

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062723\
 Data File : BF133996.D
 Acq On : 27 Jun 2023 17:54
 Operator : CG\JU
 Sample : 03357-11
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-SCAN-03

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-BF062623.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF133996.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.146	4	10	17	rVB	1070339	1344008	55.84%	6.887%
2	2.257	25	29	40	rVB	35706	54191	2.25%	0.278%
3	5.087	506	510	518	rBV	143723	168272	6.99%	0.862%
4	5.487	573	578	581	rBV	2071672	1850718	76.89%	9.483%
5	6.487	743	748	751	rBV	2052013	1940482	80.62%	9.943%
6	6.851	806	810	818	rBV	769256	609066	25.31%	3.121%
7	7.410	901	905	908	rBV	1401270	1263814	52.51%	6.476%
8	8.128	1024	1027	1035	rBV	934878	757497	31.47%	3.881%
9	9.204	1207	1210	1214	rBV	2714755	2406818	100.00%	12.332%
10	9.698	1290	1294	1298	rBV	192620	161197	6.70%	0.826%
11	9.886	1323	1326	1331	rVB	990433	774560	32.18%	3.969%
12	10.004	1340	1346	1348	rVB2	60877	56148	2.33%	0.288%
13	10.222	1380	1383	1387	rVB	122001	90268	3.75%	0.463%
14	10.604	1445	1448	1451	rBV	166265	129028	5.36%	0.661%
15	10.680	1457	1461	1467	rBV	1081760	956065	39.72%	4.899%
16	10.792	1477	1480	1487	rBV	365395	321934	13.38%	1.650%
17	10.939	1499	1505	1508	rBV3	60924	106028	4.41%	0.543%
18	11.275	1558	1562	1565	rBV	150514	181448	7.54%	0.930%
19	11.380	1577	1580	1584	rBV	811392	708310	29.43%	3.629%
20	11.563	1609	1611	1614	rBV3	63137	75331	3.13%	0.386%
21	11.627	1620	1622	1625	rBV3	62864	51553	2.14%	0.264%
22	11.910	1668	1670	1673	rBV2	116813	96571	4.01%	0.495%
23	11.980	1679	1682	1684	rBV3	154848	206186	8.57%	1.056%
24	12.063	1693	1696	1699	rBV3	168217	200922	8.35%	1.030%
25	12.092	1699	1701	1703	rVB	191975	171211	7.11%	0.877%
26	12.410	1753	1755	1758	rBV3	216931	204536	8.50%	1.048%
27	12.592	1783	1786	1788	rBV3	311881	362179	15.05%	1.856%
28	12.986	1849	1853	1856	rBV	1845468	2205272	91.63%	11.300%
29	16.321	2417	2420	2435	rVB2	901090	2062634	85.70%	10.569%

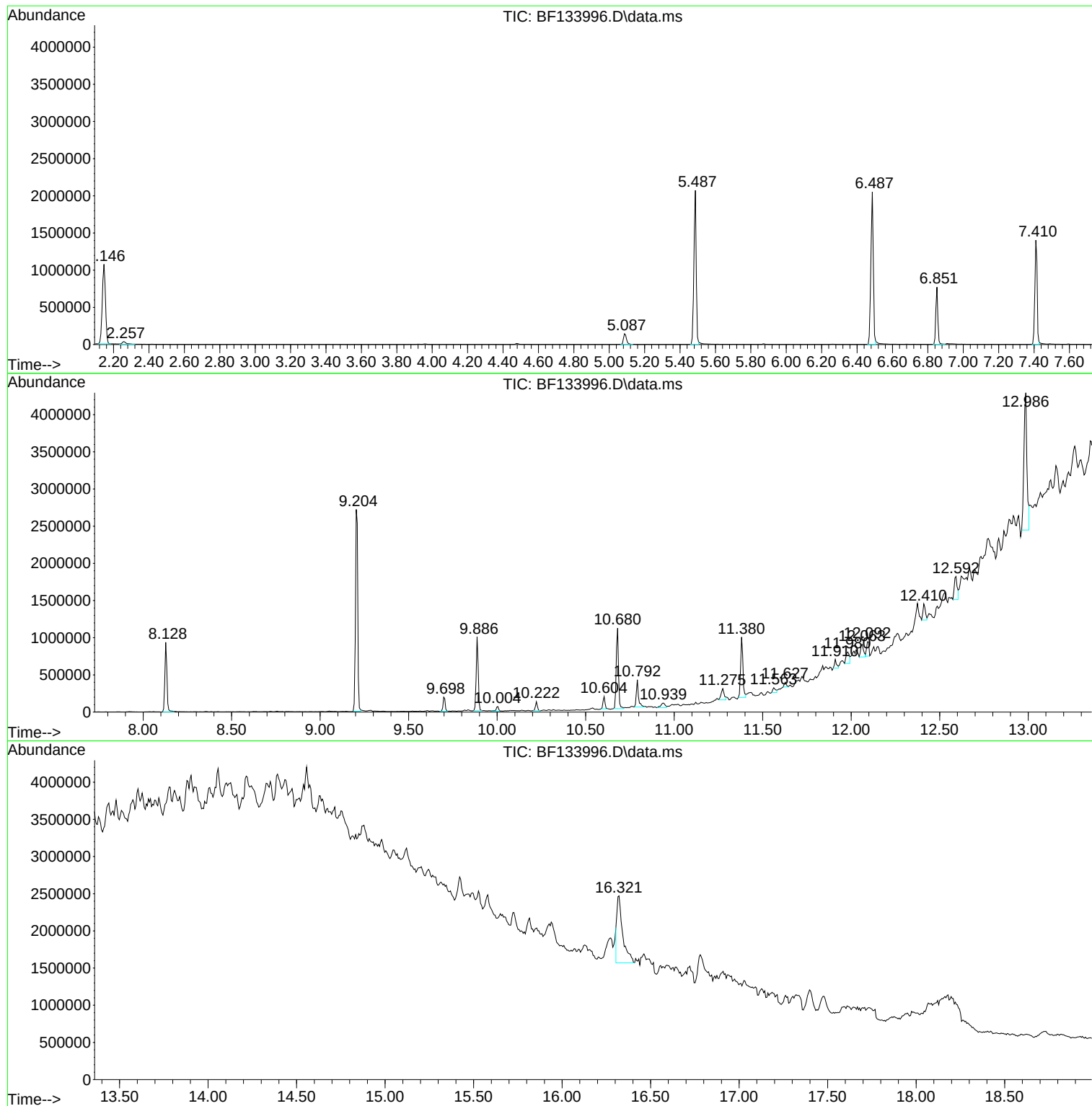
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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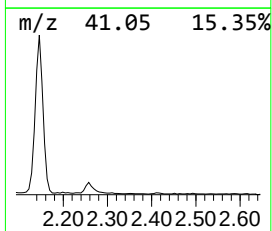
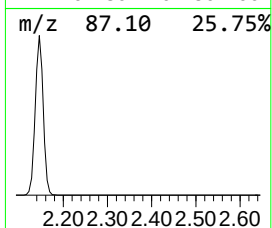
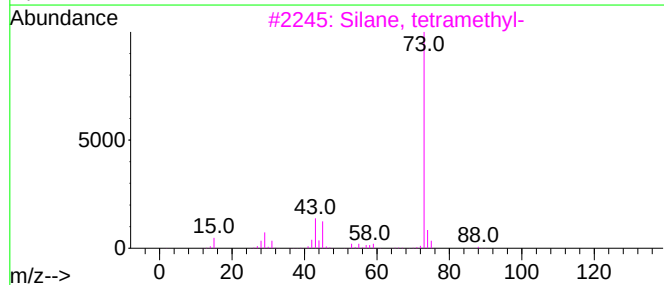
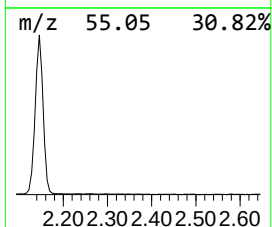
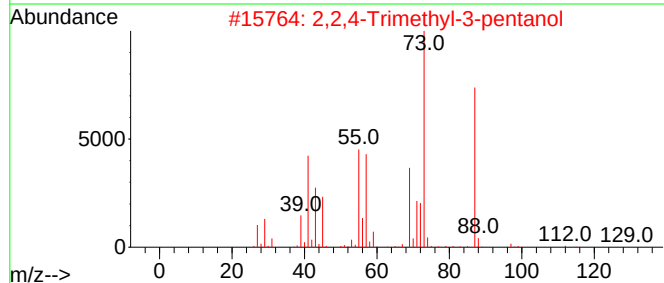
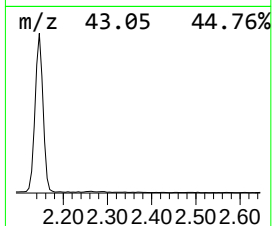
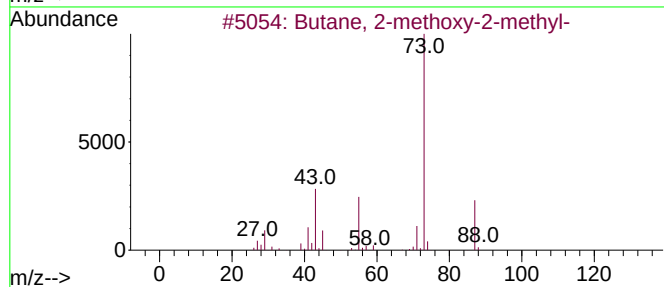
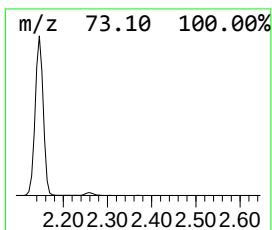
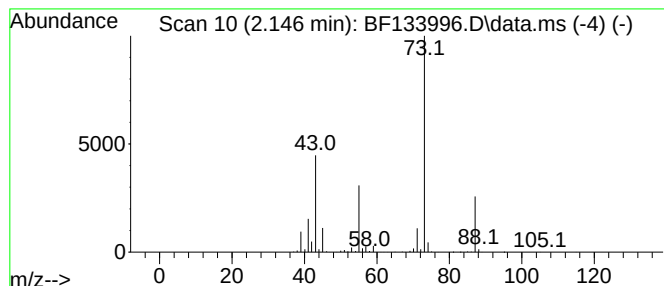
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.146	44.13 ng	1344010	1,4-Dichlorobenzene-d4	6.851

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
4			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	23
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23



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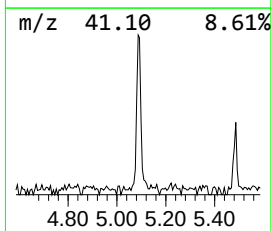
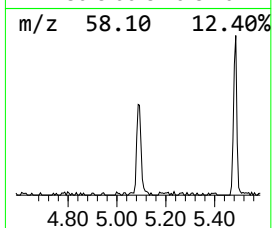
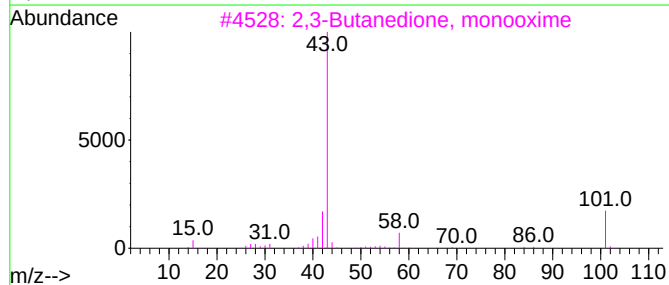
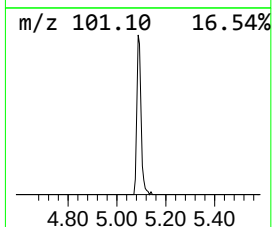
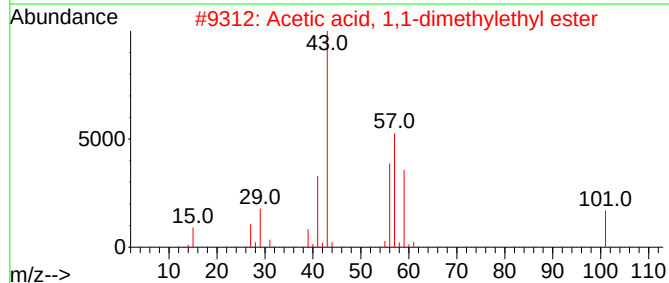
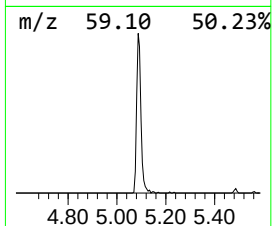
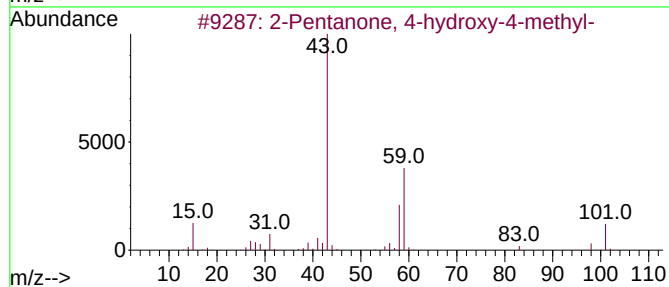
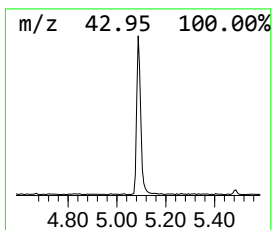
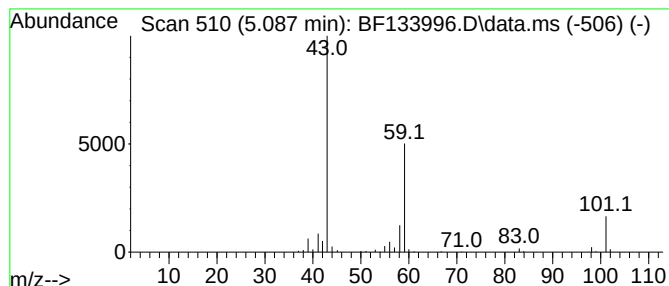
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.087	5.53 ng	168272	1,4-Dichlorobenzene-d4	6.851

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
5		Propanenitrile, 3-ethoxy-	99	C5H9NO	002141-62-0	9



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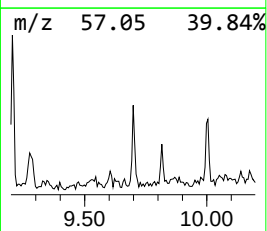
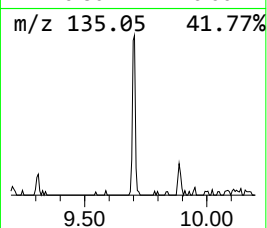
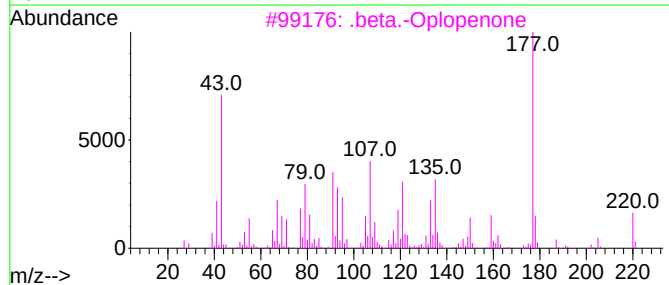
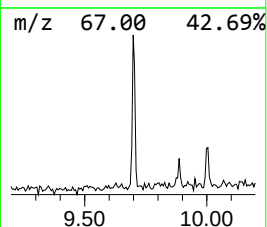
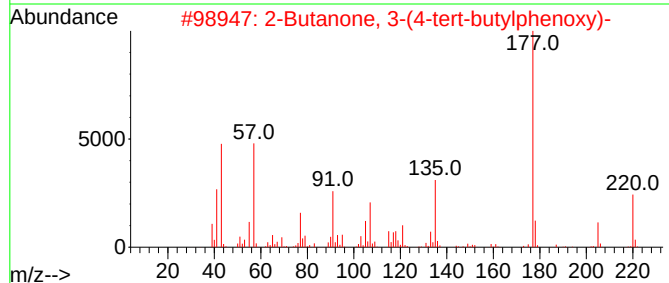
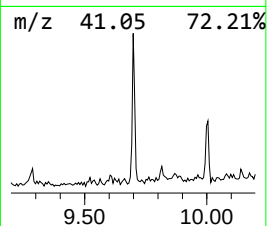
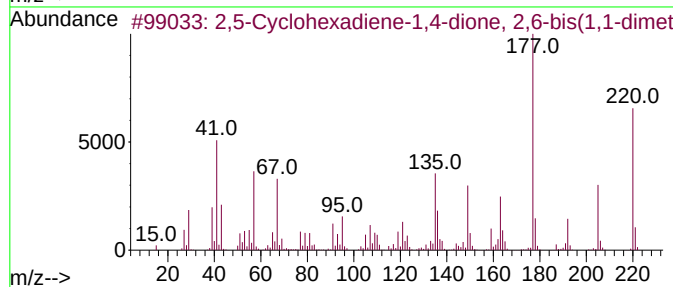
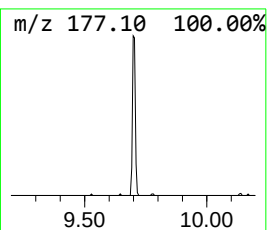
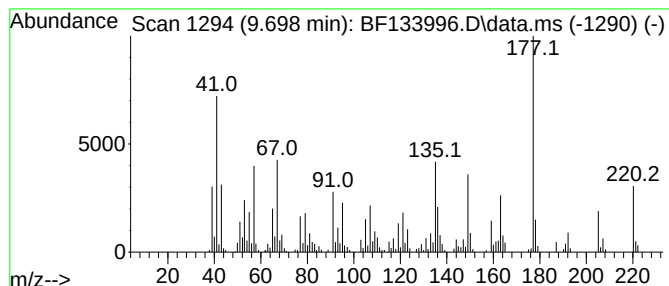
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BNA_F
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2,5-Cyclohexadiene-1,4-dion... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
9.698	4.16 ng	161197	Acenaphthene-d10		9.886	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2,5-Cyclohexadiene-1,4-dione, 2,...		220	C14H20O2	000719-22-2	98
2	2-Butanone, 3-(4-tert-butylpheno...		220	C14H20O2	160875-28-5	78
3	.beta.-Oplopenone		220	C15H24O	028305-60-4	46
4	2H-1-benzopyran-6-ol, 3,4-dihydr...		220	C14H20O2	1000396-25-4	42
5	1,3,5,7-Tetramethyl-adamantane		192	C14H24	001687-36-1	38



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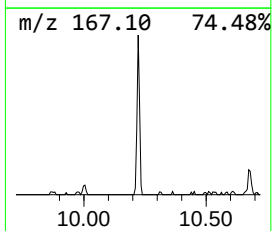
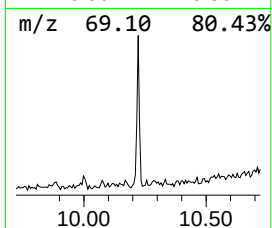
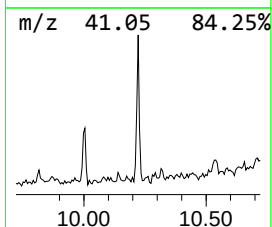
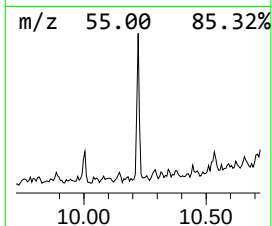
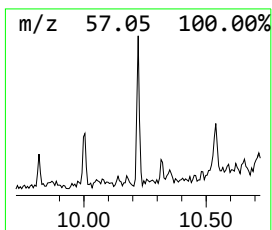
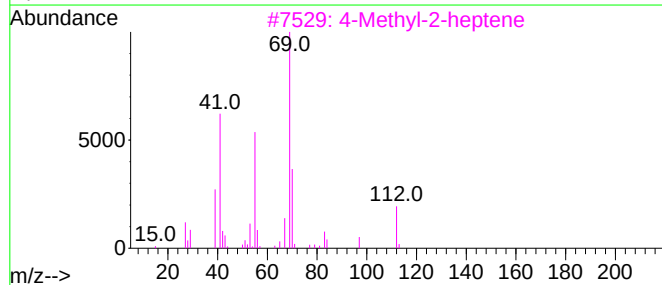
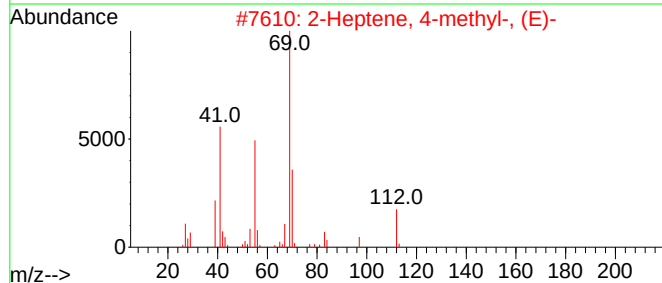
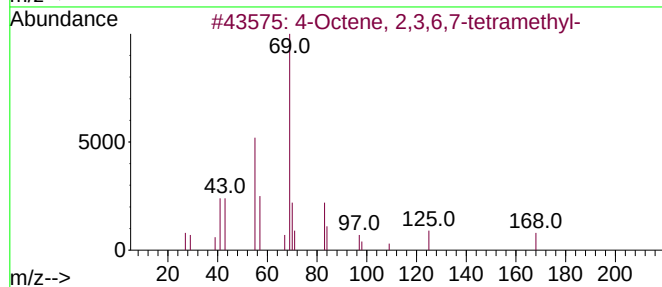
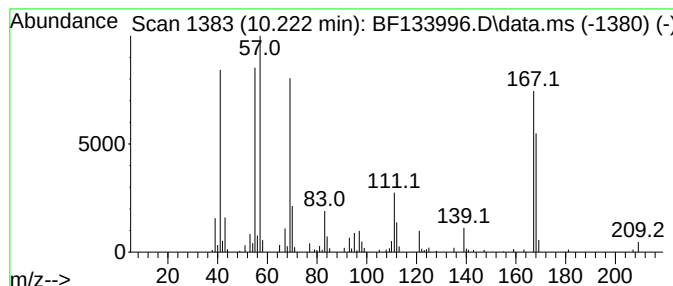
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 unknown10.222 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.222	2.33 ng	90268	Acenaphthene-d10	9.886	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Octene, 2,3,6,7-tetramethyl-	168	C12H24	063830-66-0	46
2	2-Heptene, 4-methyl-, (E)-	112	C8H16	066225-17-0	30
3	4-Methyl-2-heptene	112	C8H16	003404-56-6	30
4	1-Pentene, 3,3-dimethyl-	98	C7H14	003404-73-7	27
5	3-Heptene, 2-methyl-, (E)-	112	C8H16	000692-96-6	27



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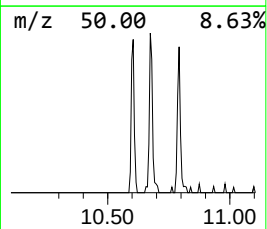
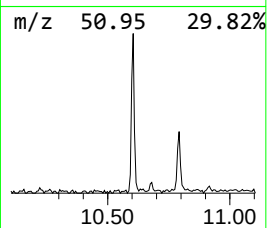
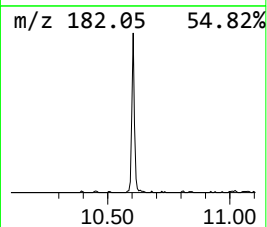
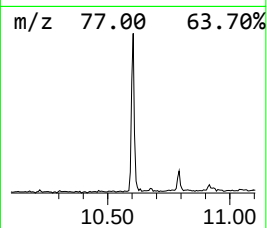
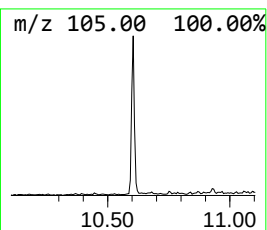
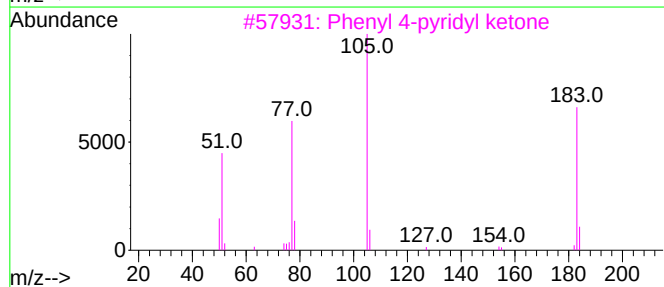
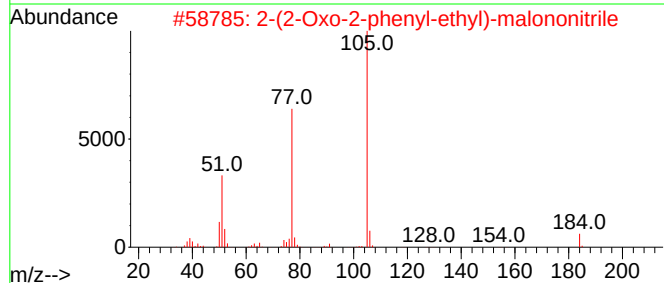
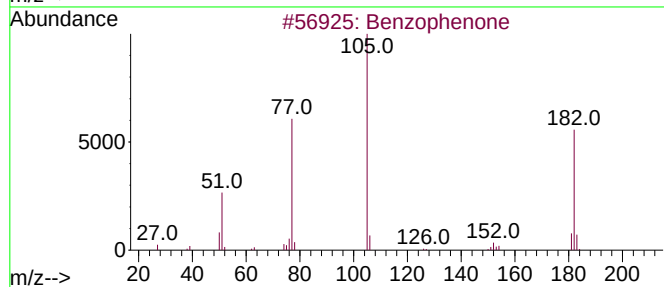
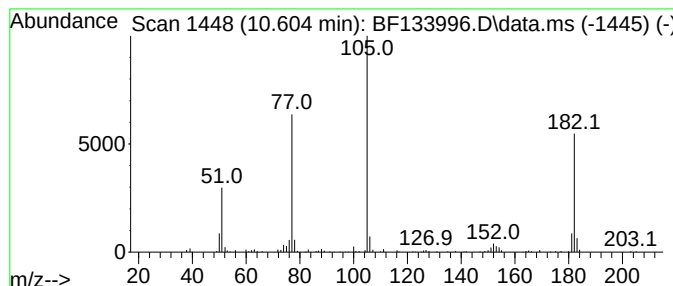
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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Peak Number 5 Benzophenone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.604	3.33 ng	129028	Acenaphthene-d10	9.886

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			2-(2-Oxo-2-phenyl-ethyl)-malonon...	184	C11H8N2O	1000296-76-9	53
3			Phenyl 4-pyridyl ketone	183	C12H9NO	014548-46-0	50
4			Azobenzene, (E)-	182	C12H10N2	017082-12-1	50
5			Methyl benzilate	242	C15H14O3	000076-89-1	47



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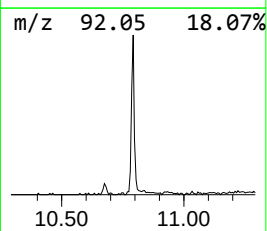
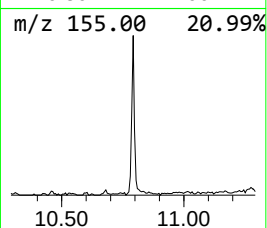
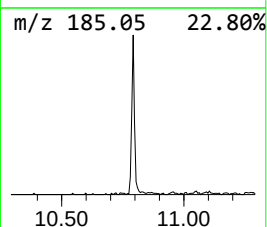
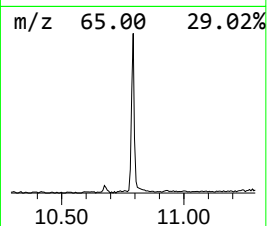
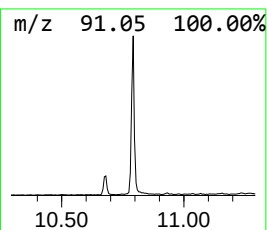
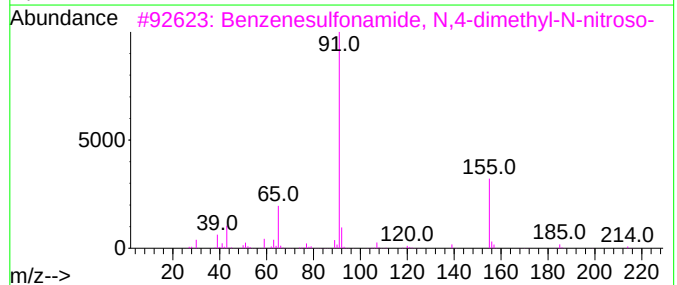
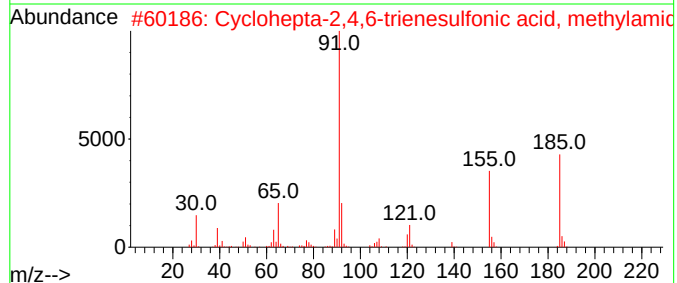
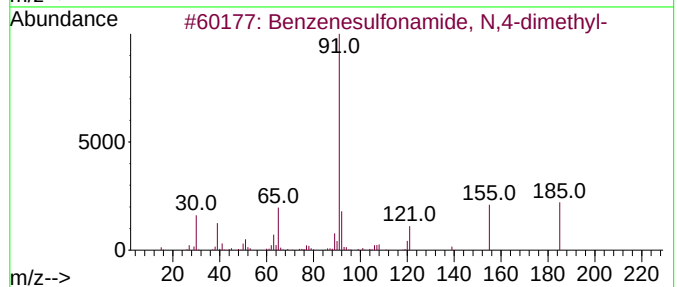
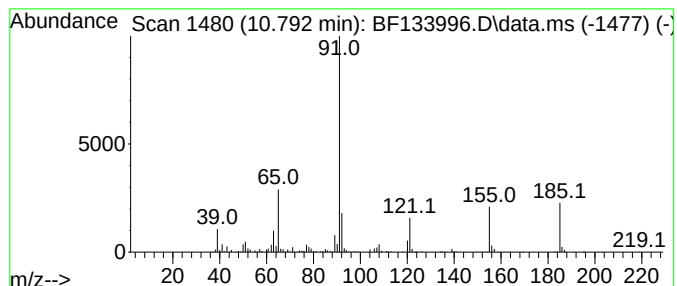
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzenesulfonamide, N,4-dim... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.792	9.09 ng	321934	Phenanthrene-d10	11.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N,4-dimethyl-	185	C8H11NO2S	000640-61-9	95
2			Cyclohepta-2,4,6-trienesulfonic ...	185	C8H11NO2S	1000187-28-7	58
3			Benzenesulfonamide, N,4-dimethyl...	214	C8H10N2O3S	000080-11-5	50
4			Ethyl(E/Z)-.alpha.-cyano-4-nitro...	416	C19H16N2O7S	1000287-26-8	50
5			Benzeneacetaldehyde	120	C8H8O	000122-78-1	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062723\
Data File : BF133996.D
Acq On : 27 Jun 2023 17:54
Operator : CG\JU
Sample : 03357-11
Misc :
ALS Vial : 15 Sample Multiplier: 1

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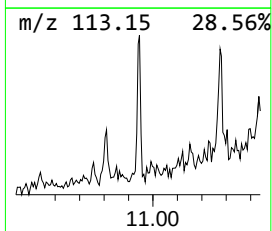
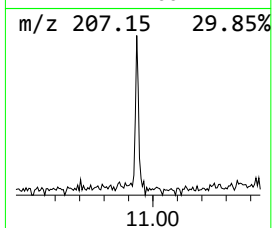
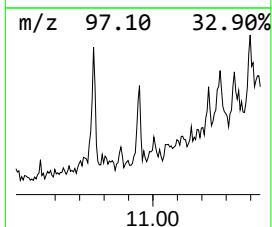
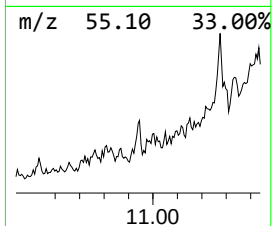
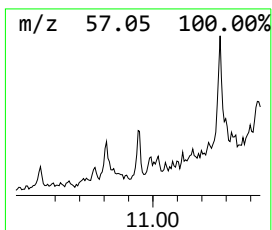
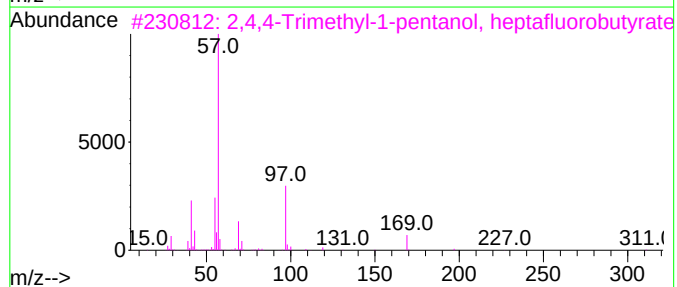
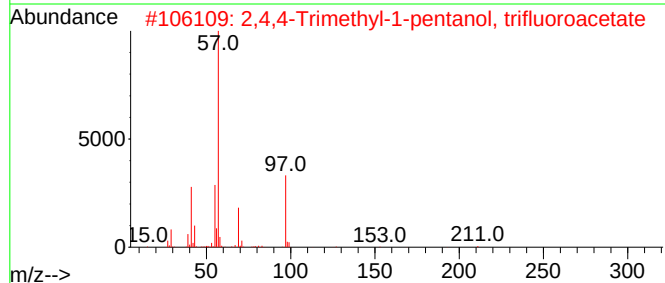
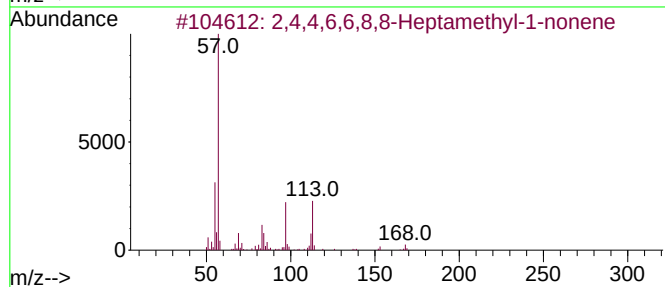
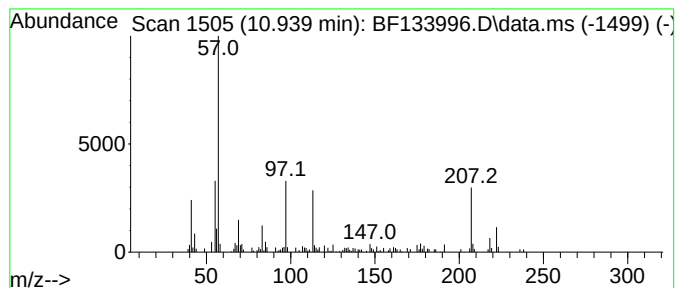
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown10.939 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.939	2.99 ng	106028	Phenanthrene-d10	11.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	37
2			2,4,4-Trimethyl-1-pentanol, trif...	226	C10H17F3O2	1000365-19-5	22
3			2,4,4-Trimethyl-1-pentanol, hept...	326	C12H17F7O2	1010365-01-4	22
4			4-Methylnonanoic acid	172	C10H20O2	045019-28-1	16
5			Sulfurous acid, 2-ethylhexyl oct...	446	C26H54O3S	1000309-20-1	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062723\
Data File : BF133996.D
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ClientSampleId :
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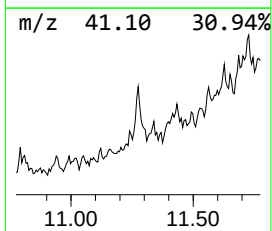
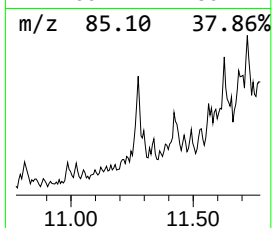
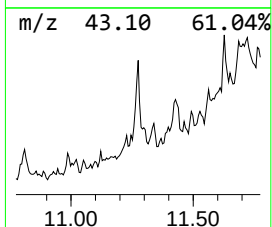
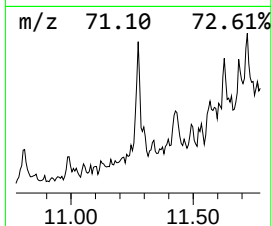
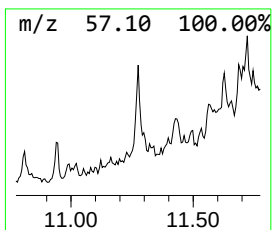
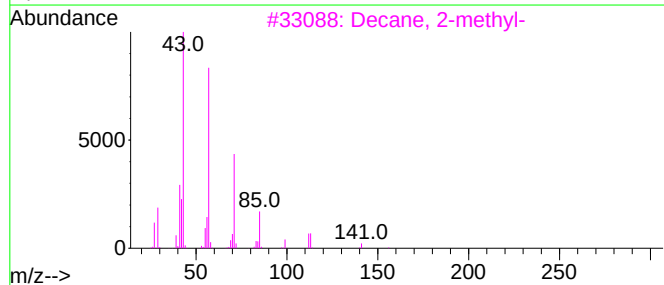
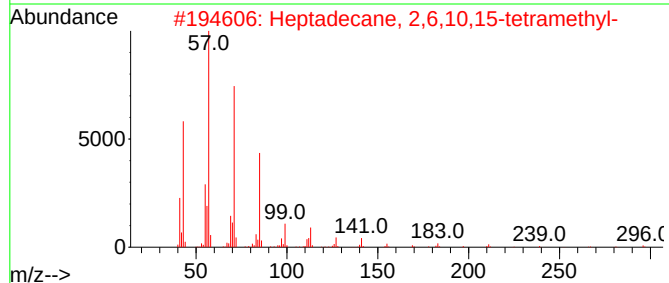
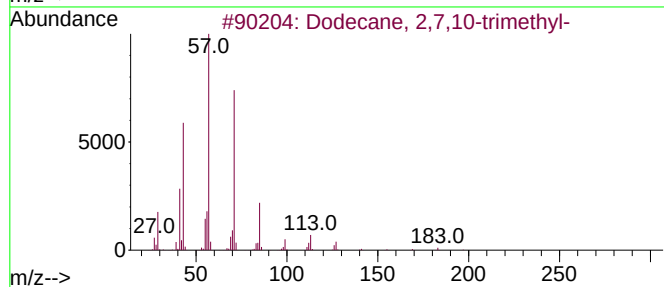
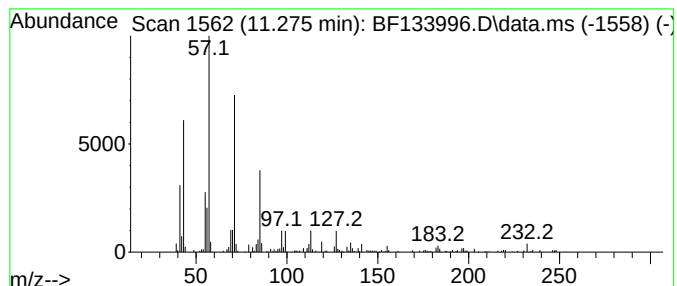
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Dodecane, 2,7,10-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.275	5.12 ng	181448	Phenanthrene-d10	11.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	80
2			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	80
3			Decane, 2-methyl-	156	C11H24	006975-98-0	74
4			Heptadecane	240	C17H36	000629-78-7	72
5			Tridecane, 4-methyl-	198	C14H30	026730-12-1	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062723\
Data File : BF133996.D
Acq On : 27 Jun 2023 17:54
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Misc :
ALS Vial : 15 Sample Multiplier: 1

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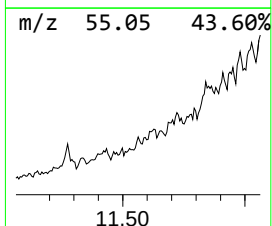
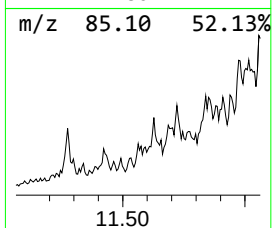
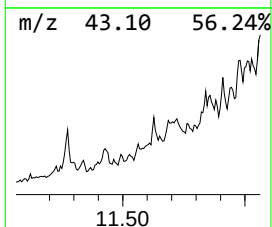
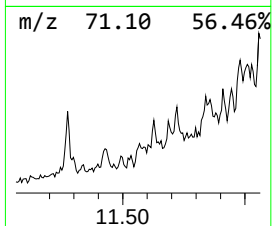
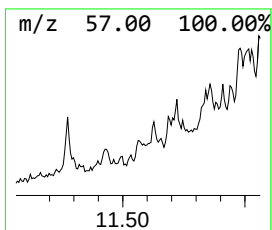
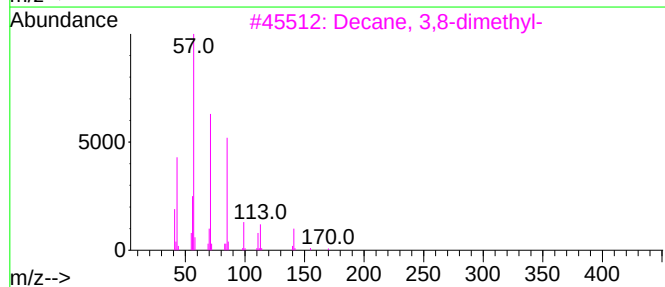
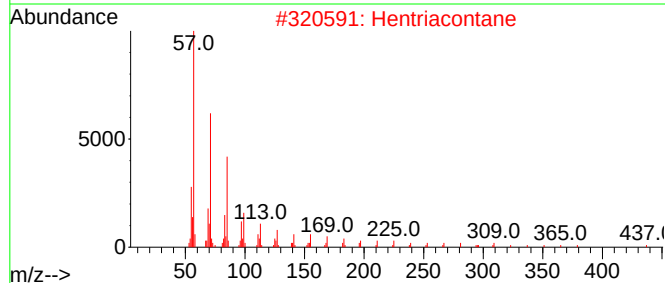
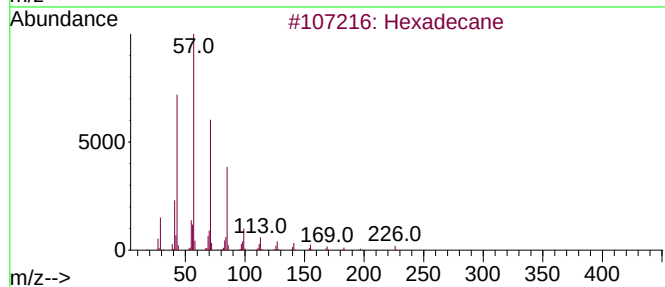
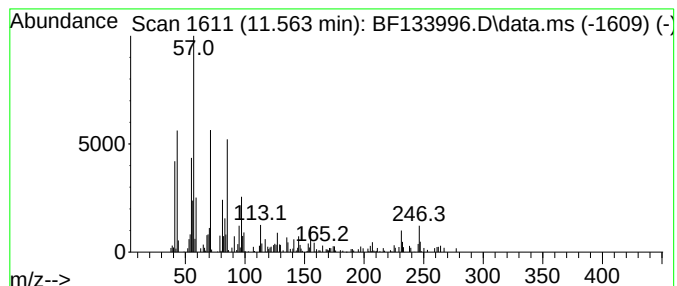
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.563	2.13 ng	75331	Phenanthrene-d10	11.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	50
2			Hentriacontane	437	C31H64	000630-04-6	49
3			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	46
4			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	46
5			2-Methyltetracosane	352	C25H52	001560-78-7	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062723\
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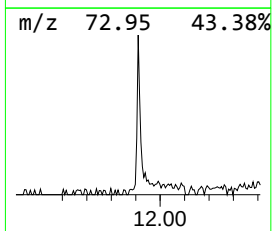
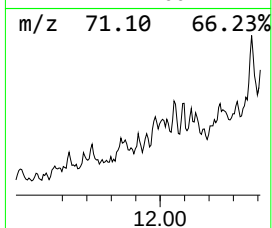
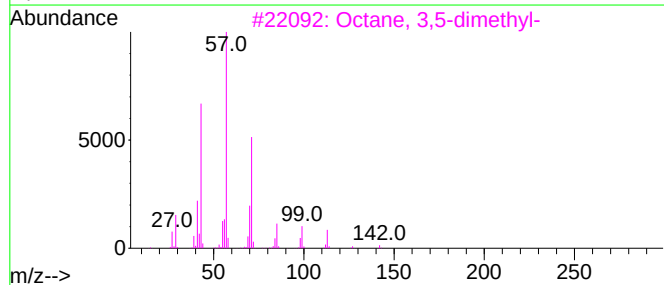
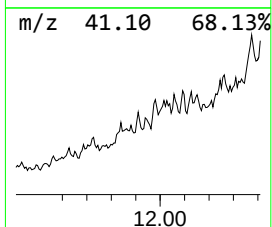
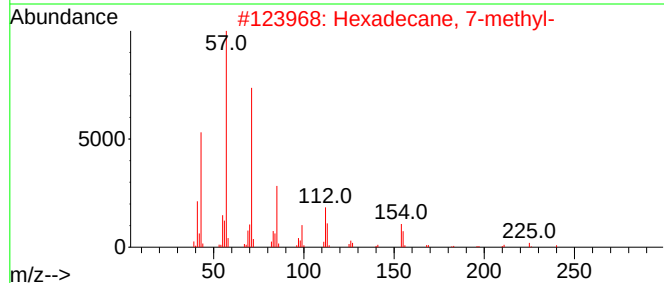
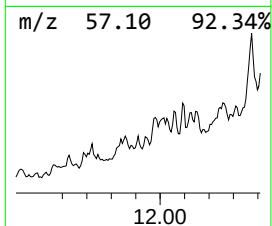
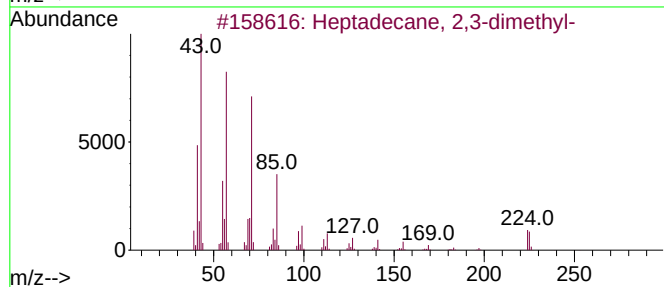
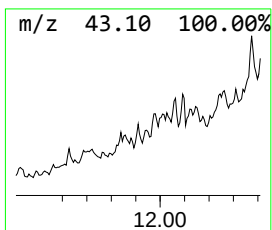
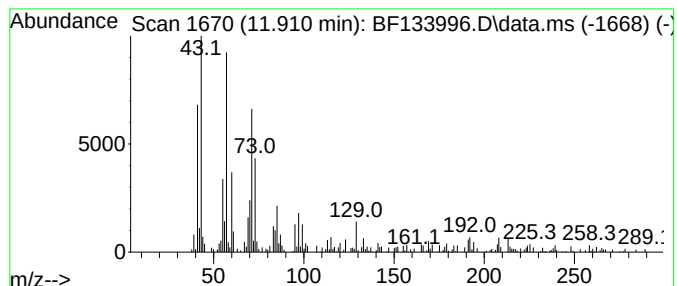
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown11.910 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.910	2.73 ng	96571	Phenanthrene-d10	11.380

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 2,3-dimethyl-	268	C19H40	061868-03-9	43
2		Hexadecane, 7-methyl-	240	C17H36	026730-20-1	43
3		Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	38
4		4-Methyl-3-heptanol, chlorodiflu...	242	C10H17ClF2O2	1000447-28-3	38
5		4-Methyl-3-heptanol, chlorodiflu...	242	C10H17ClF2O2	1000447-27-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062723\
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Sample : 03357-11
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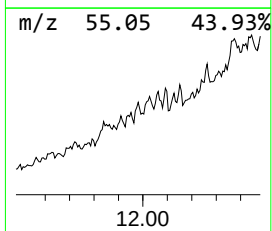
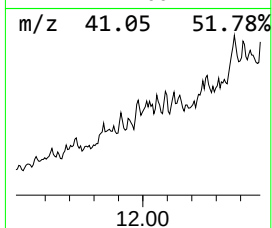
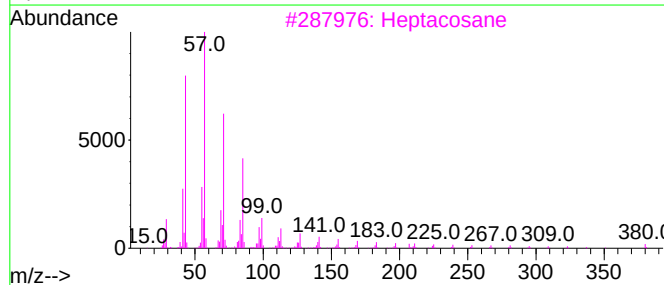
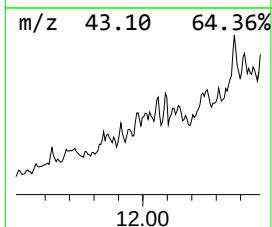
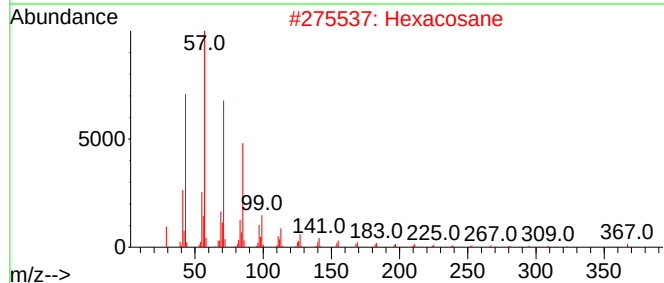
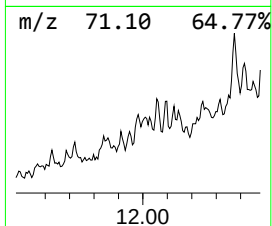
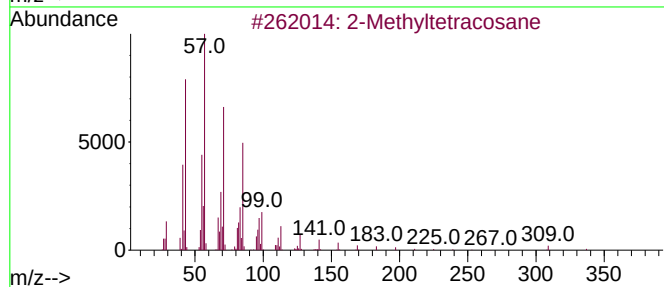
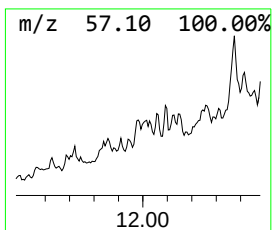
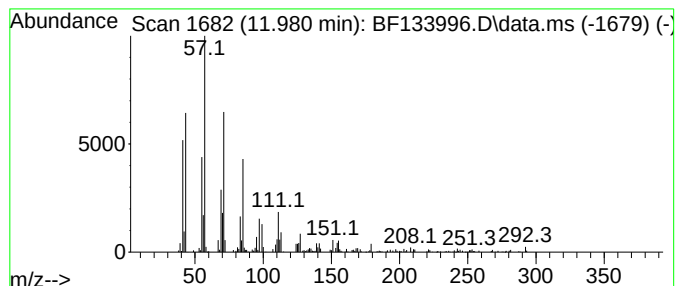
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2-Methyltetracosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.980	5.82 ng	206186	Phenanthrene-d10		11.380	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Methyltetracosane		352	C25H52	001560-78-7	87
2	Hexacosane		366	C26H54	000630-01-3	87
3	Heptacosane		380	C27H56	000593-49-7	86
4	Tetratetracontane		619	C44H90	007098-22-8	86
5	Eicosane, 7-hexyl-		366	C26H54	055333-99-8	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062723\
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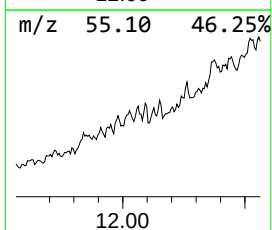
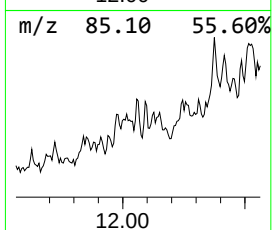
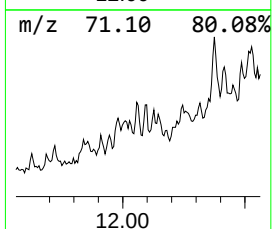
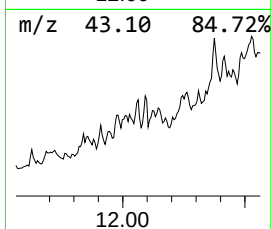
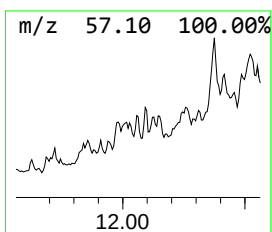
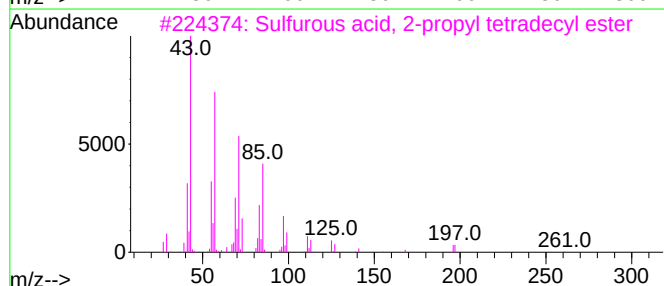
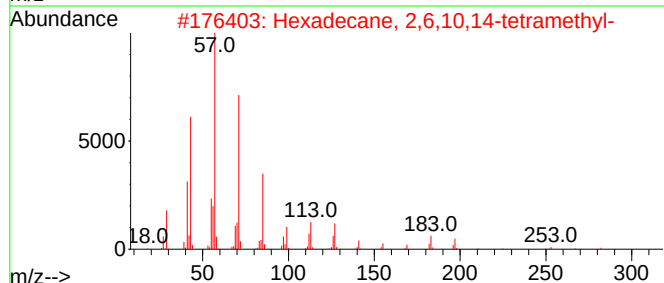
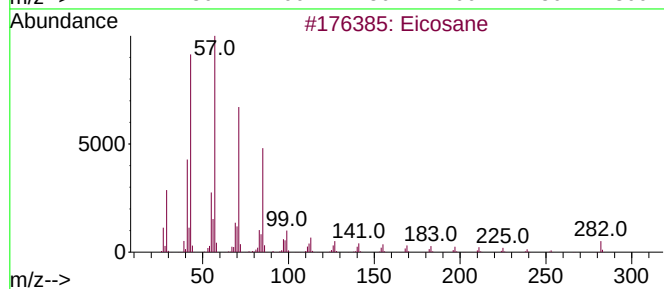
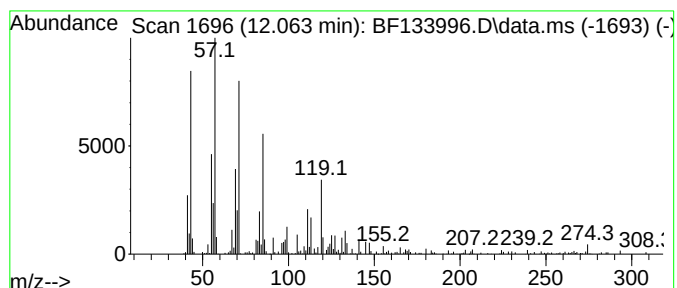
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.063	5.67 ng	200922	Phenanthrene-d10	11.380

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	64
2		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	55
3		Sulfurous acid, 2-propyl tetradecyl ester	320	C17H36O3S	1010309-12-5	53
4		Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	53
5		Nonyl tetradecyl ether	340	C23H48O	1000406-37-6	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062723\
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Misc :
ALS Vial : 15 Sample Multiplier: 1

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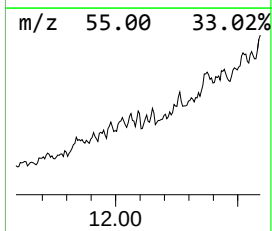
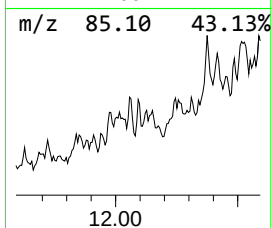
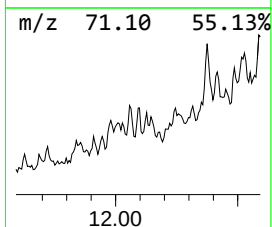
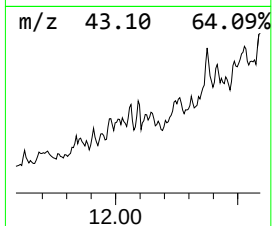
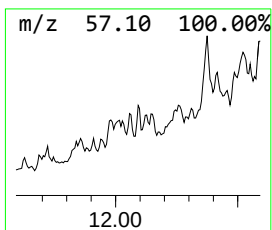
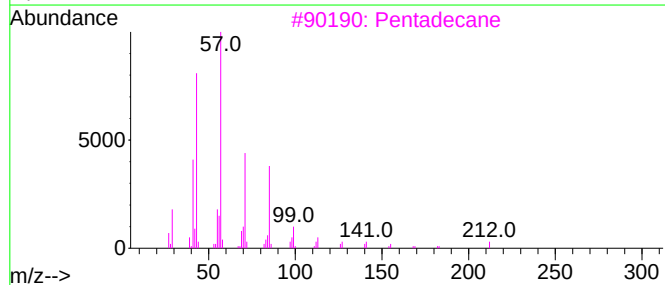
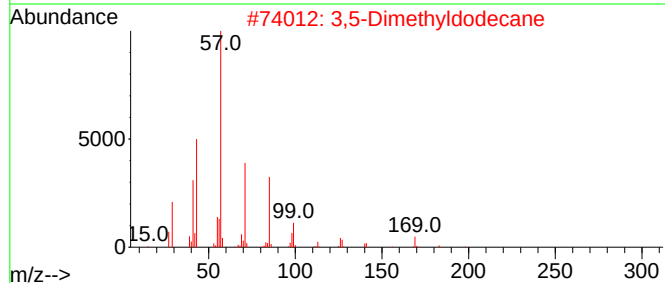
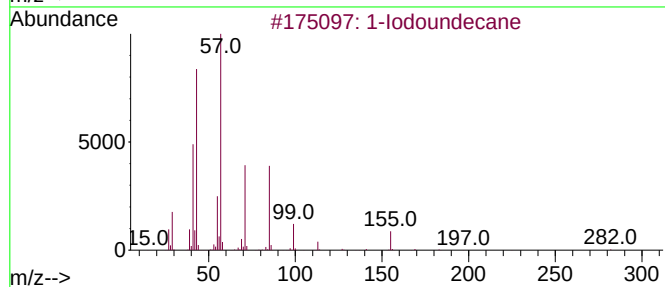
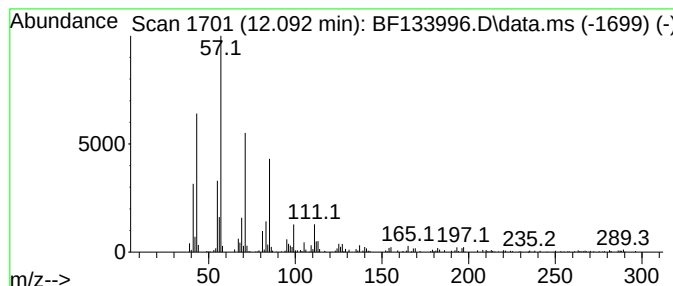
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1-Iodoundecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.092	4.83 ng	171211	Phenanthrene-d10	11.380

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Iodoundecane		282	C11H23I	004282-44-4	72
2	3,5-Dimethyldodecane		198	C14H30	107770-99-0	70
3	Pentadecane		212	C15H32	000629-62-9	64
4	Tetradecane, 1-iodo-		324	C14H29I	019218-94-1	58
5	Undecane, 5-methyl-		170	C12H26	001632-70-8	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062723\
Data File : BF133996.D
Acq On : 27 Jun 2023 17:54
Operator : CG\JU
Sample : 03357-11
Misc :
ALS Vial : 15 Sample Multiplier: 1

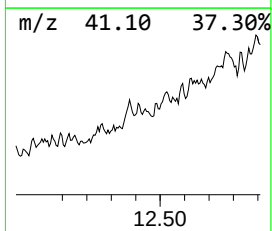
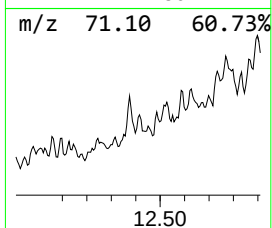
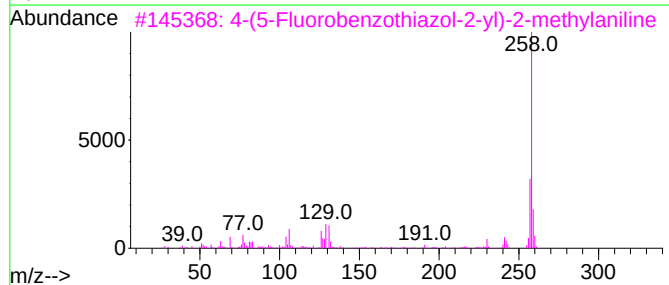
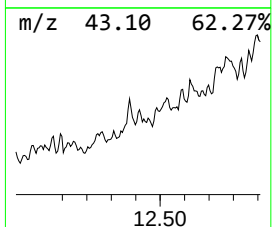
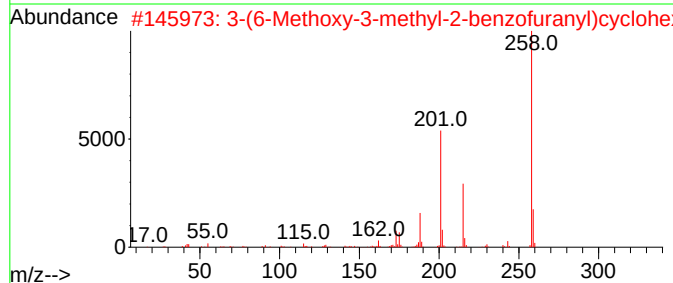
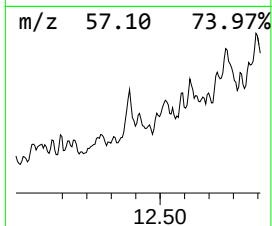
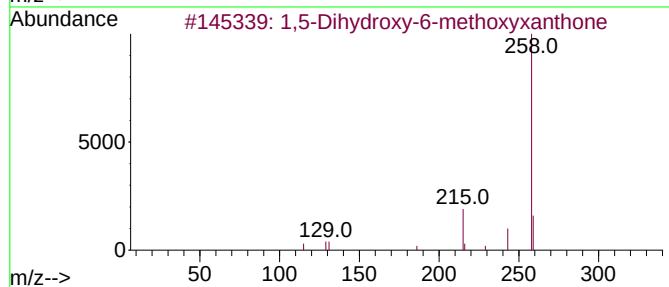
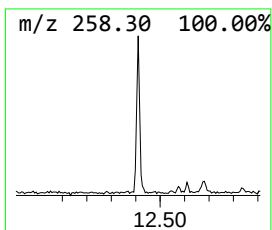
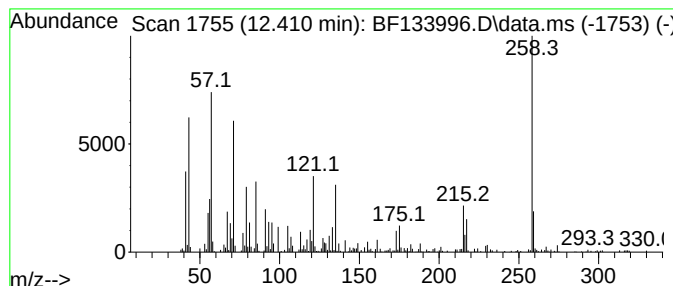
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ClientSampleId :
WC-SCAN-03

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1,5-Dihydroxy-6-methoxyxant... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.410	5.78 ng	204536	Phenanthrene-d10		11.380	
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,5-Dihydroxy-6-methoxyxanthone	258	C14H10O5	181307-35-7	53
2		3-(6-Methoxy-3-methyl-2-benzofur...	258	C16H18O3	1000239-79-6	38
3		4-(5-Fluorobenzothiazol-2-yl)-2-...	258	C14H11FN2S	260443-89-8	30
4		Benzene, 1-bromo-4-chlorodifluor...	256	C7H4BrClF2O	112556-13-5	25
5		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062723\
Data File : BF133996.D
Acq On : 27 Jun 2023 17:54
Operator : CG\JU
Sample : 03357-11
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-SCAN-03

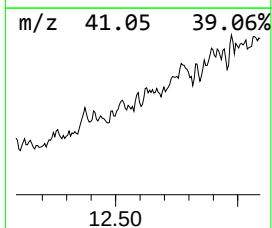
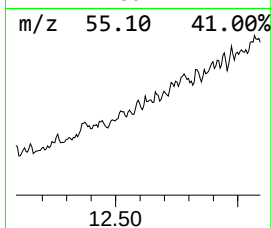
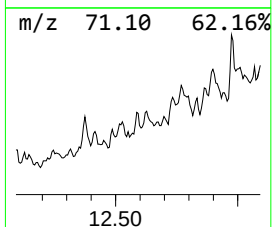
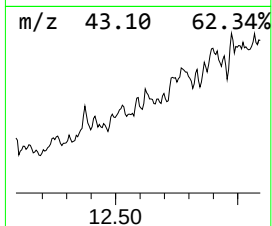
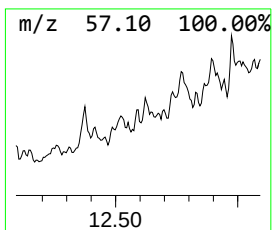
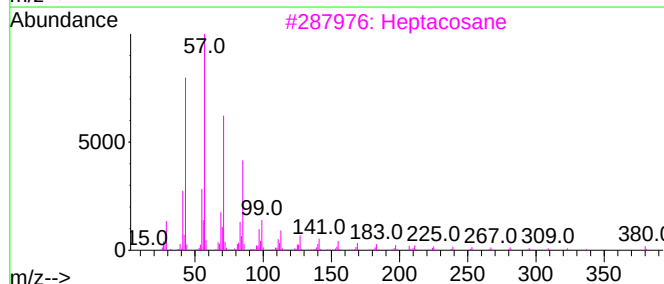
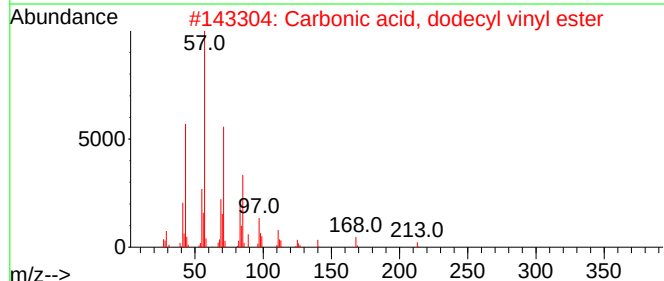
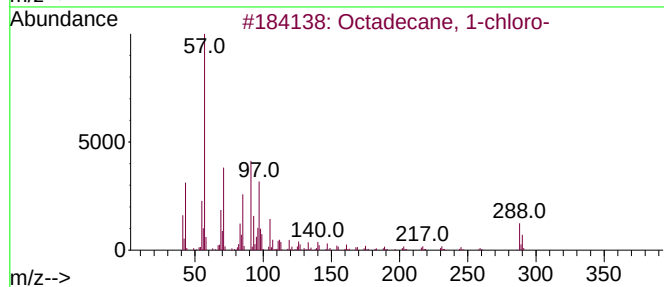
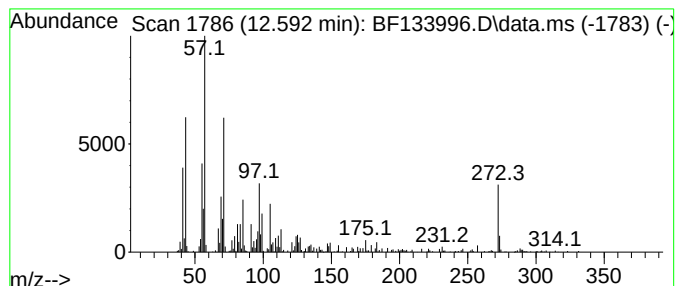
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown12.592 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.592	10.23 ng	362179	Phenanthrene-d10	11.380

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	49
2		Carbonic acid, dodecyl vinyl ester	256	C15H28O3	1000382-54-8	38
3		Heptacosane	380	C27H56	000593-49-7	38
4		Heptadecane	240	C17H36	000629-78-7	38
5		Sulfurous acid, 2-ethylhexyl hep...	432	C25H52O3S	1000309-20-0	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062723\
Data File : BF133996.D
Acq On : 27 Jun 2023 17:54
Operator : CG\JU
Sample : 03357-11
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-SCAN-03

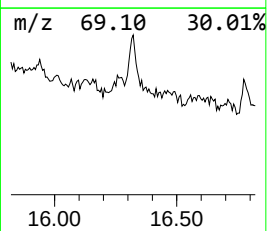
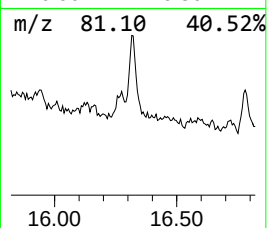
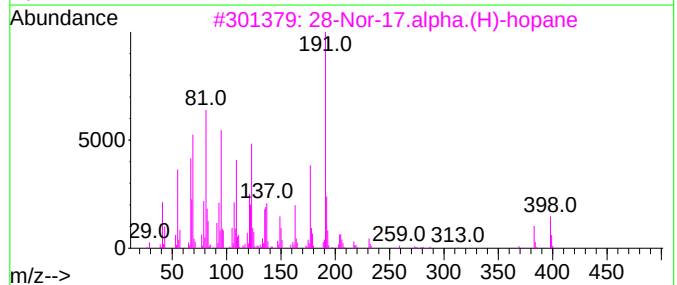
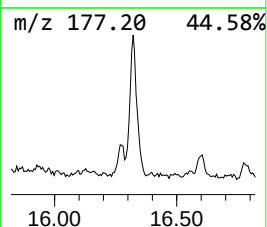
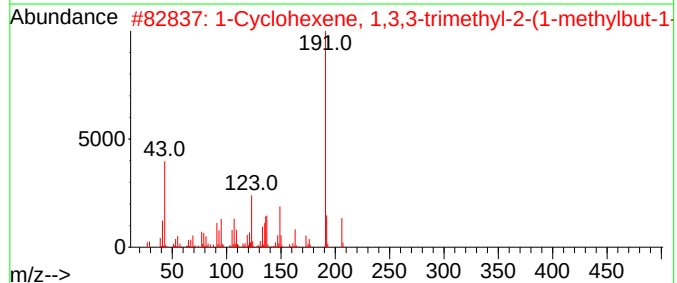
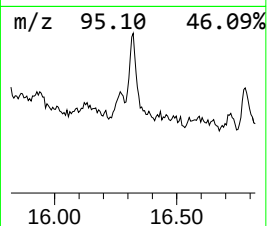
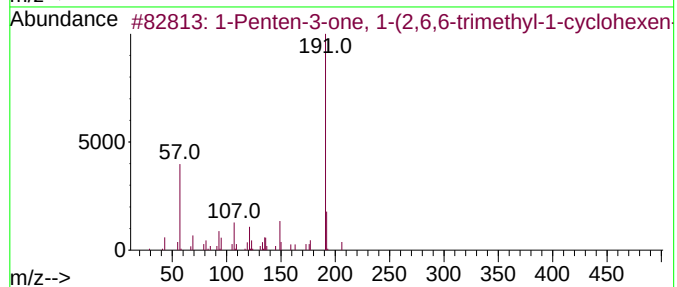
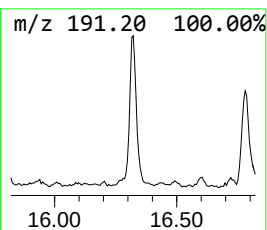
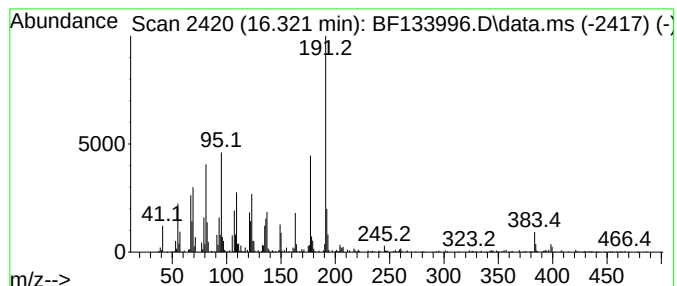
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown16.321 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.321	20.00 ng	2062630	Perylene-d12	15.574

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	45		
2	1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	37		
3	28-Nor-17.alpha.(H)-hopane	398	C29H50	053584-60-4	30		
4	1H-Pyrrolo[3,4-c]pyridine-1,3,(2...	191	C9H9N3O2	298688-32-1	27		
5	(1R,2S,5R)-2-Isopropyl-5-methylc...	374	C18H38O4Si2	1000499-39-0	27		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062723\
Data File : BF133996.D
Acq On : 27 Jun 2023 17:54
Operator : CG\JU
Sample : 03357-11
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-SCAN-03

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.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
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2-Pentanone, 4-...	5.087	5.5	ng	168272	1	6.851	609066	20.0
2,5-Cyclohexadi...	9.698	4.2	ng	161197	3	9.886	774560	20.0
unknown10.222	10.222	2.3	ng	90268	3	9.886	774560	20.0
Benzophenone	10.604	3.3	ng	129028	3	9.886	774560	20.0
Benzenesulfonam...	10.792	9.1	ng	321934	4	11.380	708310	20.0
unknown10.939	10.939	3.0	ng	106028	4	11.380	708310	20.0
Dodecane, 2,7,1...	11.275	5.1	ng	181448	4	11.380	708310	20.0
Hexadecane	11.563	2.1	ng	75331	4	11.380	708310	20.0
unknown11.910	11.910	2.7	ng	96571	4	11.380	708310	20.0
2-Methyltetraaco...	11.980	5.8	ng	206186	4	11.380	708310	20.0
Eicosane	12.063	5.7	ng	200922	4	11.380	708310	20.0
1-Iodoundecane	12.092	4.8	ng	171211	4	11.380	708310	20.0
1,5-Dihydroxy-6...	12.410	5.8	ng	204536	4	11.380	708310	20.0
unknown12.592	12.592	10.2	ng	362179	4	11.380	708310	20.0
unknown16.321	16.321	20.0	ng	2062630	6	15.574	2062630	20.0