

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071124\
 Data File : BF138537.D
 Acq On : 11 Jul 2024 22:23
 Operator : CG\JU
 Sample : P3181-01 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-02-070924

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-BF070924.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF138537.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.481	574	577	586	rVB	244301	207955	9.40%	1.184%
2	6.493	746	749	753	rBV	258533	206228	9.32%	1.174%
3	6.863	808	812	816	rBV	520782	399072	18.04%	2.271%
4	8.145	1027	1030	1032	rVB	583021	440137	19.90%	2.505%
5	8.681	1117	1121	1124	rBV4	105217	168619	7.62%	0.960%
6	8.916	1159	1161	1164	rBV3	133533	167351	7.57%	0.953%
7	8.963	1165	1169	1171	rVV4	156262	228588	10.33%	1.301%
8	9.022	1176	1179	1182	rVV5	160243	201075	9.09%	1.144%
9	9.051	1182	1184	1187	rBV3	120179	134802	6.09%	0.767%
10	9.139	1196	1199	1201	rBV3	240095	205149	9.27%	1.168%
11	9.163	1201	1203	1208	rBV2	604664	579442	26.20%	3.298%
12	9.216	1210	1212	1215	rVB3	289372	213887	9.67%	1.217%
13	9.257	1217	1219	1221	rBV2	165849	146452	6.62%	0.834%
14	9.304	1223	1227	1229	rBV2	627141	753892	34.08%	4.291%
15	9.534	1264	1266	1268	rBV2	353112	368765	16.67%	2.099%
16	9.569	1268	1272	1274	rVV2	1104858	1261005	57.01%	7.177%
17	9.610	1277	1279	1281	rBV	1015988	1151424	52.06%	6.554%
18	9.628	1281	1282	1284	rVB	1142757	688010	31.11%	3.916%
19	9.775	1303	1307	1310	rBV4	912899	1497285	67.69%	8.522%
20	9.816	1311	1314	1317	rVV	1400160	1435959	64.92%	8.173%
21	9.892	1323	1327	1328	rBV2	1907800	2211859	100.00%	12.589%
22	10.234	1382	1385	1386	rBV2	1051693	1103783	49.90%	6.282%
23	10.328	1398	1401	1404	rBV2	1380222	1752093	79.21%	9.973%
24	15.504	2278	2281	2291	rVB	375796	558684	25.26%	3.180%
25	16.198	2395	2399	2415	rVB2	317086	789713	35.70%	4.495%
26	16.633	2468	2473	2483	rVB	372568	698016	31.56%	3.973%

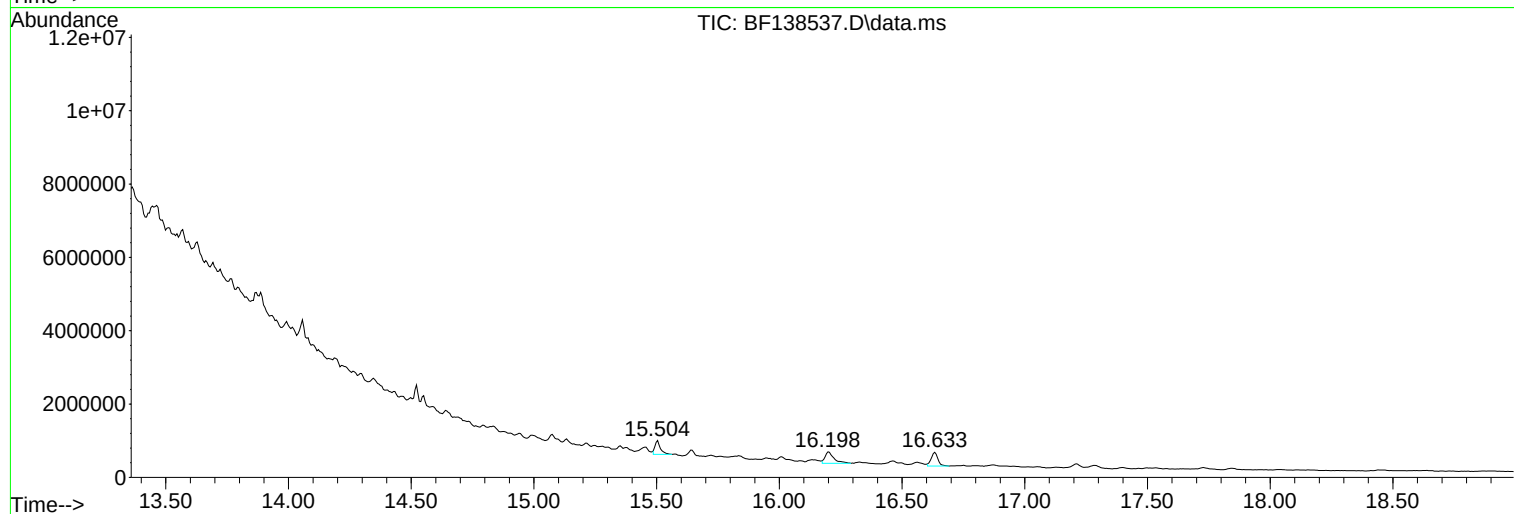
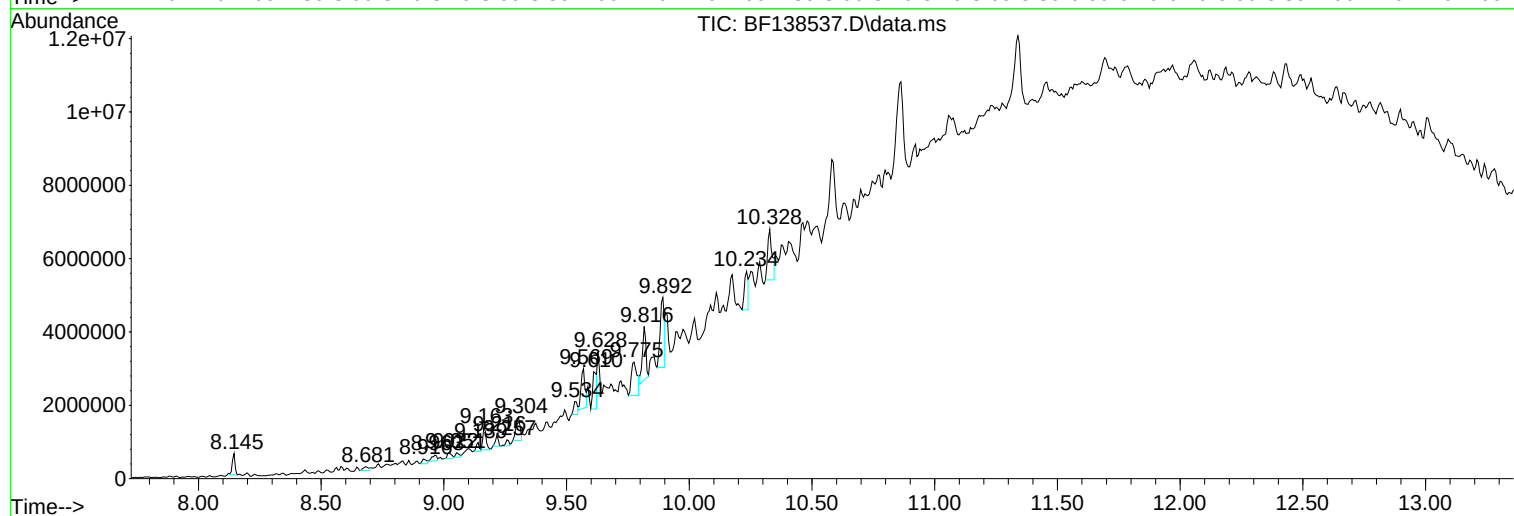
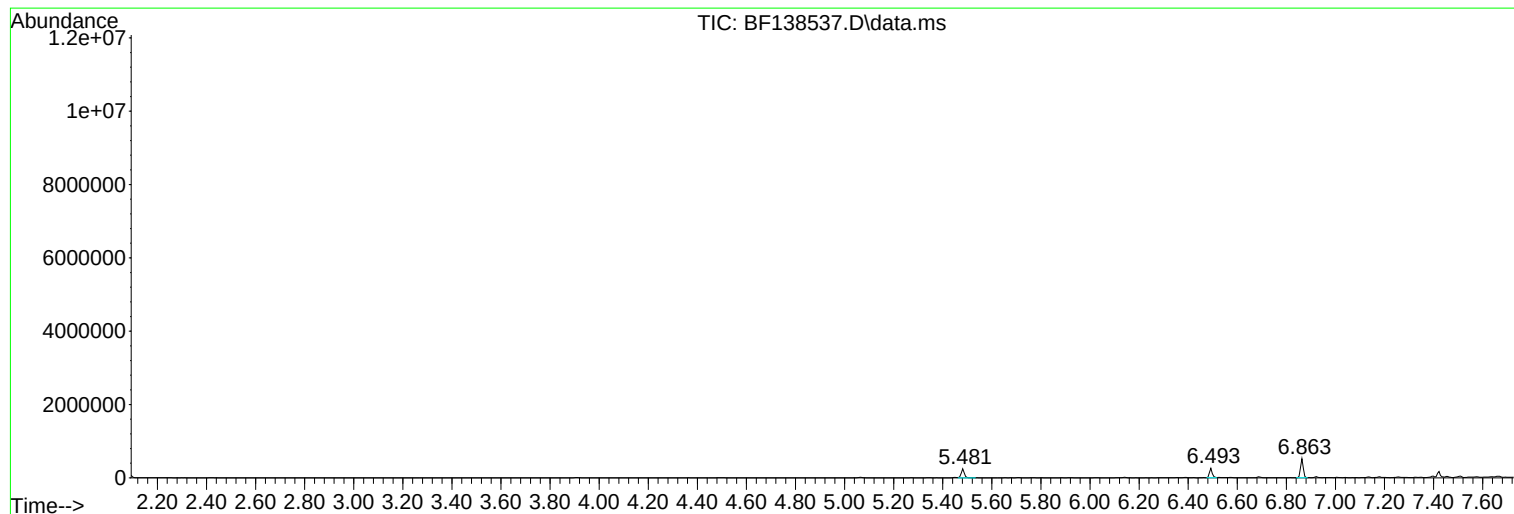
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Instrument :
BNA_F
ClientSampleId :
OR-02-070924

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF070924.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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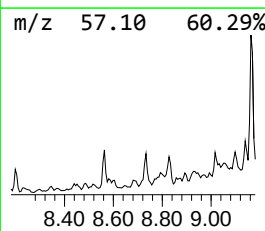
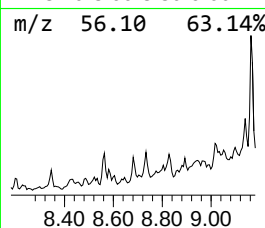
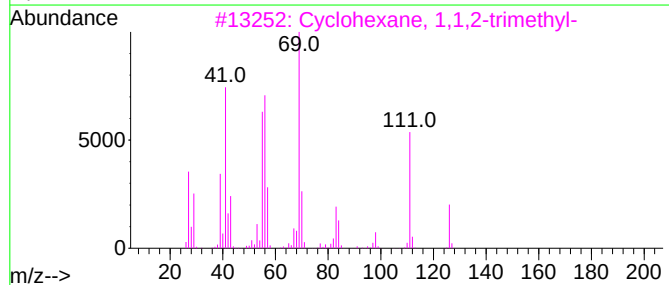
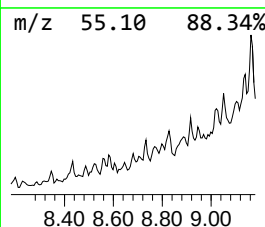
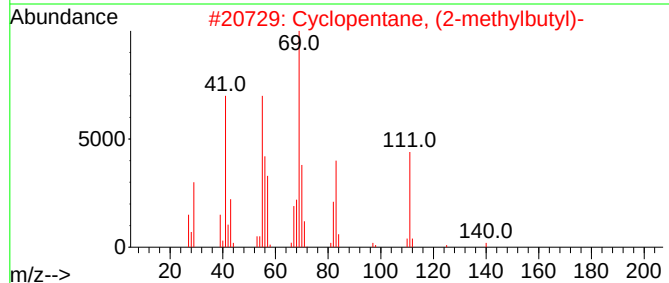
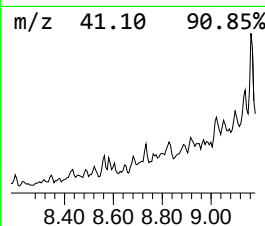
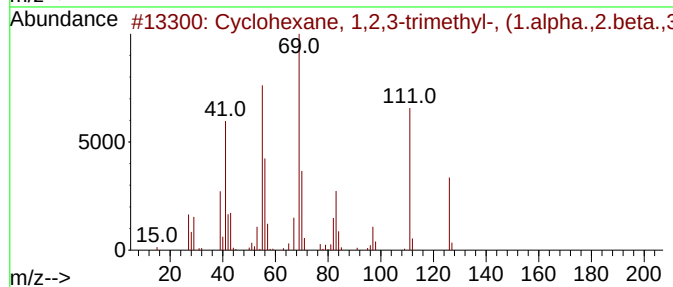
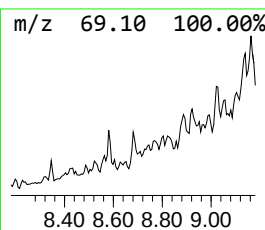
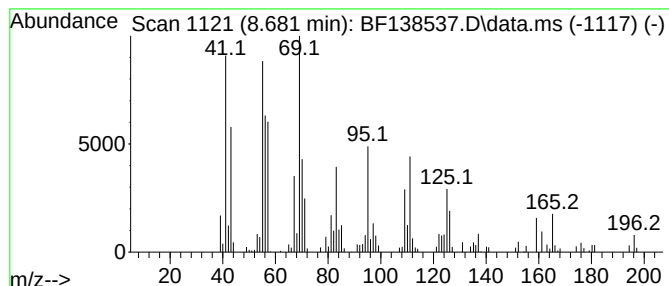
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Cyclohexane, 1,2,3-trimethyl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.681	7.66 ng	168619	Naphthalene-d8	8.145

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2,3-trimethyl-, (...)	126	C9H18	001678-81-5	70
2			Cyclopentane, (2-methylbutyl)-	140	C10H20	053366-38-4	62
3			Cyclohexane, 1,1,2-trimethyl-	126	C9H18	007094-26-0	60
4			Tridecane, 7-methylene-	196	C14H28	019780-80-4	53
5			3-Heptene, 2,6-dimethyl-	126	C9H18	002738-18-3	49



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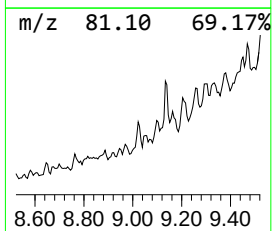
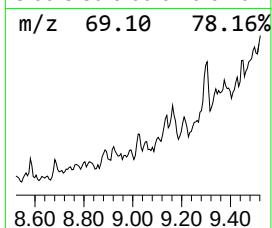
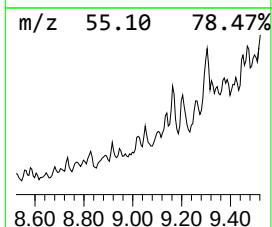
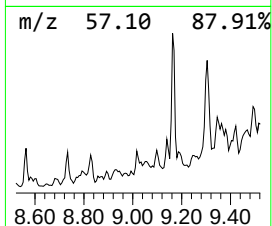
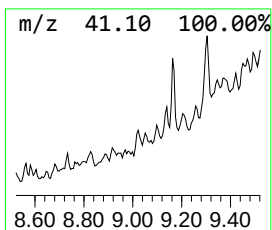
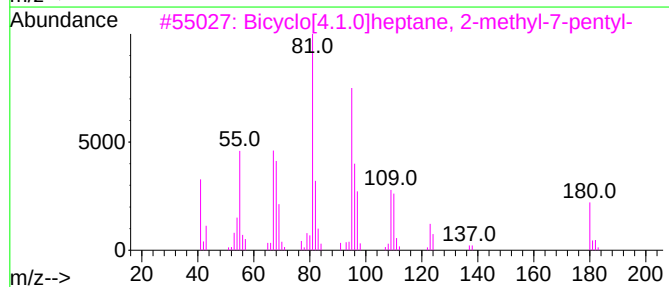
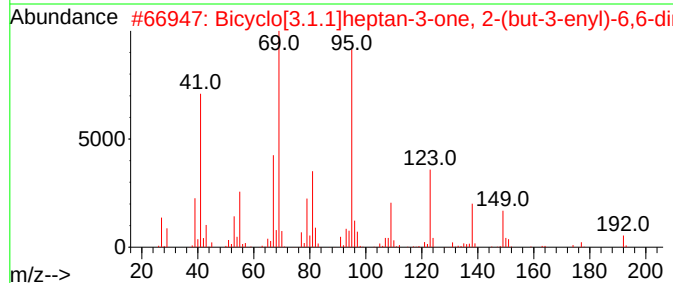
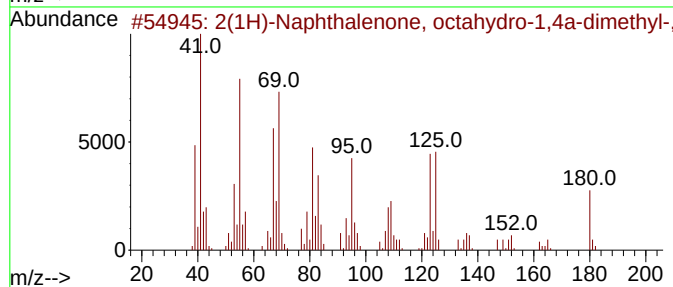
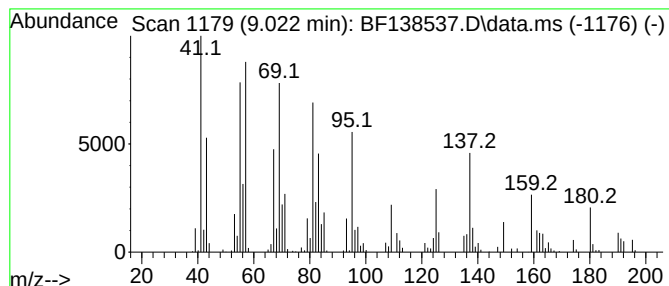
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown9.022 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.022	9.14 ng	201075	Naphthalene-d8	8.145

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Naphthalenone, octahydro-1...	180	C12H20O	022738-31-4	49		
2	Bicyclo[3.1.1]heptan-3-one, 2-(b...	192	C13H20O	1010163-96-1	46		
3	Bicyclo[4.1.0]heptane, 2-methyl-...	180	C13H24	055937-92-3	38		
4	Cyclopentane, 1,2-dibutyl-	182	C13H26	062199-52-4	35		
5	(7E,9E)-Dodecadienal	180	C12H20O	097475-10-0	35		



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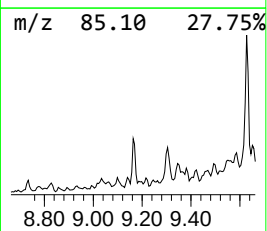
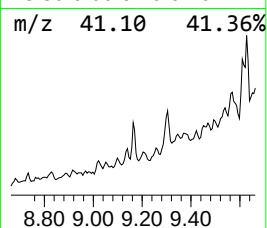
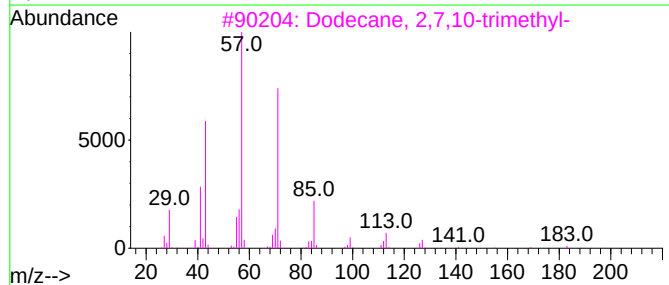
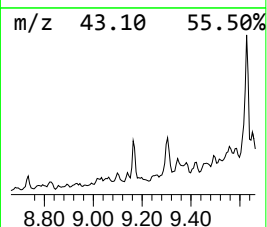
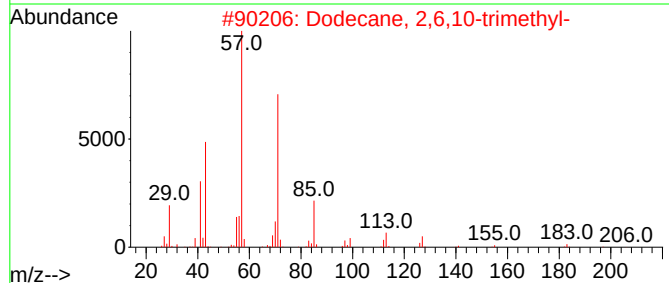
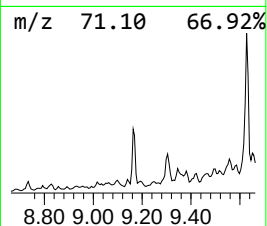
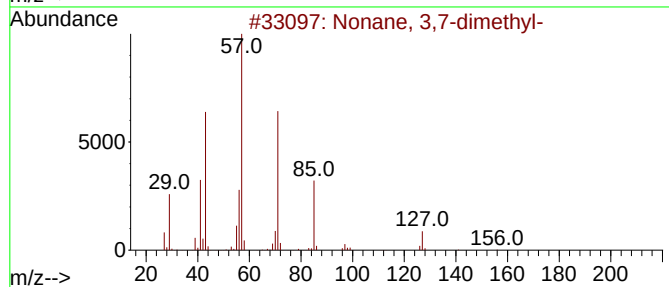
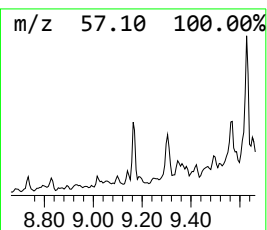
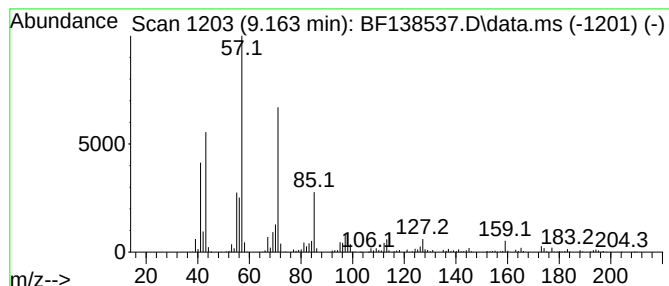
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Nonane, 3,7-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.	
9.163	5.24 ng	579442	Acenaphthene-d10		9.904	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 3,7-dimethyl-		156	C11H24	017302-32-8	81
2	Dodecane, 2,6,10-trimethyl-		212	C15H32	003891-98-3	80
3	Dodecane, 2,7,10-trimethyl-		212	C15H32	074645-98-0	64
4	Heneicosane		296	C21H44	000629-94-7	64
5	Heptacosane		380	C27H56	000593-49-7	53



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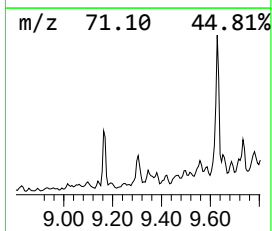
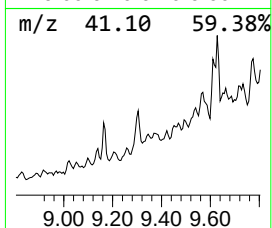
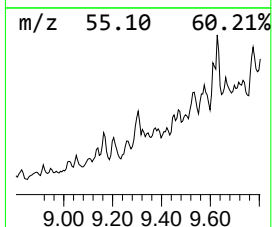
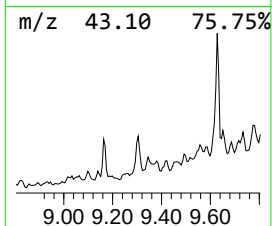
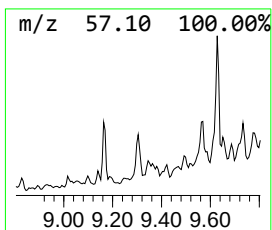
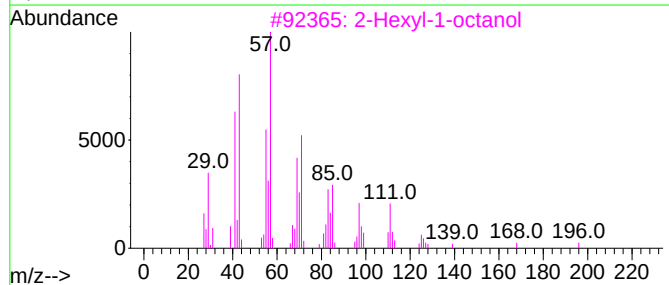
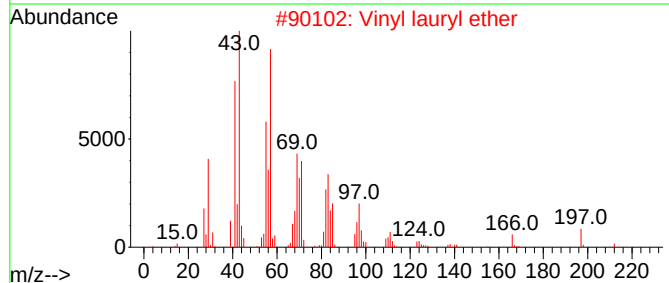
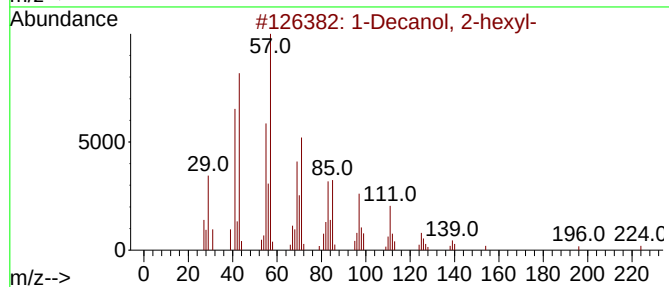
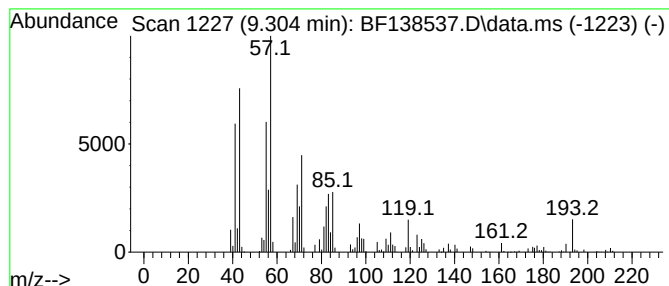
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 1-Decanol, 2-hexyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.304	6.82 ng	753892	Acenaphthene-d10	9.904	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	64
2	Vinyl lauryl ether	212	C14H28O	000765-14-0	52
3	2-Hexyl-1-octanol	214	C14H30O	019780-79-1	52
4	Carbonic acid, tetradecyl vinyl ...	284	C17H32O3	1000382-54-5	52
5	Cyclododecanol, 1-ethenyl-	210	C14H26O	006244-49-1	49



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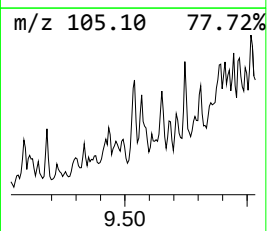
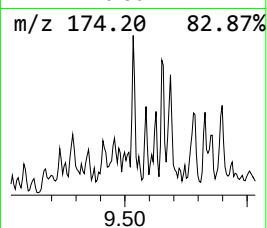
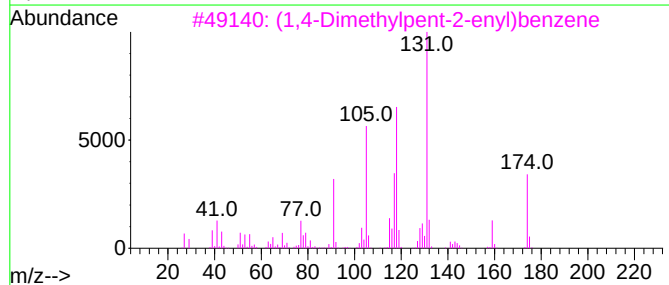
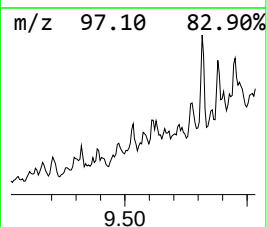
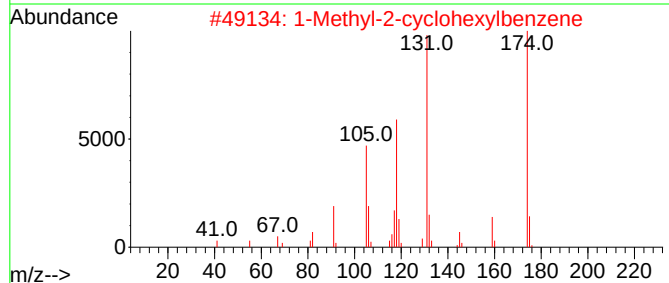
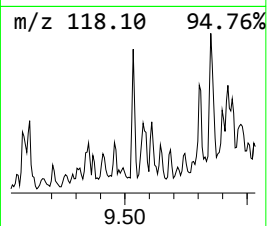
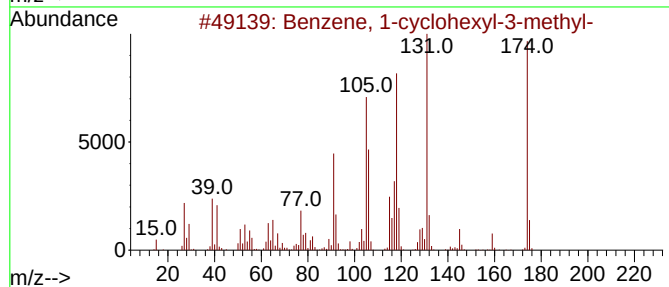
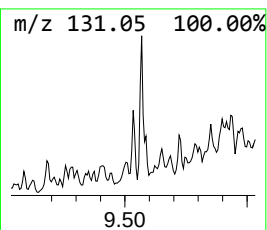
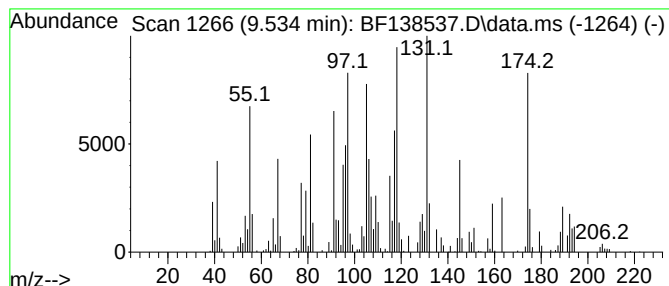
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-cyclohexyl-3-met... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.534	3.33 ng	368765	Acenaphthene-d10	9.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-cyclohexyl-3-methyl-	174	C13H18	004575-46-6	55
2			1-Methyl-2-cyclohexylbenzene	174	C13H18	004501-35-3	38
3			(1,4-Dimethylpent-2-enyl)benzene	174	C13H18	1000184-97-9	38
4			2H-1-Benzopyran-2-one 3,4-dimethyl-	174	C11H10O2	004281-39-4	27
5			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	27



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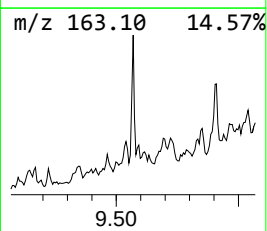
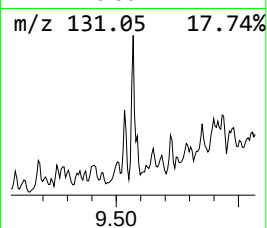
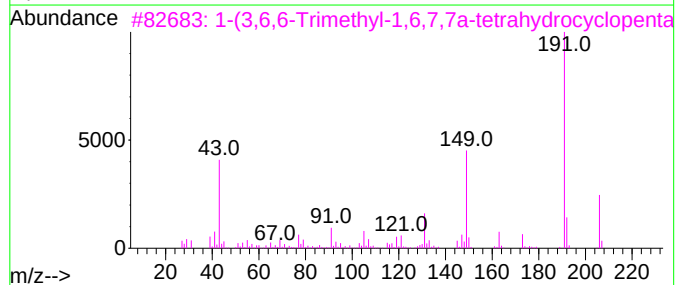
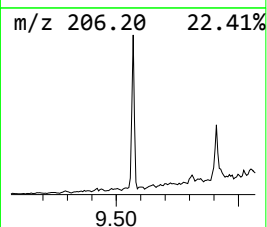
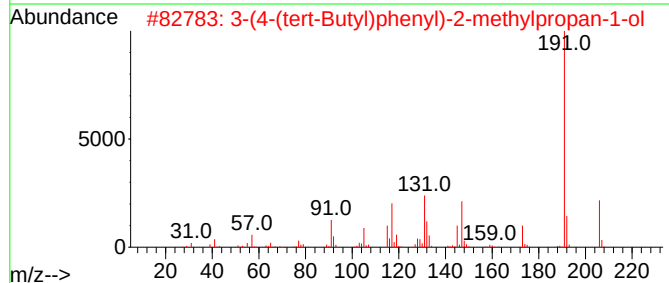
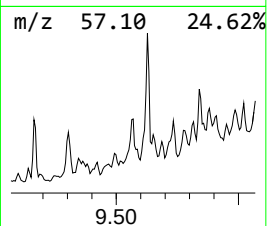
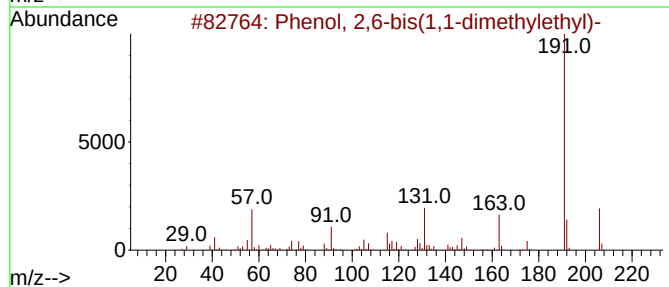
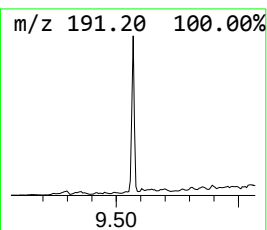
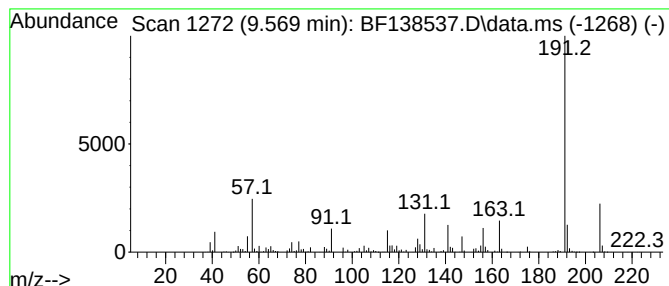
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Phenol, 2,6-bis(1,1-dimethyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.569	11.40 ng	1261010	Acenaphthene-d10	9.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,6-bis(1,1-dimethylethyl)-	206	C14H22O	000128-39-2	87
2			3-(4-(tert-Butyl)phenyl)-2-methy...	206	C14H22O	056107-04-1	68
3			1-(3,6,6-Trimethyl-1,6,7,7a-tetr...	206	C13H18O2	1000194-97-2	68
4			Phenol, 3,5-bis(1,1-dimethylethyl)-	206	C14H22O	001138-52-9	59
5			3-(4-tert-Butylphenyl)propanoic ...	206	C13H18O2	001208-64-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071124\
Data File : BF138537.D
Acq On : 11 Jul 2024 22:23
Operator : CG\JU
Sample : P3181-01 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OR-02-070924

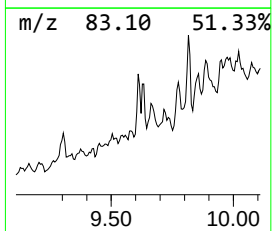
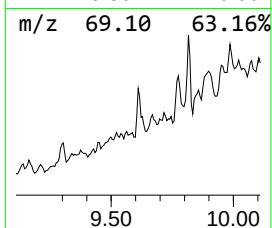
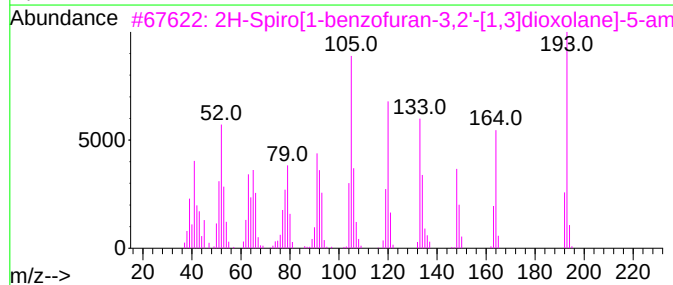
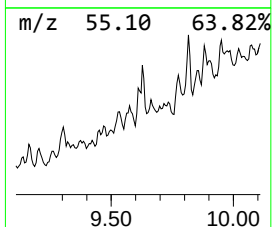
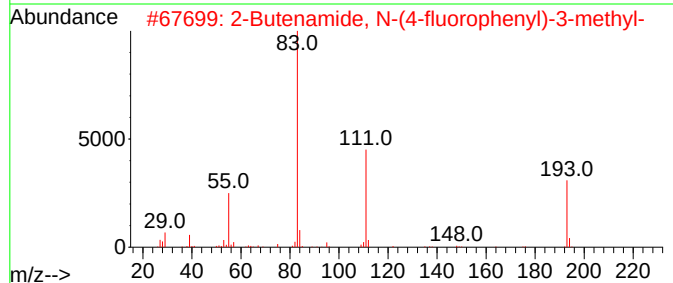
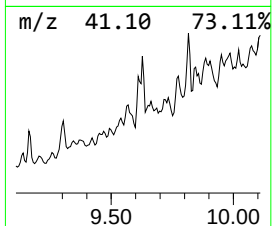
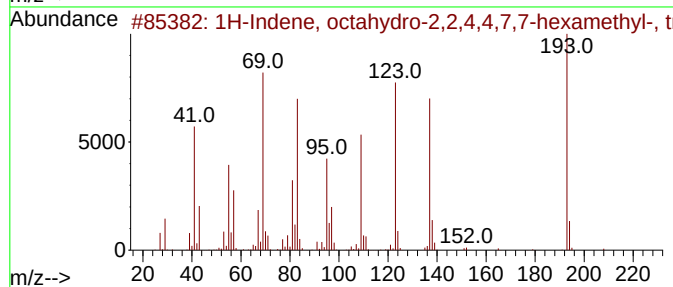
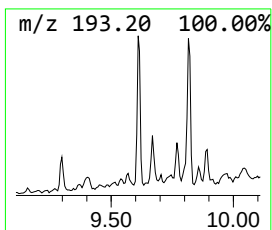
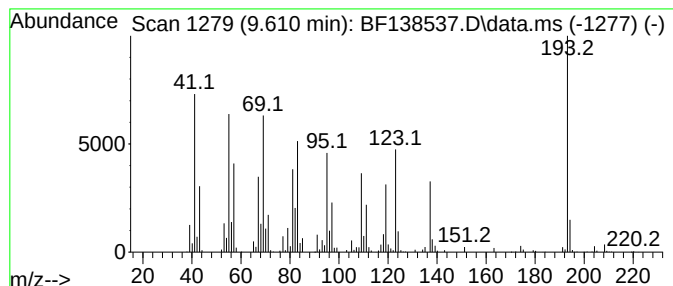
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1H-Indene, octahydro-2,2,4,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.610	10.41 ng	1151420	Acenaphthene-d10	9.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	50		
2	2-Butenamide, N-(4-fluorophenyl)...	193	C11H12FNO	1000307-26-7	30		
3	2H-Spiro[1-benzofuran-3,2'-[1,3]...	193	C10H11NO3	1000387-33-4	25		
4	9-Vinylcarbazole	193	C14H11N	001484-13-5	14		
5	Iminostilbene	193	C14H11N	000256-96-2	14		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071124\
Data File : BF138537.D
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Sample : P3181-01 5X
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ALS Vial : 23 Sample Multiplier: 1

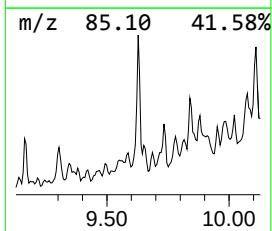
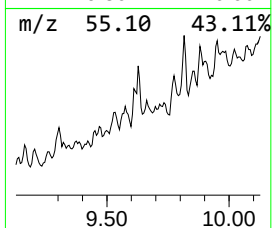
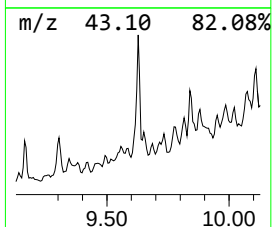
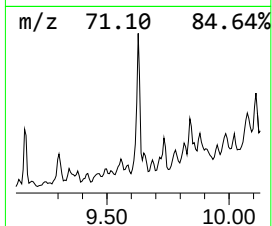
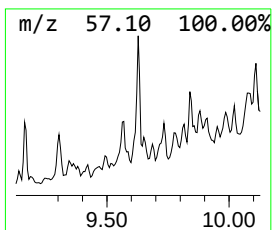
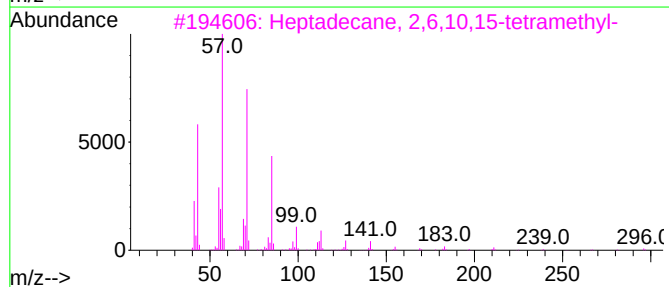
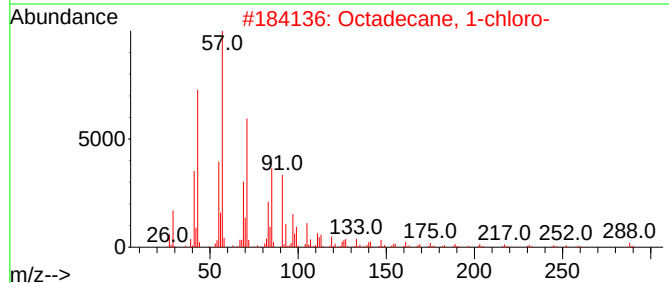
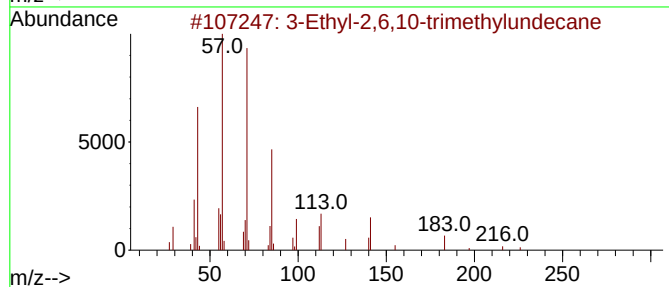
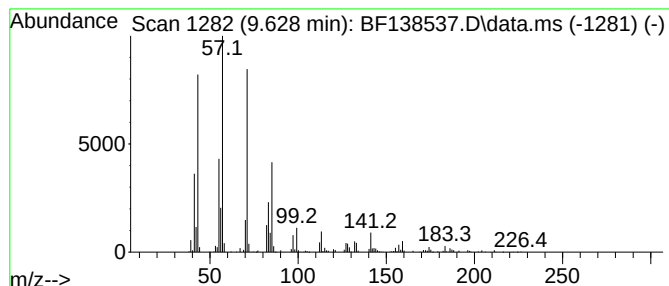
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 3-Ethyl-2,6,10-trimethylund... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
9.628	6.22 ng	688010	Acenaphthene-d10		9.904	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	3-Ethyl-2,6,10-trimethylundecane		226	C16H34	1000432-25-9	91
2	Octadecane, 1-chloro-		288	C18H37Cl	003386-33-2	86
3	Heptadecane, 2,6,10,15-tetramethyl-		296	C21H44	054833-48-6	80
4	Tetradecane, 1-iodo-		324	C14H29I	019218-94-1	80
5	Decane, 2,3,5,8-tetramethyl-		198	C14H30	192823-15-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071124\
Data File : BF138537.D
Acq On : 11 Jul 2024 22:23
Operator : CG\JU
Sample : P3181-01 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

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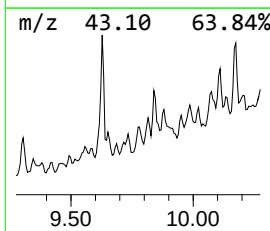
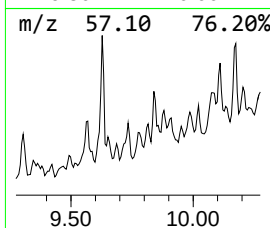
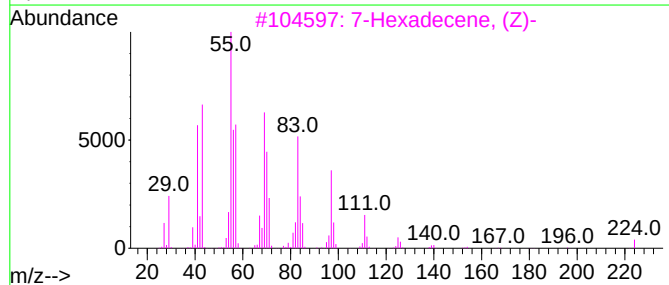
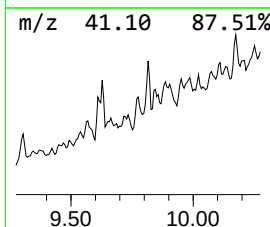
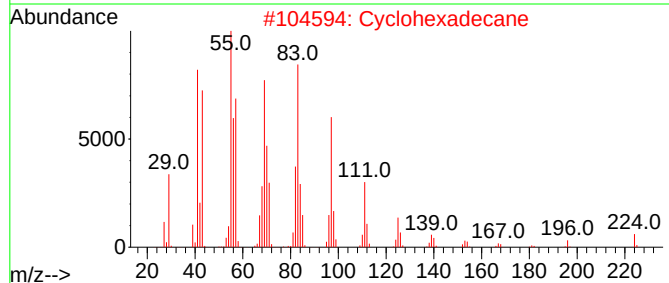
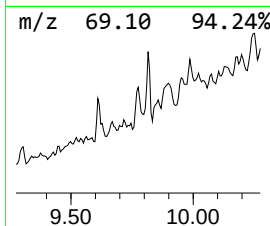
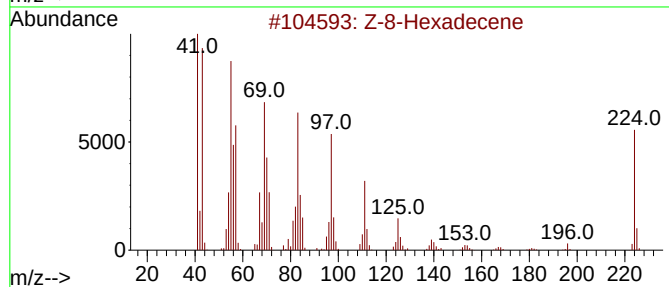
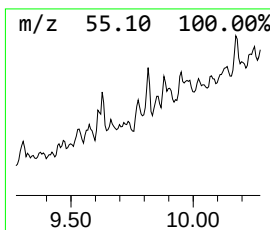
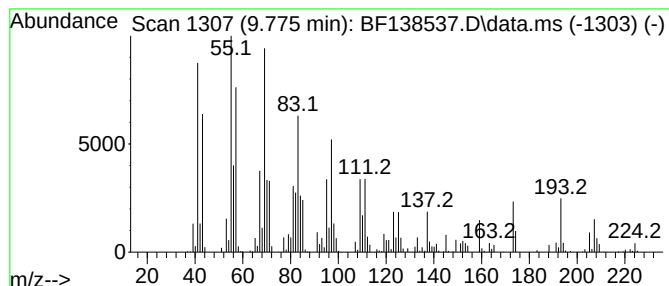
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Z-8-Hexadecene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.775	13.54 ng	1497290	Acenaphthene-d10	9.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Z-8-Hexadecene	224	C16H32	1000130-87-5	83
2			Cyclohexadecane	224	C16H32	000295-65-8	83
3			7-Hexadecene, (Z)-	224	C16H32	035507-09-6	83
4			2-Heptafluorobutyropentadecane	424	C19H31F7O2	1010245-50-0	62
5			3-Eicosene, (E)-	280	C20H40	074685-33-9	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071124\
Data File : BF138537.D
Acq On : 11 Jul 2024 22:23
Operator : CG\JU
Sample : P3181-01 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

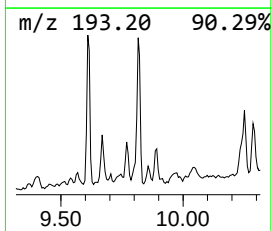
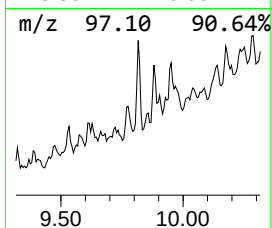
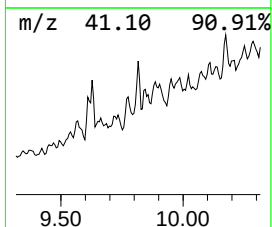
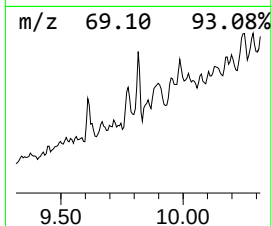
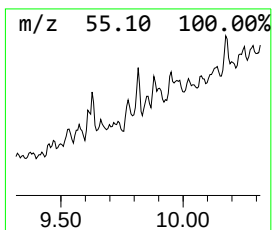
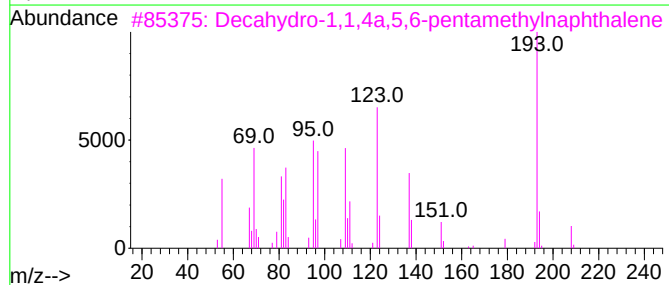
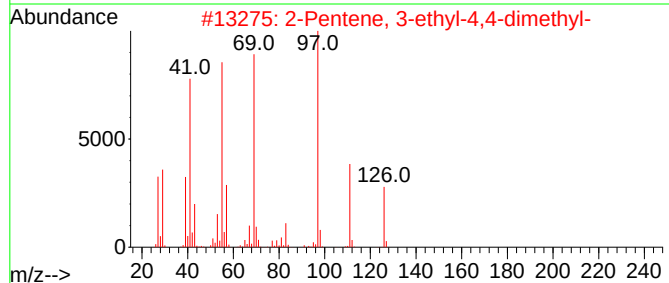
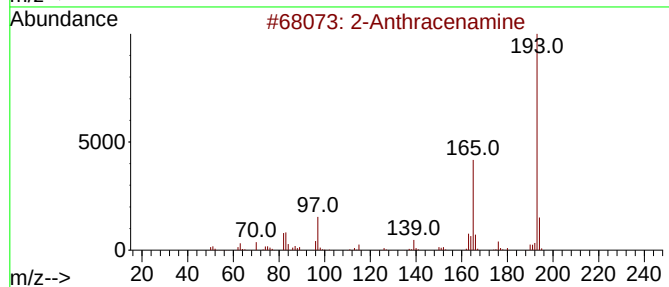
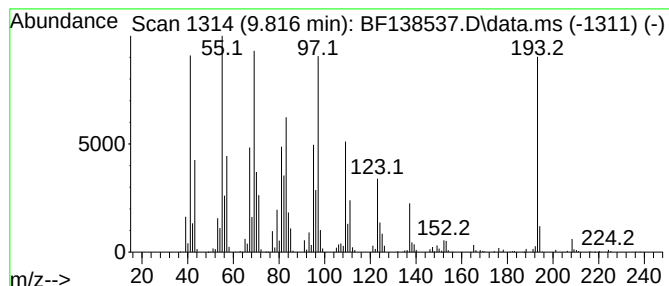
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown9.816 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.816	12.98 ng	1435960	Acenaphthene-d10	9.904	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Anthracenamine	193	C14H11N	000613-13-8	46
2	2-Pentene, 3-ethyl-4,4-dimethyl-	126	C9H18	053907-59-8	38
3	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	30
4	Cyclohexane, 1-(cyclohexylmethyl...	208	C15H28	054934-92-8	30
5	Cyclododecanone, 2-methylene-	194	C13H22O	003045-76-9	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071124\
Data File : BF138537.D
Acq On : 11 Jul 2024 22:23
Operator : CG\JU
Sample : P3181-01 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
OR-02-070924

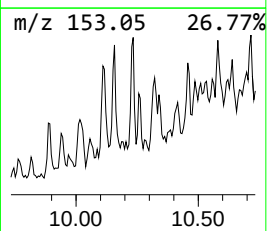
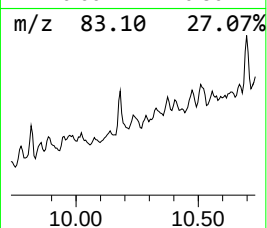
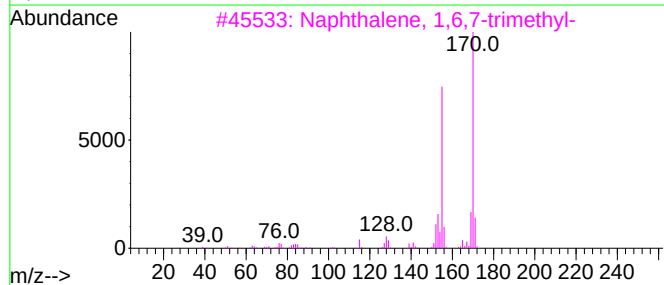
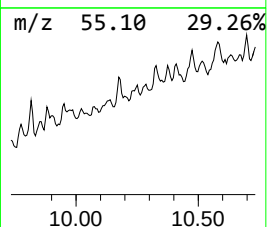
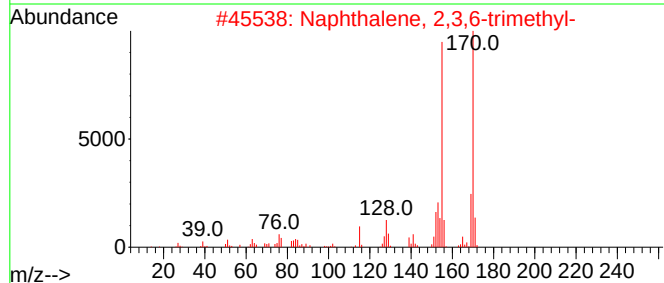
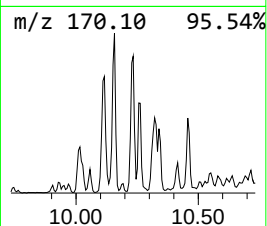
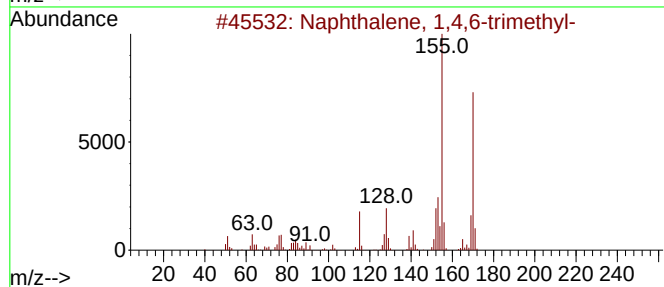
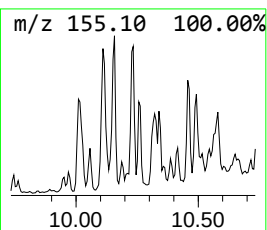
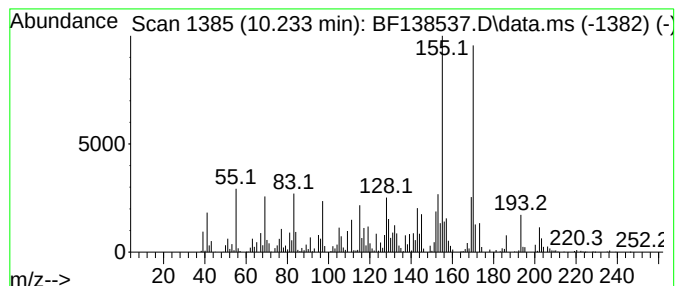
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Naphthalene, 1,4,6-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.233	9.98 ng	1103780	Acenaphthene-d10	9.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	97
2			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
3			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
4			3-(2-Methyl-propenyl)-1H-indene	170	C13H14	1000187-78-5	95
5			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071124\
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Sample : P3181-01 5X
Misc :
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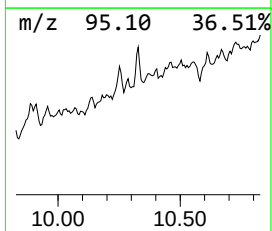
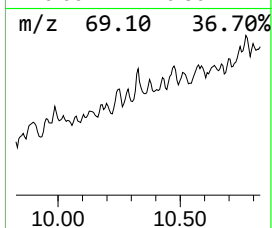
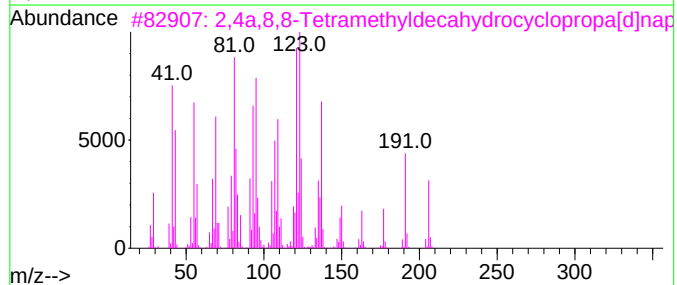
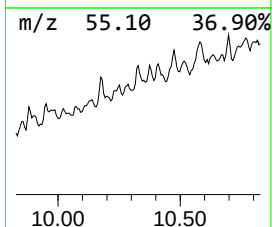
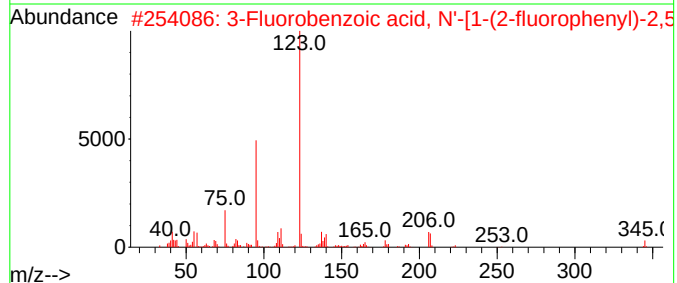
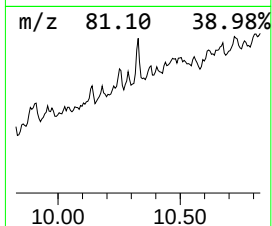
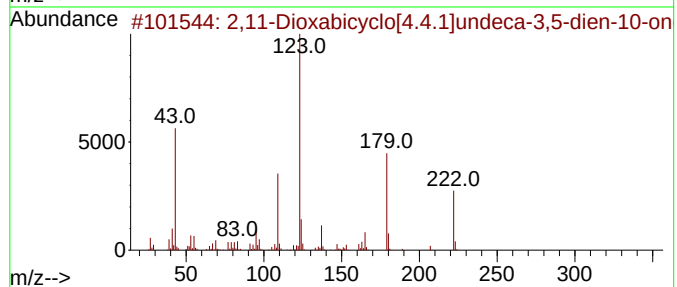
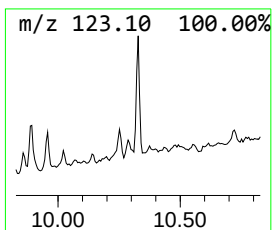
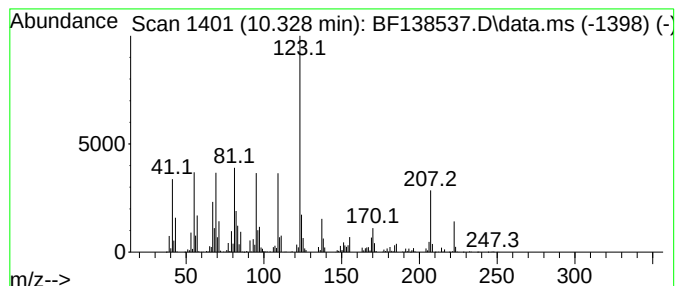
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2,11-Dioxabicyclo[4.4.1]und... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.328	15.84 ng	1752090	Acenaphthene-d10	9.904		
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,11-Dioxabicyclo[4.4.1]undeca-3...	222	C13H18O3		070412-52-1	53
2	3-Fluorobenzoic acid, N'-[1-(2-f...	345	C17H13F2N3O3		1000304-52-3	50
3	2,4a,8,8-Tetramethyldecahydrocyc...	206	C15H26		074022-04-1	50
4	Thiazolidine-2-thione, 4-(4-fluo...	320	C15H13FN2OS2		1000264-97-7	47
5	Bromo-4-fluoroacetophenone	216	C8H6BrFO		000403-29-2	47



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Data File : BF138537.D
Acq On : 11 Jul 2024 22:23
Operator : CG\JU
Sample : P3181-01 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OR-02-070924

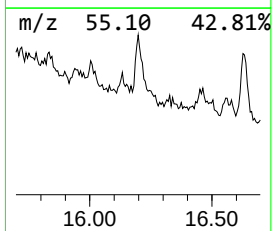
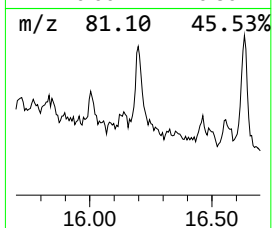
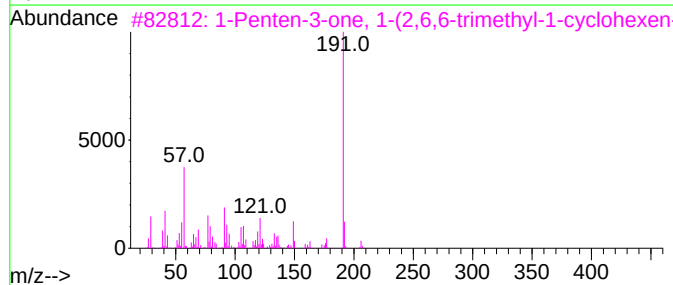
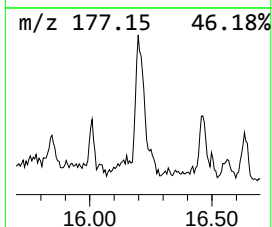
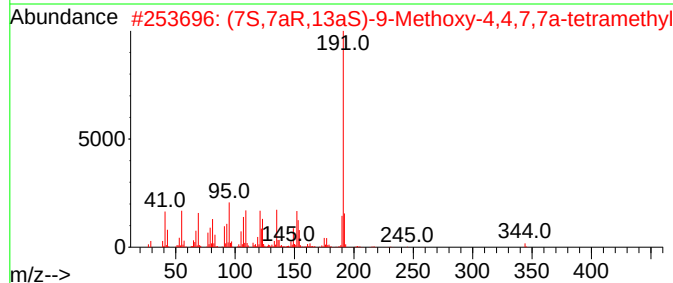
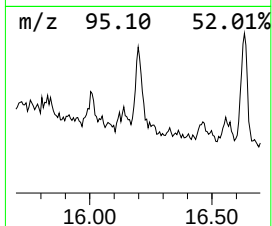
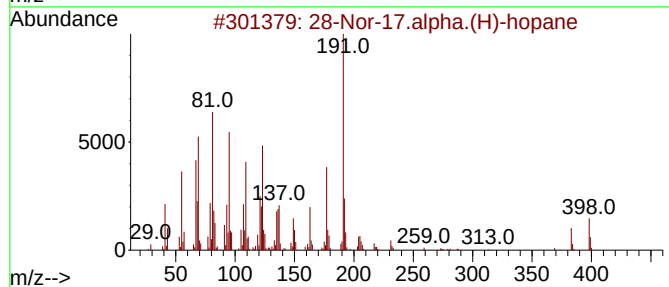
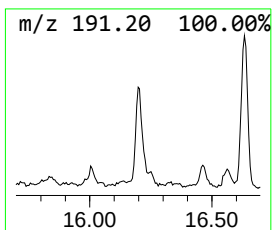
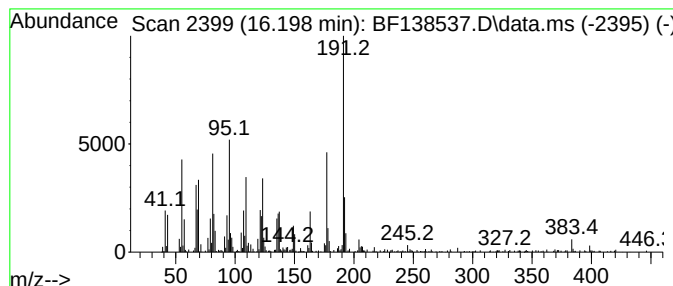
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 28-Nor-17.alpha.(H)-hopane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.198	28.27 ng	789713	Perylene-d12	15.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			28-Nor-17.alpha.(H)-hopane	398	C29H50	053584-60-4	70
2			(7S,7aR,13aS)-9-Methoxy-4,4,7,7a...	344	C22H32O3	1000482-45-3	50
3			1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	35
4			.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	30
5			4,5-Dihydrothiazol-4-one, 2-(2-m...	222	C10H10N2O2S	022476-61-5	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071124\
Data File : BF138537.D
Acq On : 11 Jul 2024 22:23
Operator : CG\JU
Sample : P3181-01 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OR-02-070924

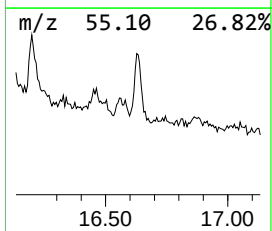
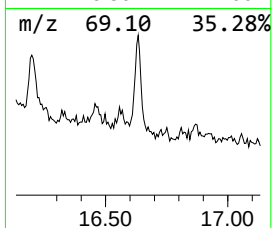
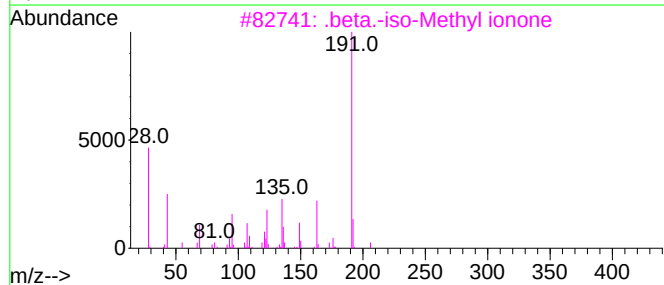
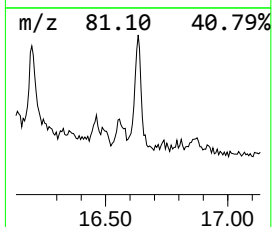
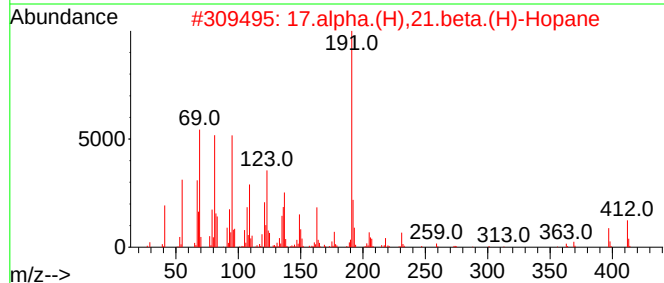
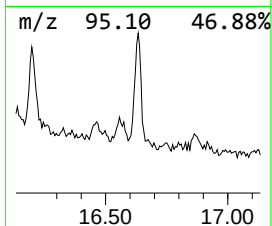
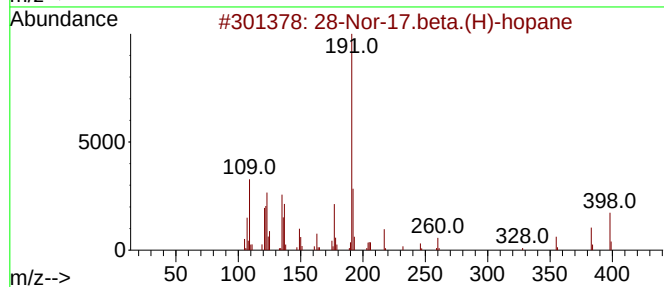
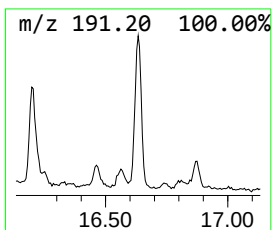
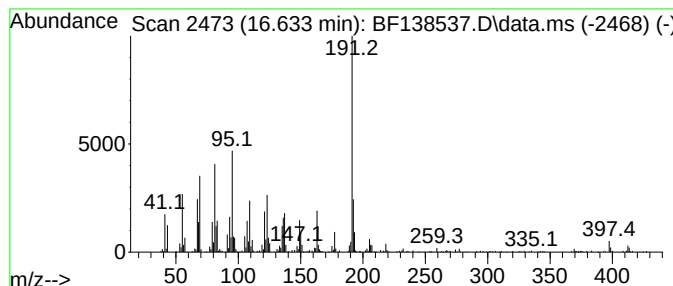
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 28-Nor-17.beta.(H)-hopane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.633	24.99 ng	698016	Perylene-d12	15.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	59
2			17.alpha.(H),21.beta.(H)-Hopane	412	C30H52	013849-96-2	50
3			.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	49
4			28-Nor-17.alpha.(H)-hopane	398	C29H50	053584-60-4	49
5			(4aS,8aS)-8-Isopentyl-4,4,7,8a-t...	262	C19H34	220766-79-0	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071124\
Data File : BF138537.D
Acq On : 11 Jul 2024 22:23
Operator : CG\JU
Sample : P3181-01 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OR-02-070924

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF070924.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
Cyclohexane, 1,...	8.681	7.7	ng	168619	2	8.145	440137	20.0
unknown9.022	9.022	9.1	ng	201075	2	8.145	440137	20.0
Nonane, 3,7-dim...	9.163	5.2	ng	579442	3	9.904	2211860	20.0
1-Decanol, 2-he...	9.304	6.8	ng	753892	3	9.904	2211860	20.0
Benzene, 1-cycl...	9.534	3.3	ng	368765	3	9.904	2211860	20.0
Phenol, 2,6-bis...	9.569	11.4	ng	1261010	3	9.904	2211860	20.0
1H-Indene, octa...	9.610	10.4	ng	1151420	3	9.904	2211860	20.0
3-Ethyl-2,6,10-...	9.628	6.2	ng	688010	3	9.904	2211860	20.0
Z-8-Hexadecene	9.775	13.5	ng	1497290	3	9.904	2211860	20.0
unknown9.816	9.816	13.0	ng	1435960	3	9.904	2211860	20.0
Naphthalene, 1,...	10.233	10.0	ng	1103780	3	9.904	2211860	20.0
2,11-Dioxabicyc...	10.328	15.8	ng	1752090	3	9.904	2211860	20.0
28-Nor-17.alpha...	16.198	28.3	ng	789713	6	15.504	558684	20.0
28-Nor-17.beta....	16.633	25.0	ng	698016	6	15.504	558684	20.0