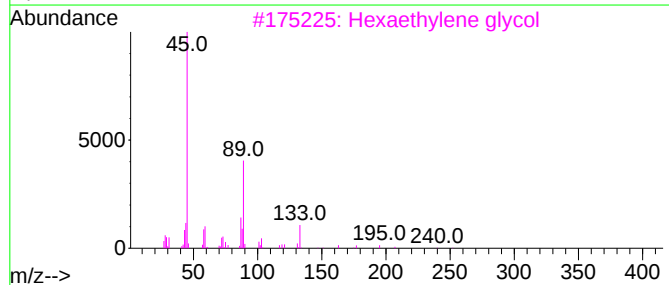
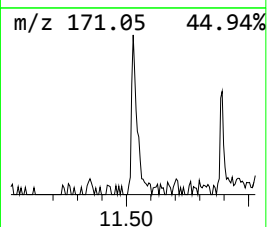
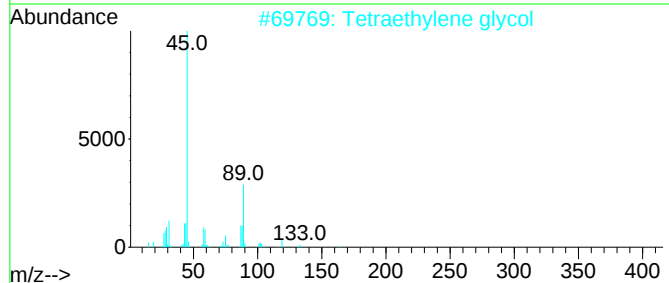
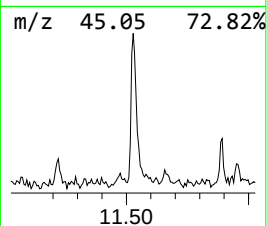
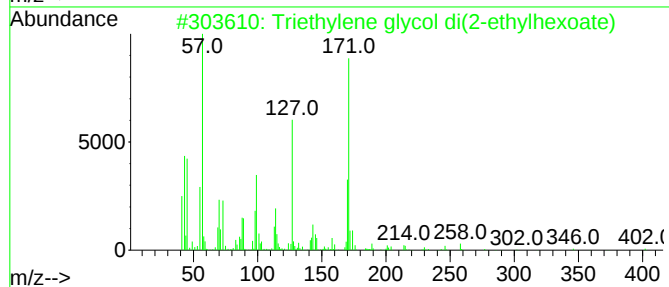
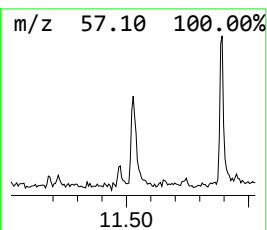
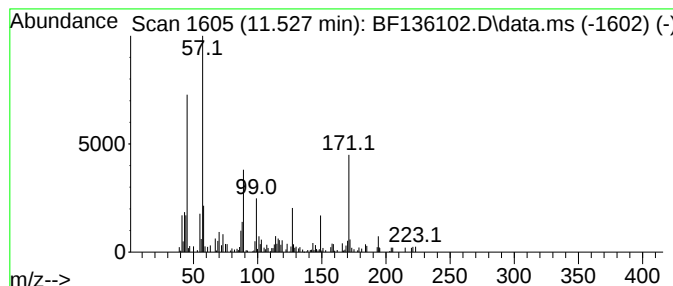
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Triethylene glycol di(2-eth... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.528	2.35 ng	87190	Phenanthrene-d10	11.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triethylene glycol di(2-ethylhex...	402	C22H42O6	000094-28-0	50
2			Tetraethylene glycol	194	C8H18O5	000112-60-7	27
3			Hexaethylene glycol	282	C12H26O7	002615-15-8	16
4			Heptaethylene glycol	326	C14H30O8	005617-32-3	16
5			2-{2-[2-(Hexyloxy)ethoxy]ethoxy}...	234	C12H26O4	025961-89-1	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110223\
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ALS Vial : 13 Sample Multiplier: 1

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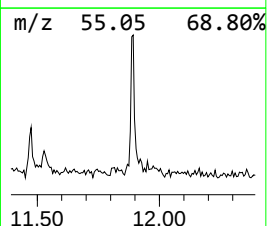
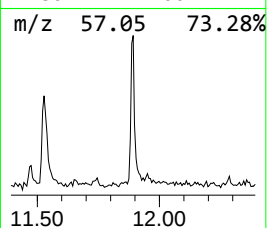
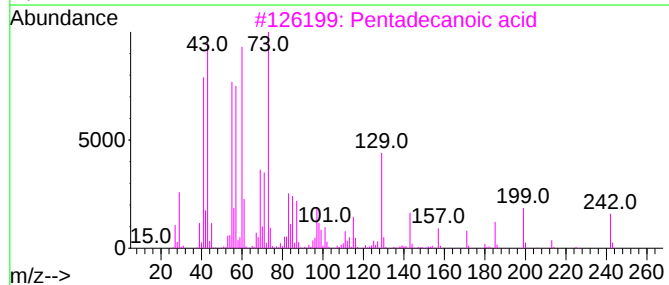
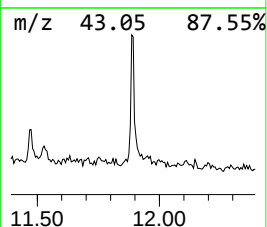
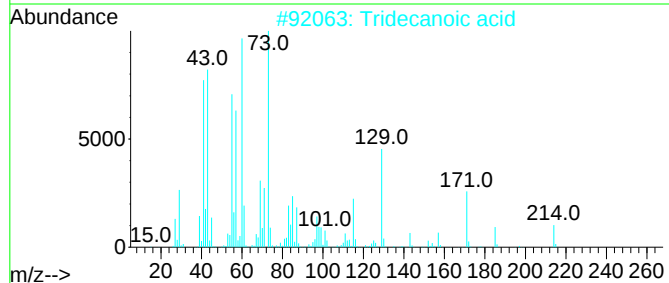
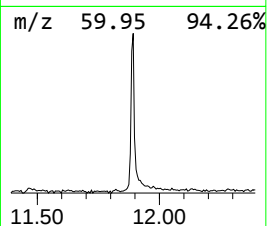
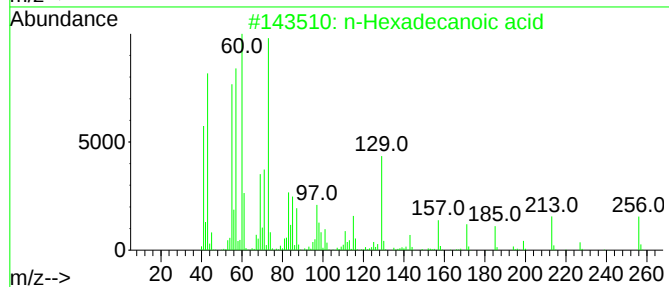
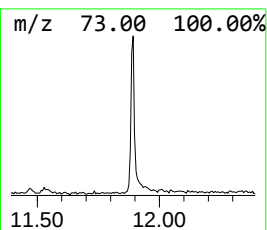
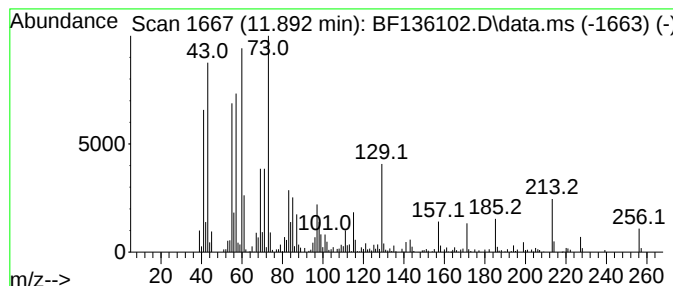
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.892	6.29 ng	233049	Phenanthrene-d10	11.351

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	87
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5			n-Decanoic acid	172	C10H20O2	000334-48-5	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110223\
Data File : BF136102.D
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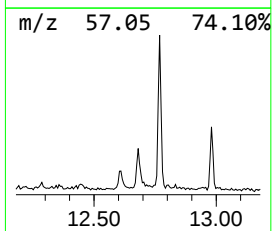
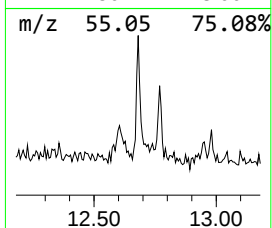
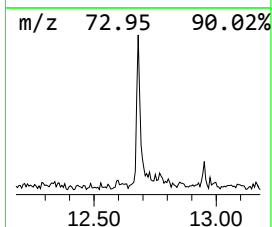
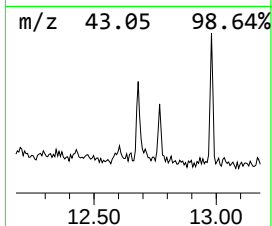
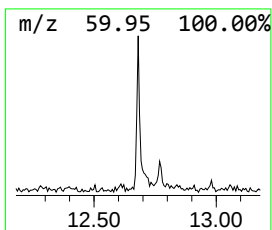
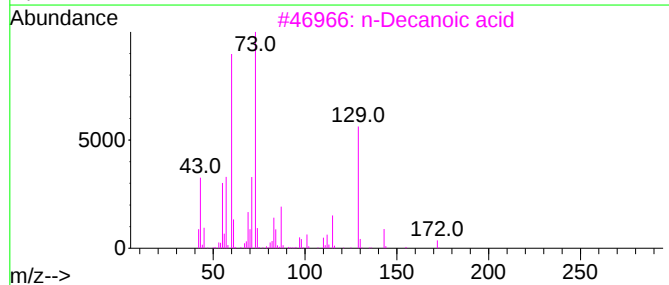
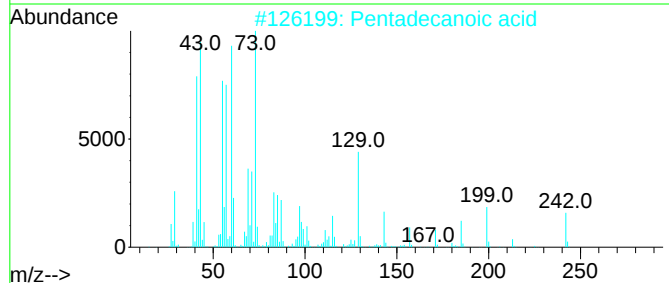
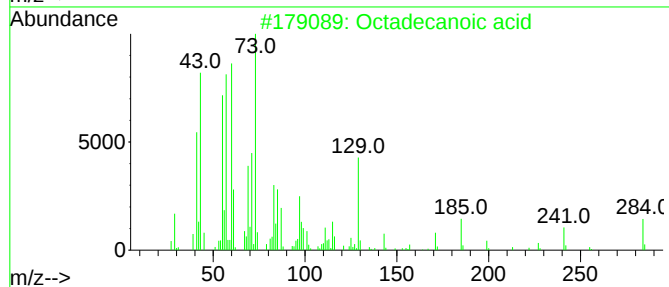
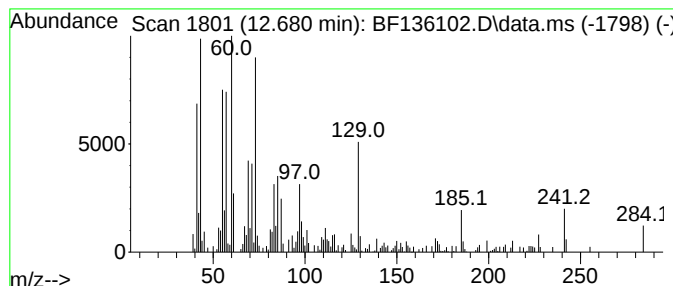
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.680	4.07 ng	109366	Chrysene-d12	14.004

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	95
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3		n-Decanoic acid	172	C10H20O2	000334-48-5	86
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	74
5		Undecanoic acid	186	C11H22O2	000112-37-8	64



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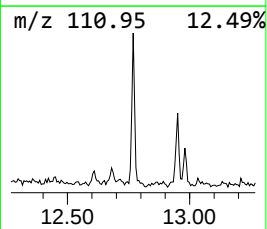
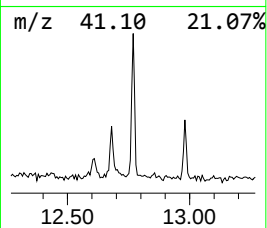
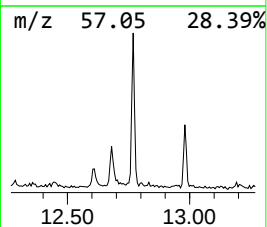
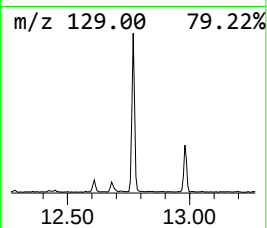
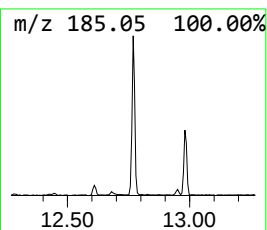
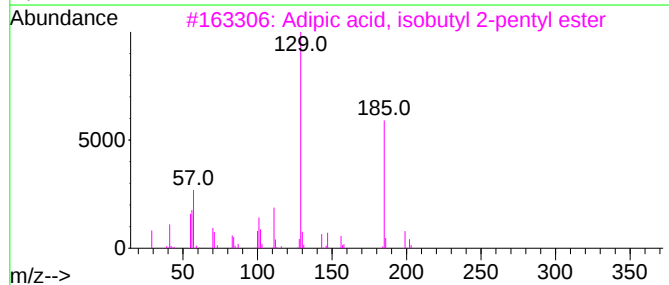
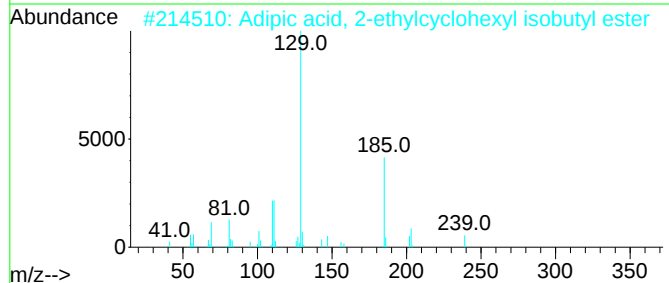
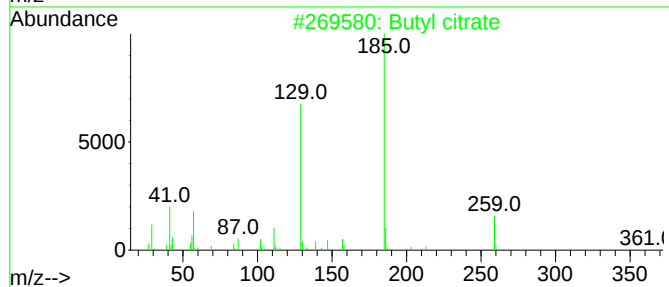
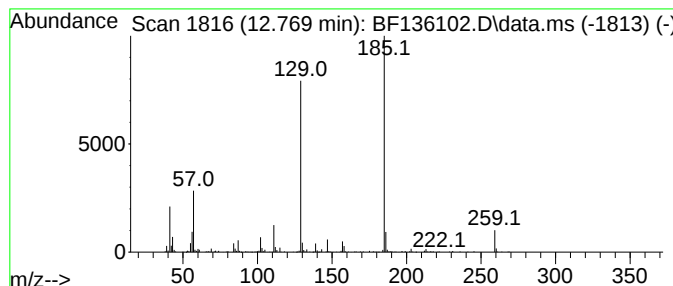
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Butyl citrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.769	8.65 ng	232483	Chrysene-d12	14.004

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butyl citrate	360	C18H32O7	000077-94-1	91
2			Adipic acid, 2-ethylcyclohexyl i...	312	C18H32O4	1000324-22-5	72
3			Adipic acid, isobutyl 2-pentyl e...	272	C15H28O4	1000324-15-6	64
4			Hexanedioic acid, bis(2-methylpr...	258	C14H26O4	000141-04-8	56
5			Adipic acid, 2-decyl isobutyl ester	342	C20H38O4	1000353-59-6	56



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ALS Vial : 13 Sample Multiplier: 1

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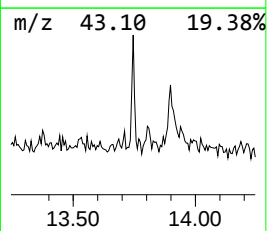
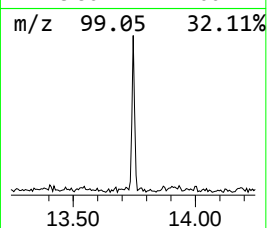
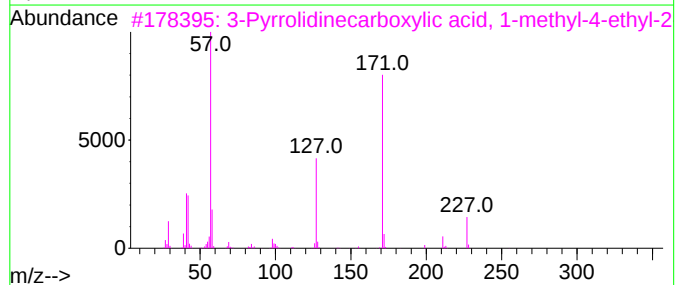
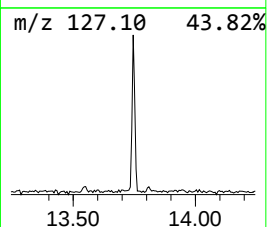
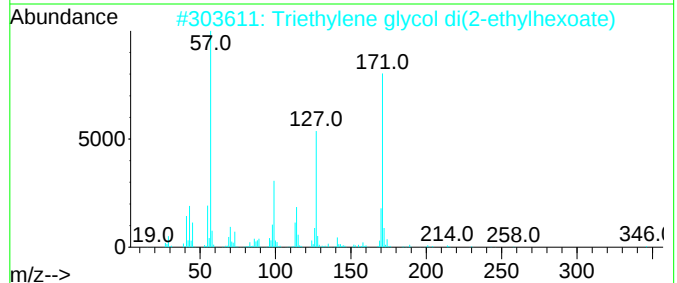
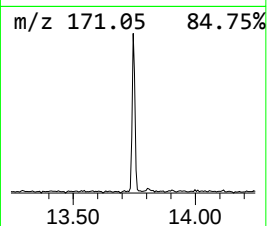
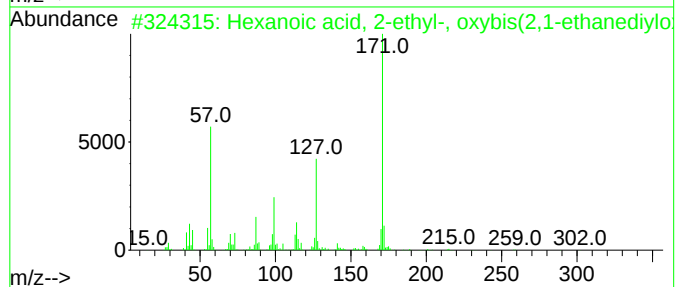
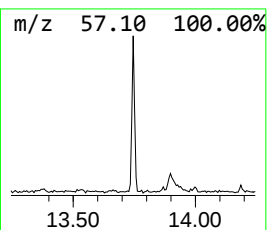
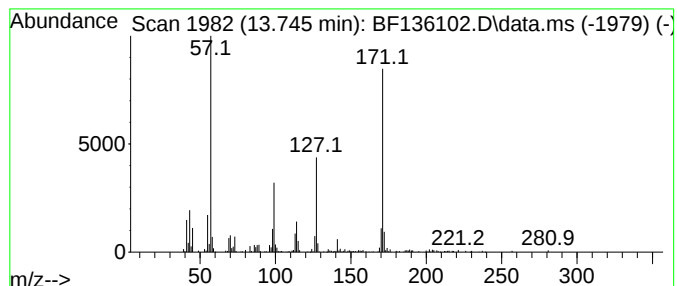
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexanoic acid, 2-ethyl-, ox... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.745	4.80 ng	128974	Chrysene-d12	14.004

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-, oxybis(...	446	C24H46O7	018268-70-7	91
2			Triethylene glycol di(2-ethylhex...	402	C22H42O6	000094-28-0	53
3			3-Pyrrolidinecarboxylic acid, 1-...	284	C15H28N2O3	1000196-95-3	45
4			1H-pyrrole-3-carboxaldehyde, 1-p...	171	C11H9NO	1000404-60-6	35
5			Vinyl 2-ethylhexanoate	170	C10H18O2	000094-04-2	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110223\
Data File : BF136102.D
Acq On : 02 Nov 2023 18:30
Operator : CG\JU
Sample : 05049-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B1

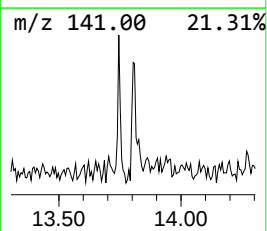
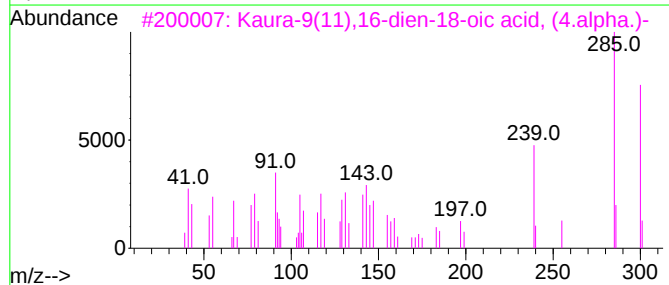
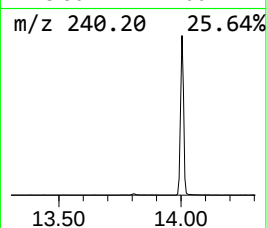
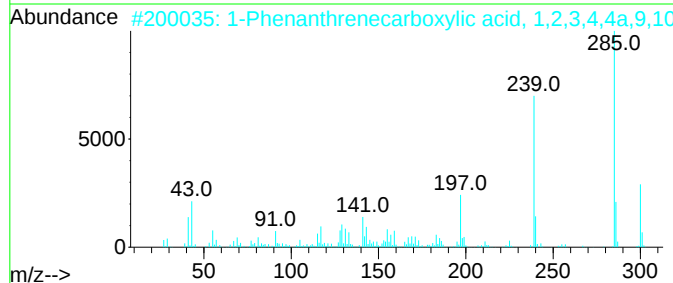
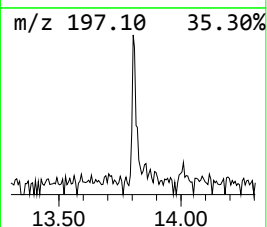
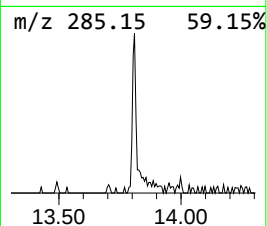
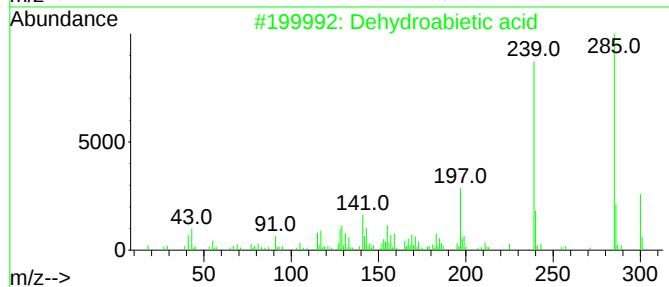
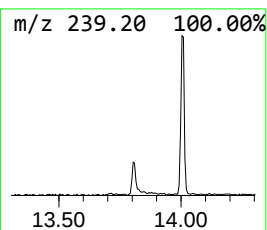
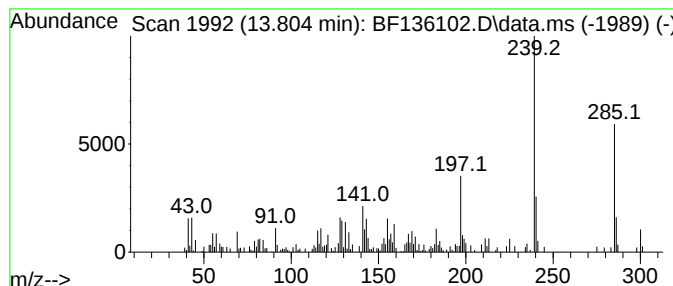
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Dehydroabiatic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.804	2.17 ng	58370	Chrysene-d12	14.004

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dehydroabiatic acid	300	C20H28O2	001740-19-8	96
2			1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	93
3			Kaura-9(11),16-dien-18-oic acid,...	300	C20H28O2	022338-67-6	64
4			Ethyl 4-hydroxy-7-trifluoromethy...	285	C13H10F3NO3	000391-02-6	52
5			Butanoic acid, 2-(cyano)(2,4,6-t...	300	C17H20N2O3	1000267-71-7	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110223\
Data File : BF136102.D
Acq On : 02 Nov 2023 18:30
Operator : CG\JU
Sample : 05049-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B1

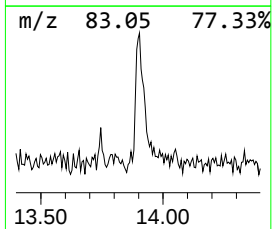
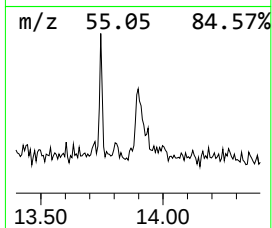
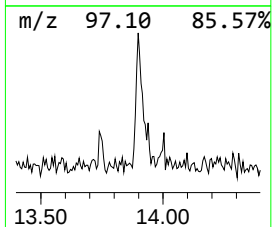
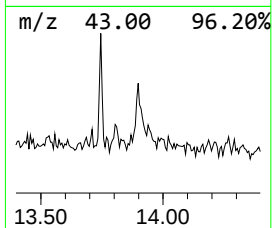
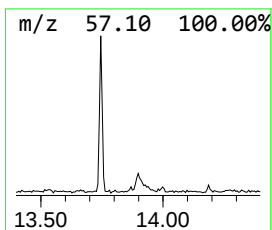
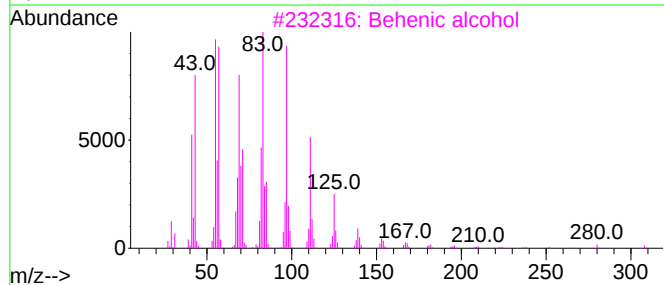
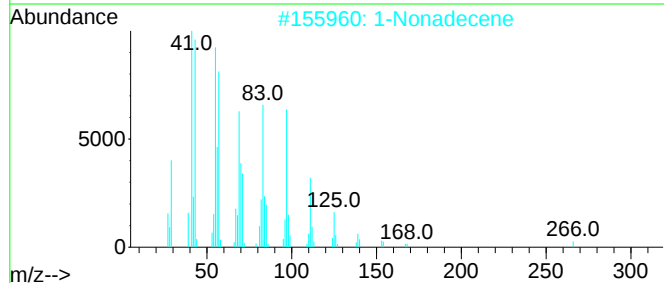
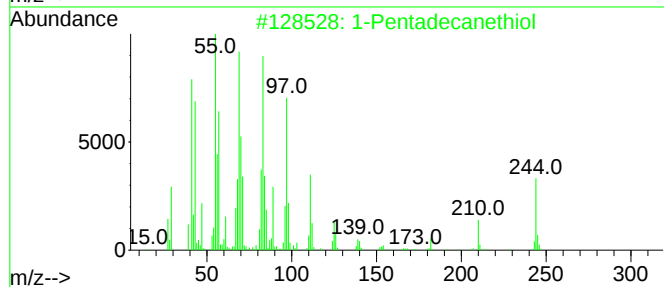
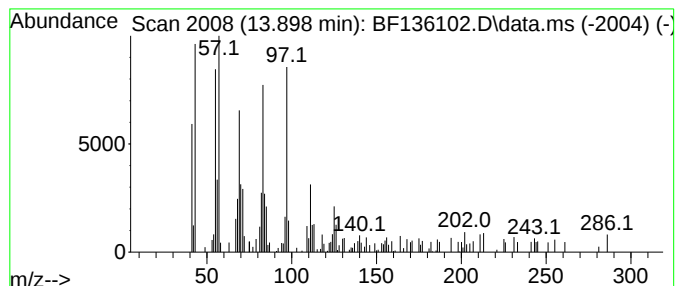
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1-Pentadecanethiol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.898	2.42 ng	64950	Chrysene-d12	14.004

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Pentadecanethiol	244	C15H32S	025276-70-4	95
2		1-Nonadecene	266	C19H38	018435-45-5	90
3		Behenic alcohol	326	C22H46O	000661-19-8	70
4		1-Octadecanol	270	C18H38O	000112-92-5	70
5		1-Heneicosanol	312	C21H44O	015594-90-8	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110223\
Data File : BF136102.D
Acq On : 02 Nov 2023 18:30
Operator : CG\JU
Sample : 05049-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF103023.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-...	5.040	2.1	ng	66613	1	6.816	626764	20.0
1-Hexanol, 2-et...	6.887	2.6	ng	81558	1	6.816	626764	20.0
Hexanoic acid, ...	7.481	2.9	ng	122024	2	8.098	828898	20.0
(+)-2-Bornanone	7.845	4.0	ng	163920	2	8.098	828898	20.0
Nonanoic acid	8.451	5.1	ng	212048	2	8.098	828898	20.0
unknown10.275	10.275	2.1	ng	92938	3	9.857	878632	20.0
unknown11.469	11.469	2.6	ng	95790	4	11.351	740862	20.0
Triethylene gly...	11.528	2.4	ng	87190	4	11.351	740862	20.0
n-Hexadecanoic ...	11.892	6.3	ng	233049	4	11.351	740862	20.0
Octadecanoic acid	12.680	4.1	ng	109366	5	14.004	537225	20.0
Butyl citrate	12.769	8.7	ng	232483	5	14.004	537225	20.0
Hexanoic acid, ...	13.745	4.8	ng	128974	5	14.004	537225	20.0
Dehydroabietic ...	13.804	2.2	ng	58370	5	14.004	537225	20.0
1-Pentadecanethiol	13.898	2.4	ng	64950	5	14.004	537225	20.0