

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF062718\
 Data File : BF106877.D
 Acq On : 27 Jun 2018 17:35
 Operator : JU/SJ
 Sample : J3570-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20180374-AMENDED-COMPOSITE-

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

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-BF061918.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.878	429	432	441	rBV	18864	23001	1.01%	0.272%
2	5.231	480	492	494	rBV	1013210	2268387	100.00%	26.778%
3	6.454	697	700	707	rBV	23986	29798	1.31%	0.352%
4	6.595	721	724	733	rBV2	113167	131638	5.80%	1.554%
5	6.731	745	747	751	rBV	43847	34353	1.51%	0.406%
6	6.954	781	785	788	rBV	296676	254436	11.22%	3.004%
7	7.107	807	811	815	rVB	802754	717597	31.63%	8.471%
8	7.519	876	881	884	rBV	610136	580564	25.59%	6.853%
9	7.672	904	907	912	rBV	29275	27809	1.23%	0.328%
10	8.236	999	1003	1006	rBV	388493	321693	14.18%	3.798%
11	8.378	1023	1027	1034	rBV	51043	78071	3.44%	0.922%
12	9.307	1181	1185	1189	rBV	1066930	1024429	45.16%	12.093%
13	9.383	1194	1198	1201	rBV3	33595	43100	1.90%	0.509%
14	9.701	1247	1252	1257	rBV	141538	144414	6.37%	1.705%
15	9.854	1274	1278	1280	rBV3	23713	25491	1.12%	0.301%
16	9.901	1283	1286	1288	rVB2	27815	27772	1.22%	0.328%
17	9.966	1294	1297	1299	rBV2	24248	25448	1.12%	0.300%
18	9.995	1299	1302	1305	rVB	402183	343664	15.15%	4.057%
19	10.395	1368	1370	1374	rVB2	38177	39063	1.72%	0.461%
20	10.789	1434	1437	1439	rBV3	30611	26772	1.18%	0.316%
21	10.878	1449	1452	1457	rBV2	25587	49259	2.17%	0.581%
22	10.919	1457	1459	1462	rVB	35653	30095	1.33%	0.355%
23	11.354	1530	1533	1540	rVB6	26024	42130	1.86%	0.497%
24	11.489	1552	1556	1559	rBV	387925	349959	15.43%	4.131%
25	12.707	1760	1763	1765	rBV	22472	23673	1.04%	0.279%
26	12.936	1800	1802	1807	rVB	44787	53048	2.34%	0.626%
27	13.071	1821	1825	1828	rBV	902871	770125	33.95%	9.091%
28	13.948	1971	1974	1980	rBV2	54485	63843	2.81%	0.754%
29	14.142	2003	2007	2010	rBV	315495	292513	12.90%	3.453%
30	15.354	2209	2213	2219	rVB8	15479	31753	1.40%	0.375%
31	15.612	2254	2257	2261	rVV2	30708	43928	1.94%	0.519%
32	15.683	2263	2269	2284	rVB	265664	399473	17.61%	4.716%
33	16.124	2342	2344	2350	rVB6	16544	23178	1.02%	0.274%
34	16.412	2389	2393	2395	rBV2	21152	30055	1.32%	0.355%

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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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	16.871	2465	2471	2478	rvB3	26456	51667	2.28%	0.610%
36	17.483	2571	2575	2582	rvB6	25504	48975	2.16%	0.578%

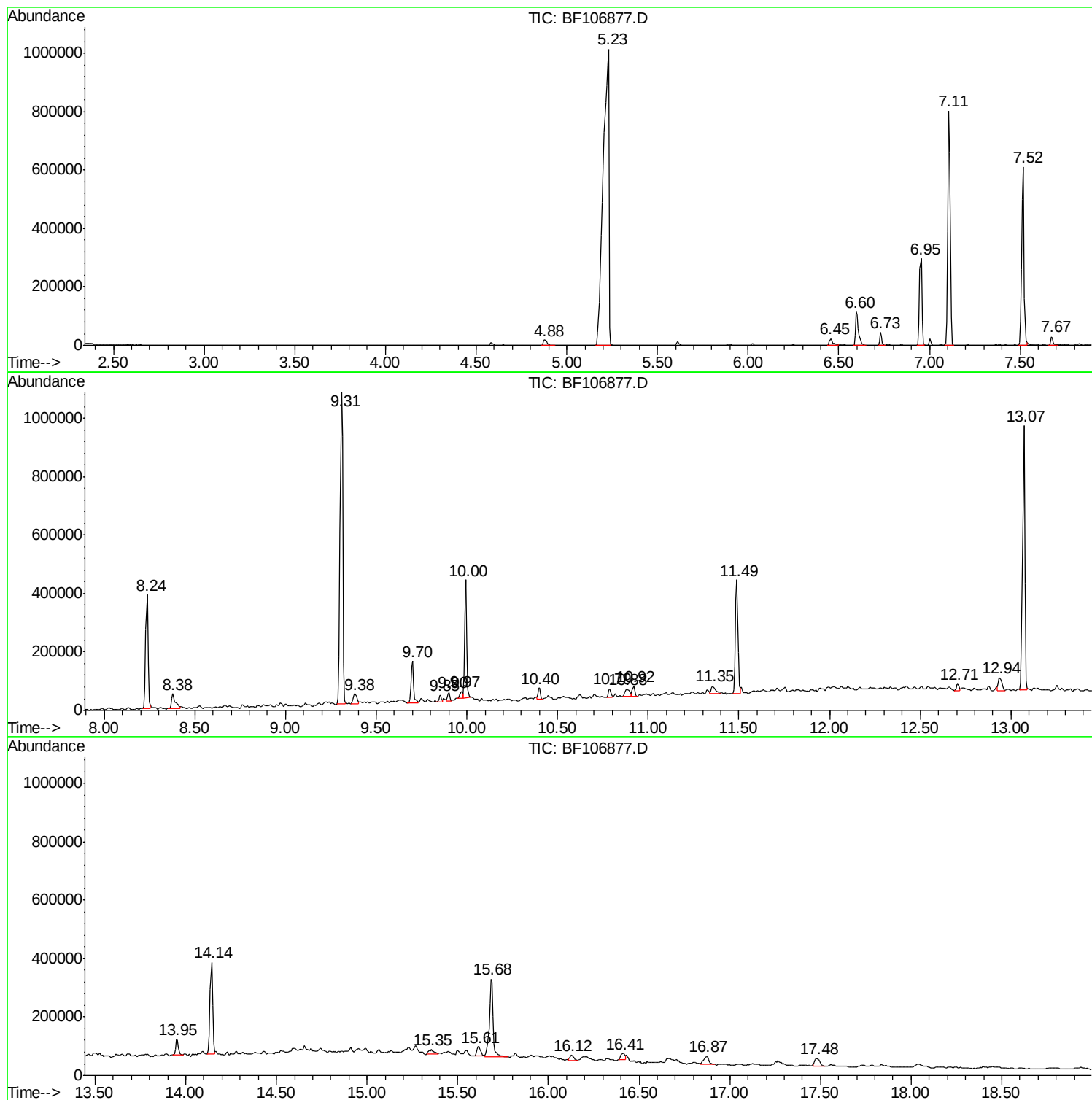
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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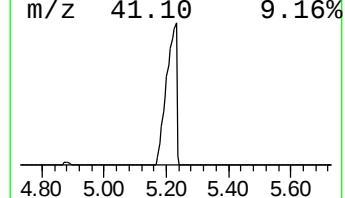
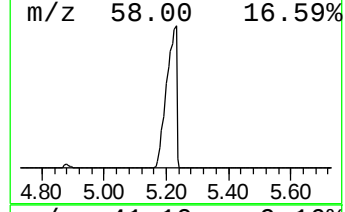
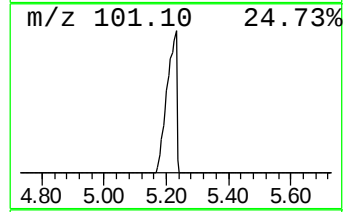
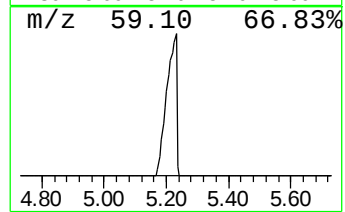
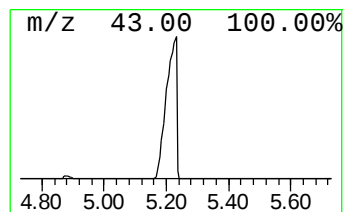
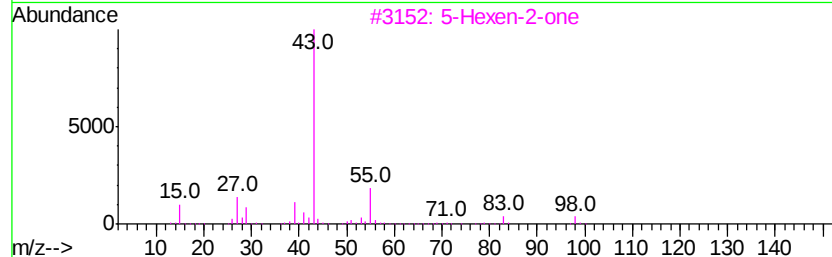
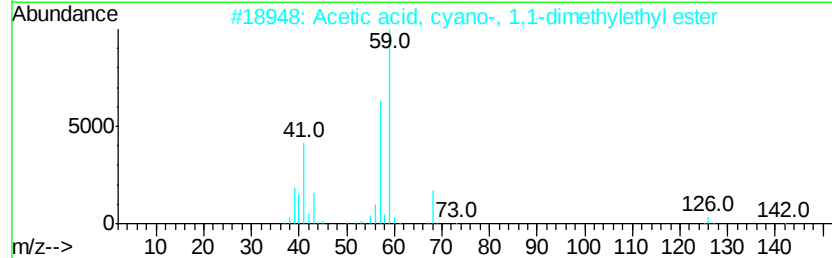
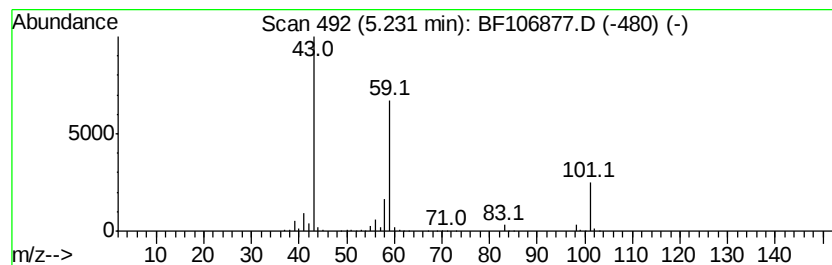
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.23	178.31 ng	2268390	1,4-Dichlorobenzene-d4	6.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		5-Hexen-2-one	98	C6H10O	000109-49-9	9
4		Guanidine	59	CH5N3	000113-00-8	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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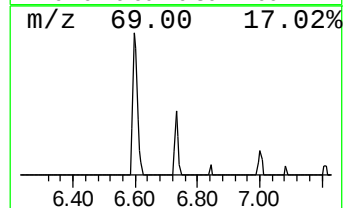
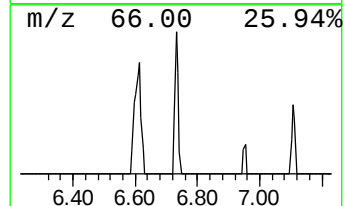
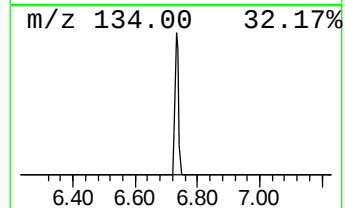
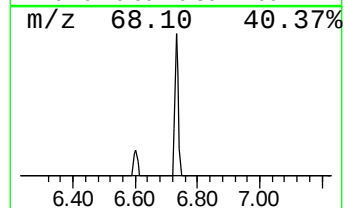
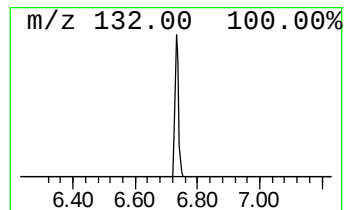
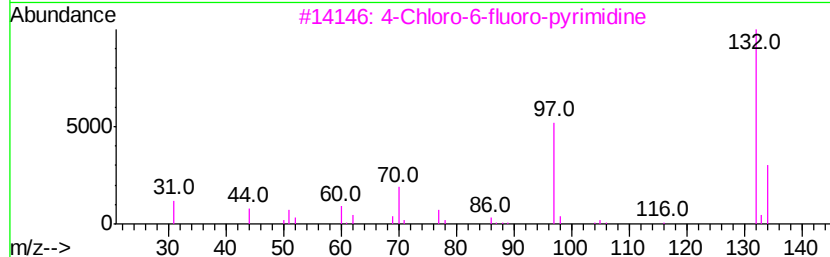
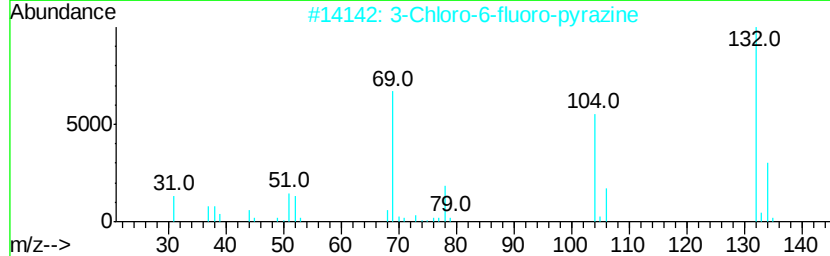
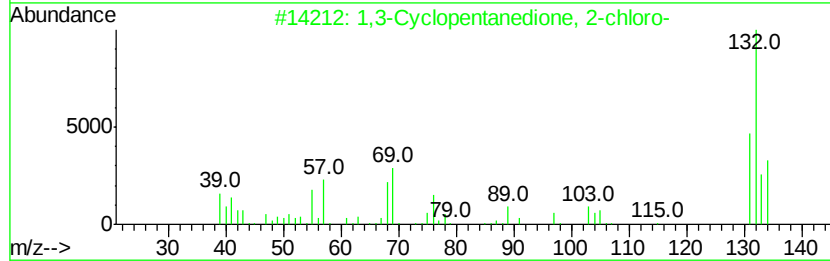
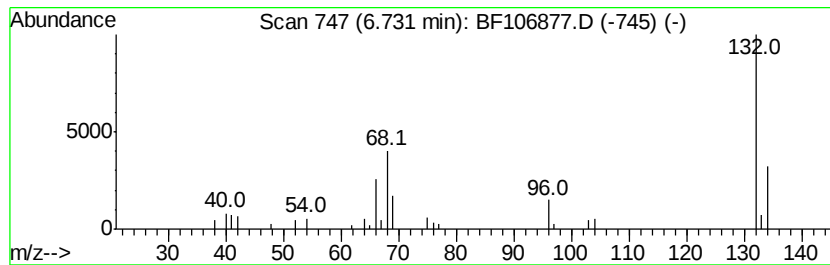
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.73 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	2.70 ng	34353	1,4-Dichlorobenzene-d4	6.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	43
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	9



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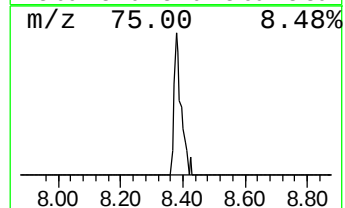
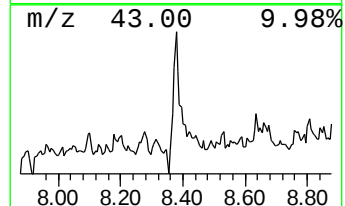
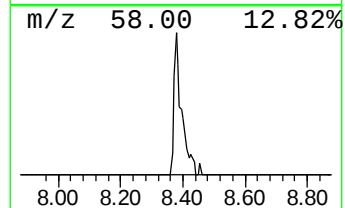
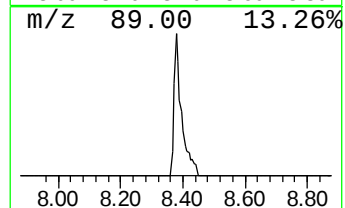
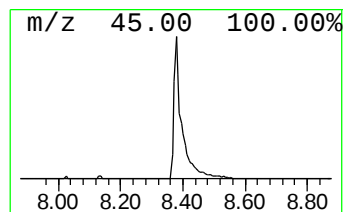
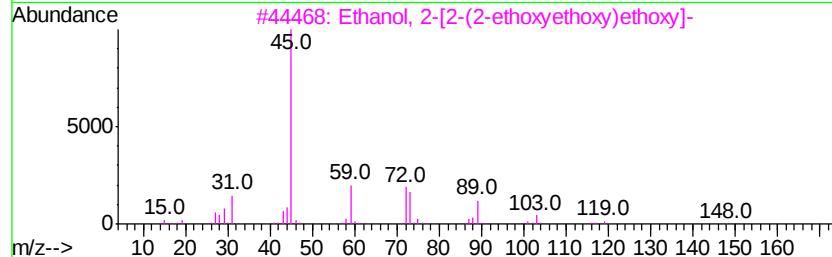
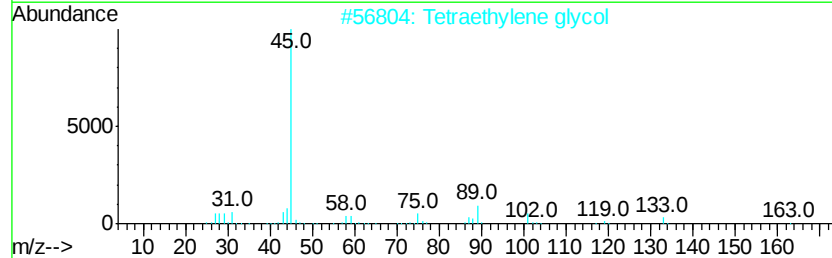
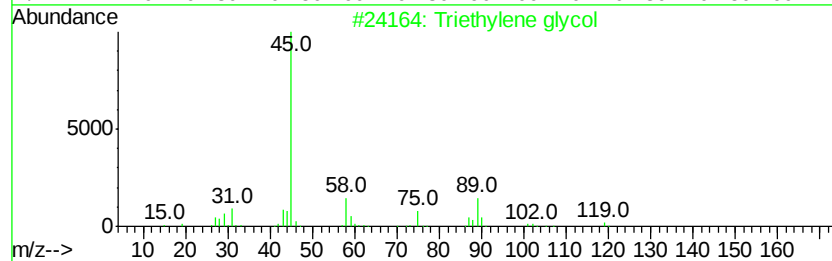
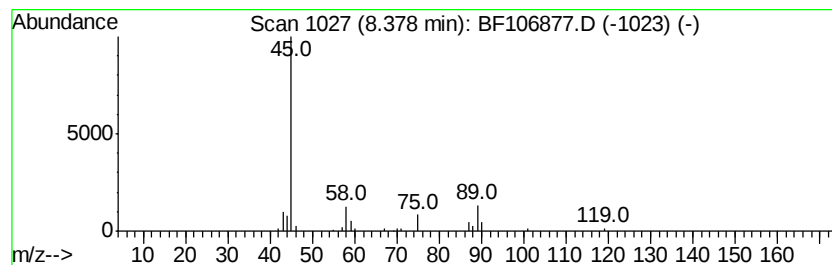
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Triethylene glycol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.38	4.85 ng	78071	Naphthalene-d8	8.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triethylene glycol	150	C6H14O4	000112-27-6	90
2		Tetraethylene glycol	194	C8H18O5	000112-60-7	72
3		Ethanol, 2-[2-(2-ethoxyethoxy)et...	178	C8H18O4	000112-50-5	50
4		Isopropyl Alcohol	60	C3H8O	000067-63-0	42
5		2-Hexanol	102	C6H14O	000626-93-7	36



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ClientSampleId :
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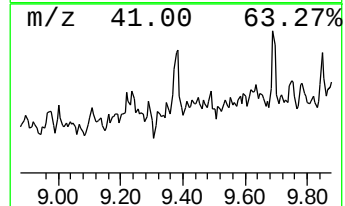
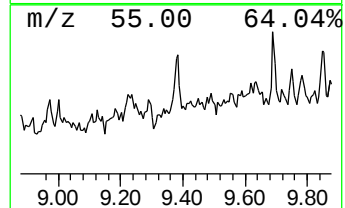
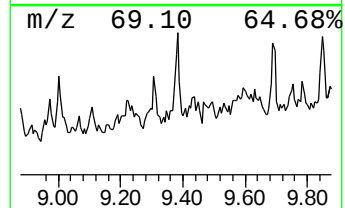
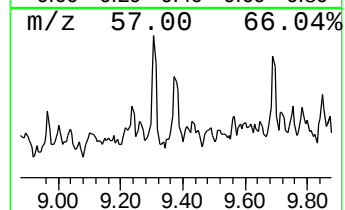
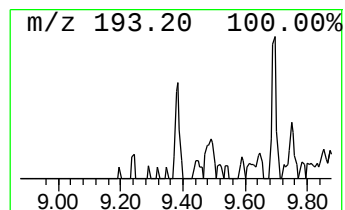
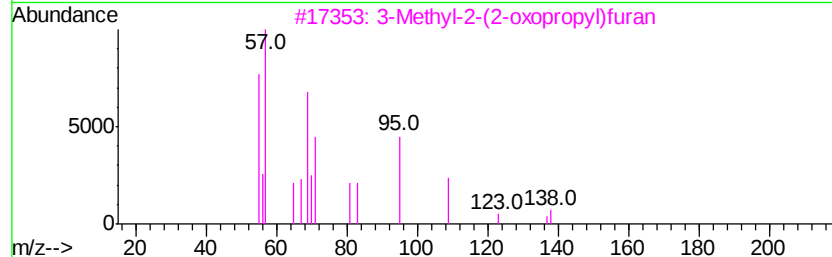
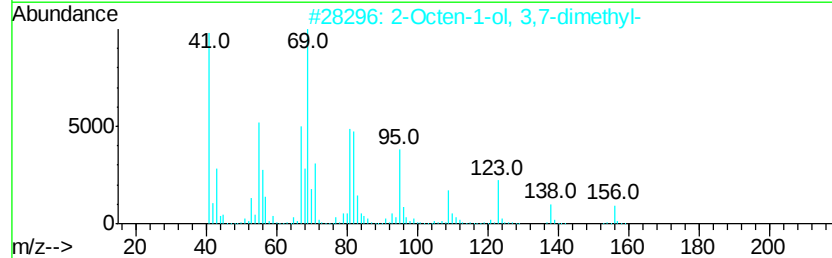
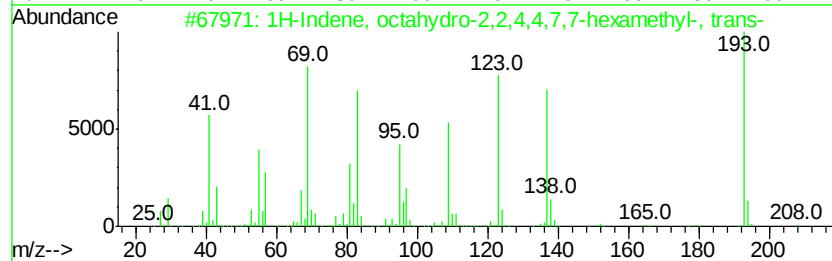
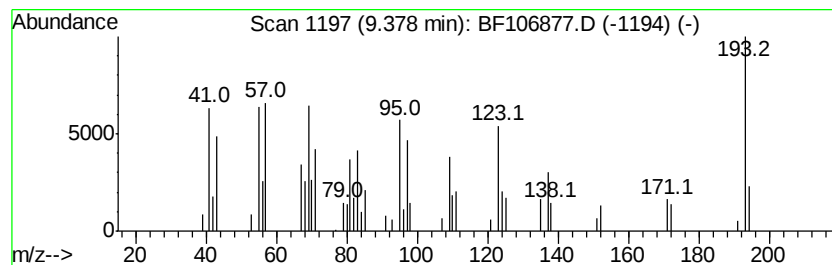
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1H-Indene, octahydro-2,2,4,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.38	2.51 ng	43100	Acenaphthene-d10	10.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	53
2		2-Octen-1-ol, 3,7-dimethyl-	156	C10H20O	040607-48-5	45
3		3-Methyl-2-(2-oxopropyl)furan	138	C8H10O2	087773-62-4	38
4		Bicyclo[3.1.1]heptan-3-one, 6,6-...	194	C13H22O	1000163-97-9	22
5		1-Nonanol, 9-bromo-	222	C9H19BrO	055362-80-6	14



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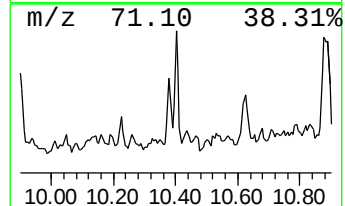
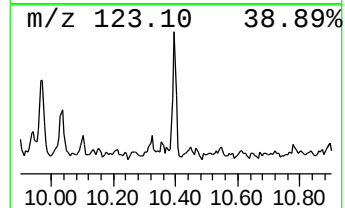
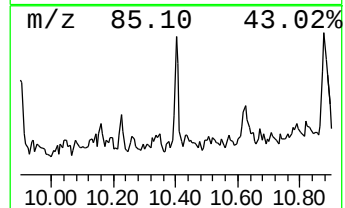
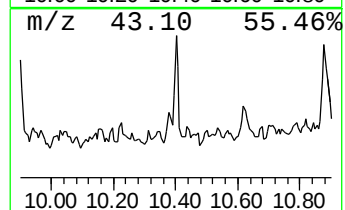
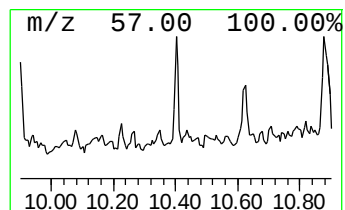
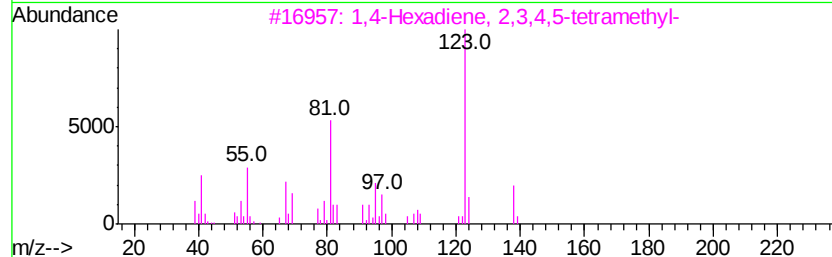
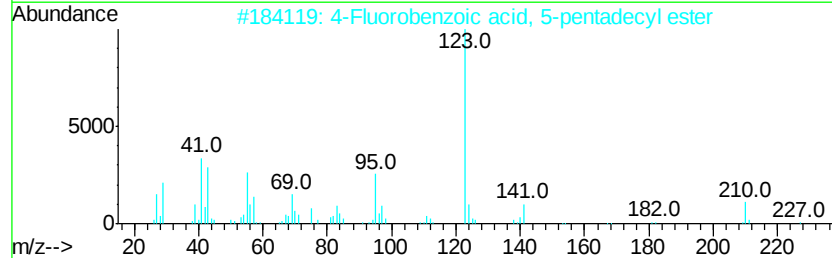
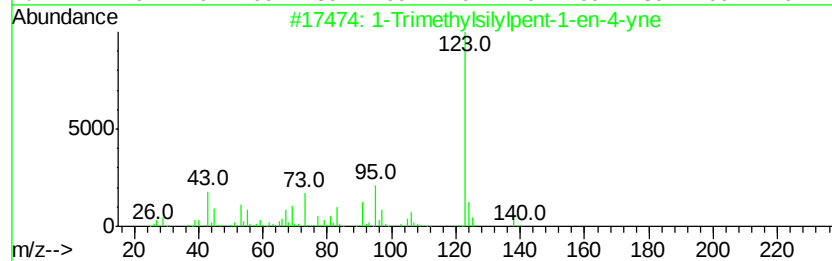
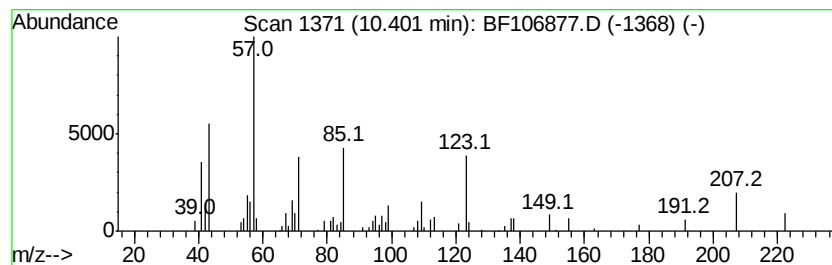
Instrument :
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1-Trimethylsilylpent-1-en-4... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.40	2.27 ng	39063	Acenaphthene-d10	10.00	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Trimethylsilylpent-1-en-4-yne	138	C8H14Si	079516-25-9	50
2	4-Fluorobenzoic acid, 5-pentadec...	350	C22H35FO2	1000280-68-6	47
3	1,4-Hexadiene, 2,3,4,5-tetramethyl-	138	C10H18	051504-54-2	47
4	trans-2-(1-Hydroxycyclohexyl)-furan	166	C10H14O2	115754-87-5	47
5	Benzene, 1,4-dimethoxy-	138	C8H10O2	000150-78-7	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062718\
Data File : BF106877.D
Acq On : 27 Jun 2018 17:35
Operator : JU/SJ
Sample : J3570-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
20180374-AMENDED-COMPOSITE-

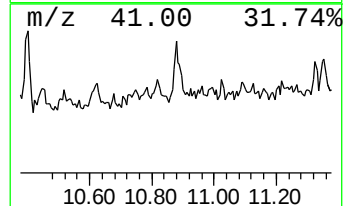
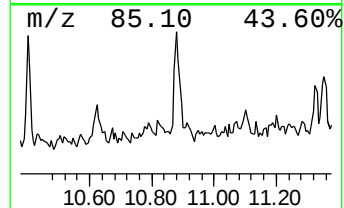
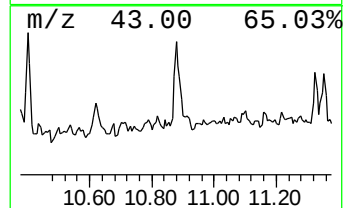
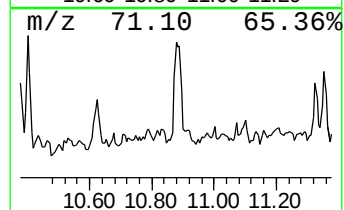
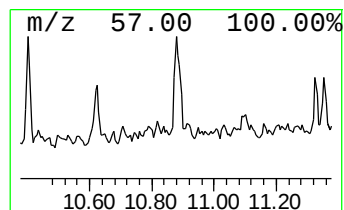
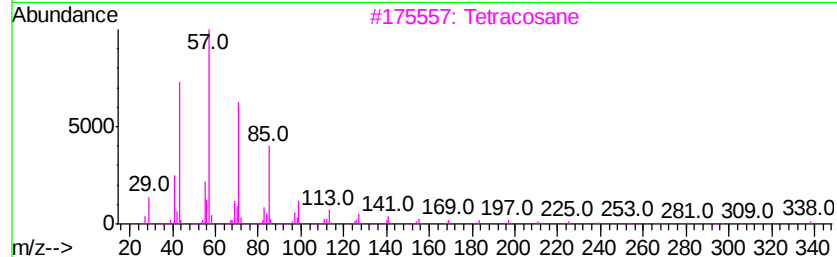
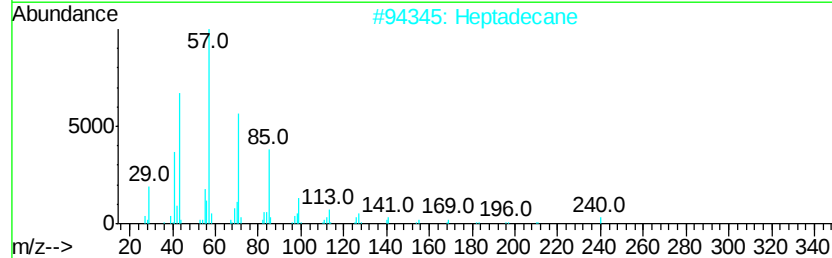
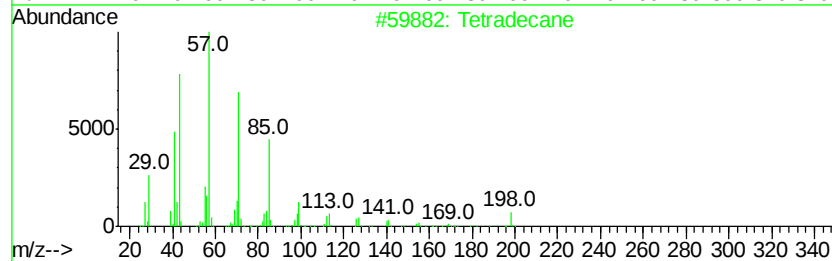
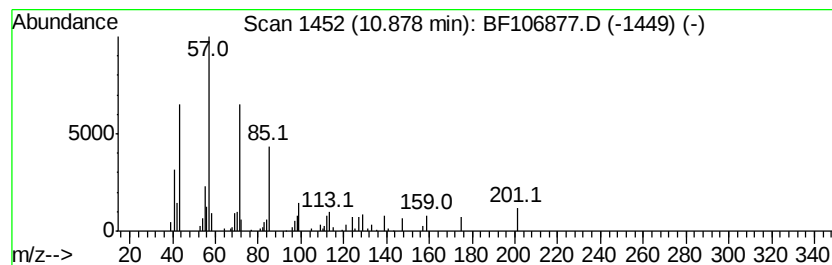
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Tetradecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.88	2.82 ng	49259	Phenanthrene-d10	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	72
2		Heptadecane	240	C17H36	000629-78-7	72
3		Tetracosane	338	C24H50	000646-31-1	72
4		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	72
5		2-Bromo dodecane	248	C12H25Br	013187-99-0	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062718\
Data File : BF106877.D
Acq On : 27 Jun 2018 17:35
Operator : JU/SJ
Sample : J3570-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
20180374-AMENDED-COMPOSITE-

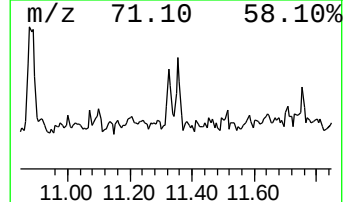
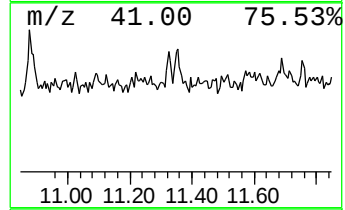
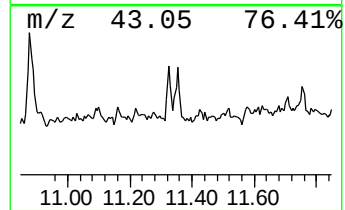
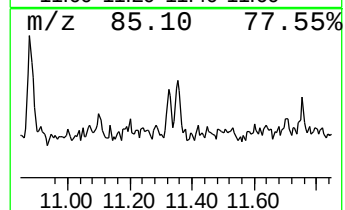
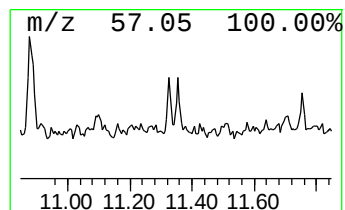
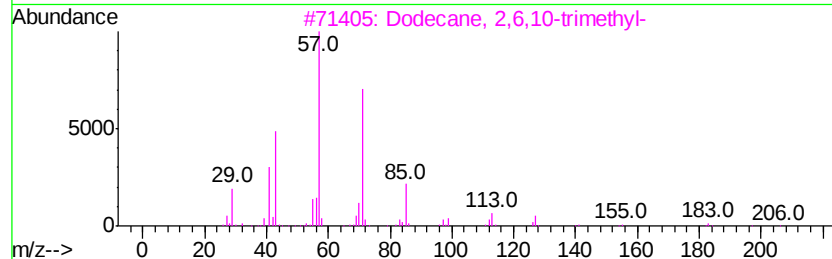
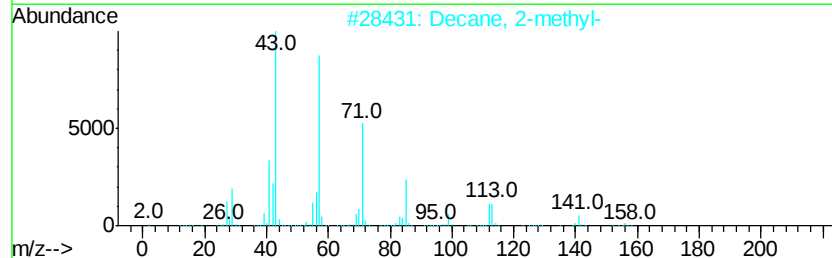
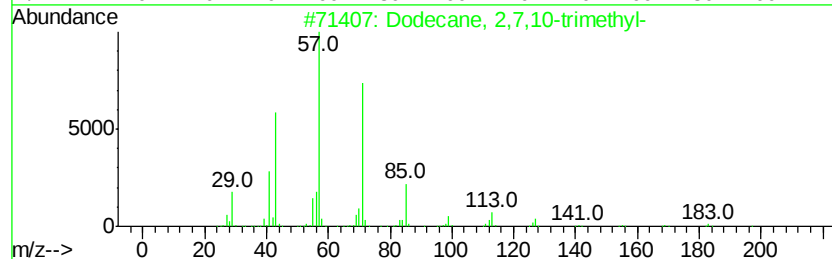
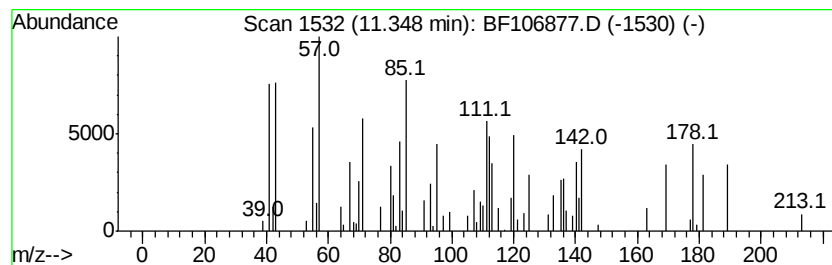
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown11.35 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.35	2.41 ng	42130	Phenanthrene-d10	11.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	38
2			Decane, 2-methyl-	156	C11H24	006975-98-0	38
3			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	38
4			Decane, 5-ethyl-5-methyl-	184	C13H28	017312-74-2	38
5			Nonane, 2-methyl-5-propyl-	184	C13H28	031081-17-1	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062718\
Data File : BF106877.D
Acq On : 27 Jun 2018 17:35
Operator : JU/SJ
Sample : J3570-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
20180374-AMENDED-COMPOSITE-

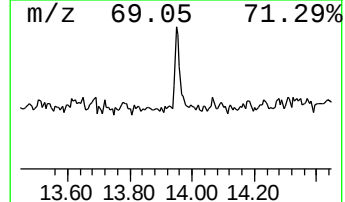
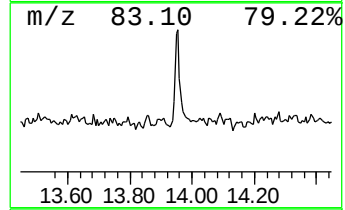
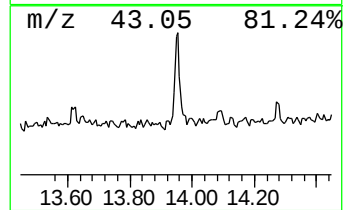
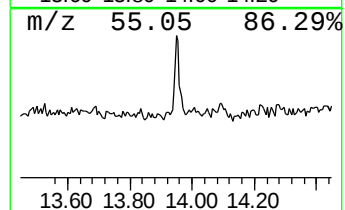
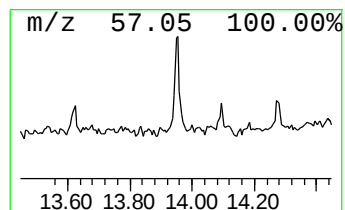
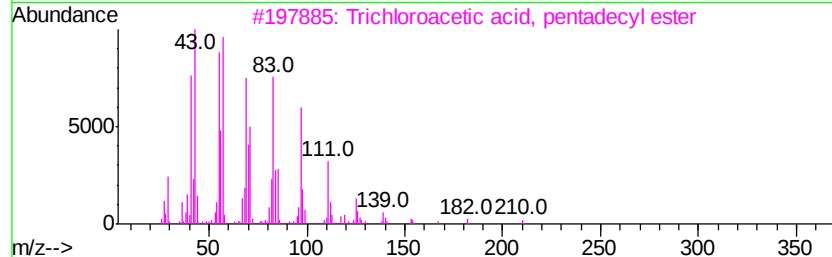
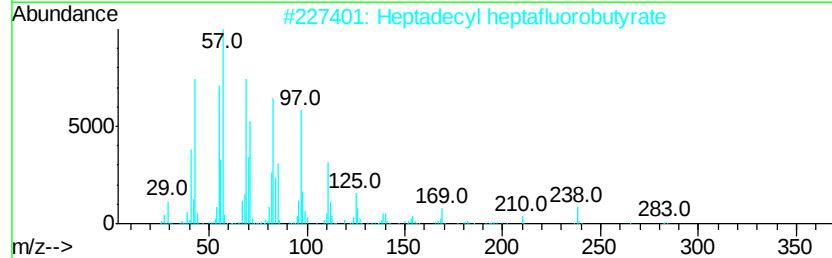
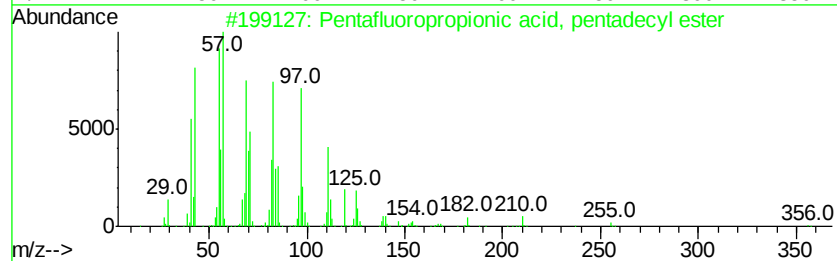
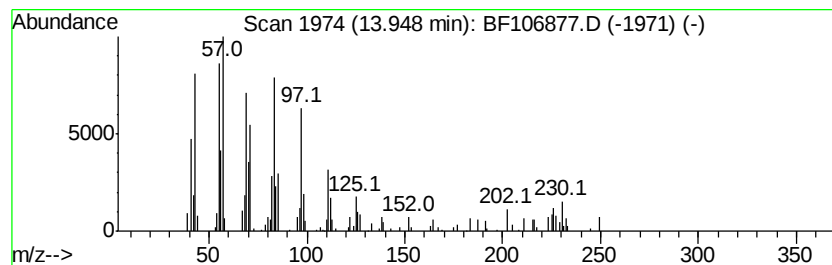
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Pentafluoropropionic acid, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	4.37 ng	63843	Chrysene-d12	14.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	93
2			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	93
3			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	93
4			Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	90
5			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062718\
Data File : BF106877.D
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Sample : J3570-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
20180374-AMENDED-COMPOSITE-

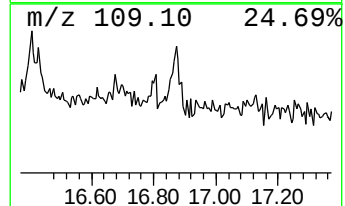
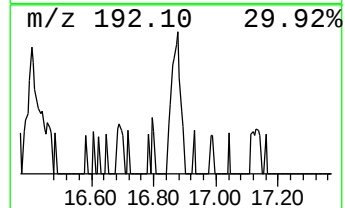
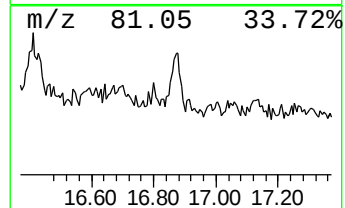
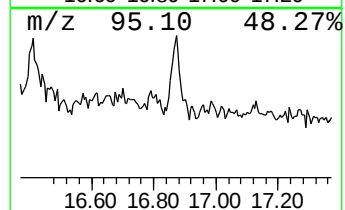
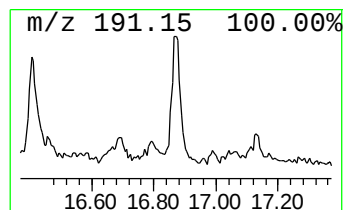
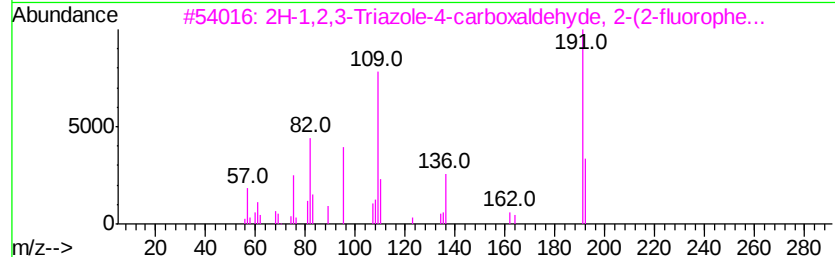
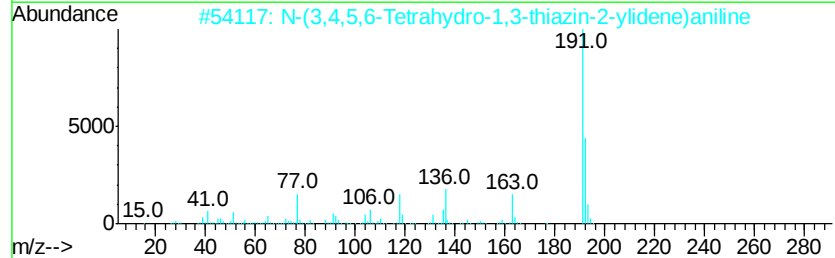
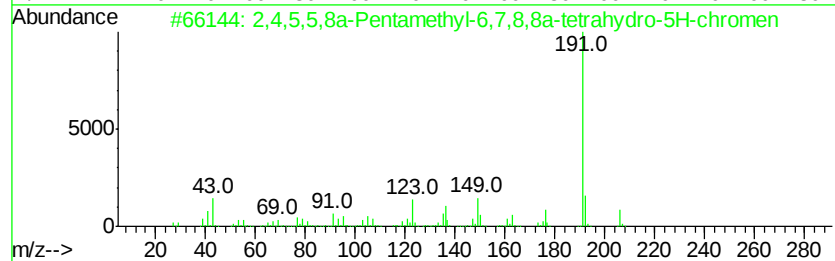
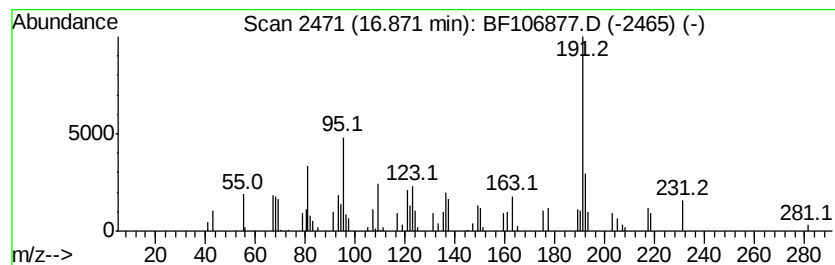
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown16.87 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.87	2.59 ng	51667	Perylene-d12	15.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	43
2		N-(3,4,5,6-Tetrahydro-1,3-thiazin...	192	C10H12N2S	003420-40-4	38
3		2H-1,2,3-Triazole-4-carboxaldehv...	191	C9H6FN3O	051306-43-5	38
4		7a-Isopropenyl-4,5-dimethyloctah...	236	C15H24O2	1000192-14-7	37
5		3-Methylthio-2-quinoxalinamine	191	C9H9N3S	090349-86-3	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062718\
Data File : BF106877.D
Acq On : 27 Jun 2018 17:35
Operator : JU/SJ
Sample : J3570-04
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Instrument :
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ClientSampleId :
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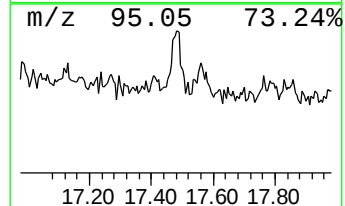
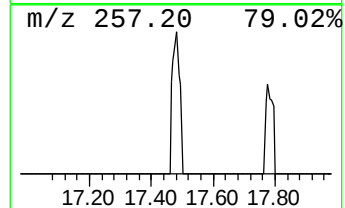
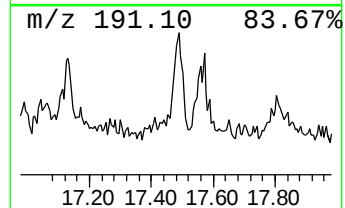
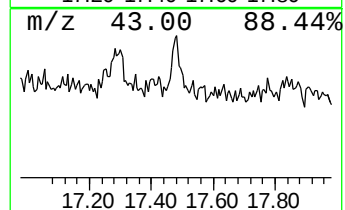
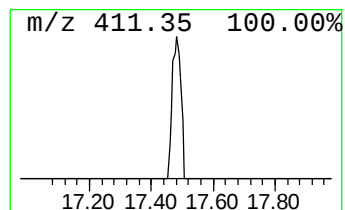
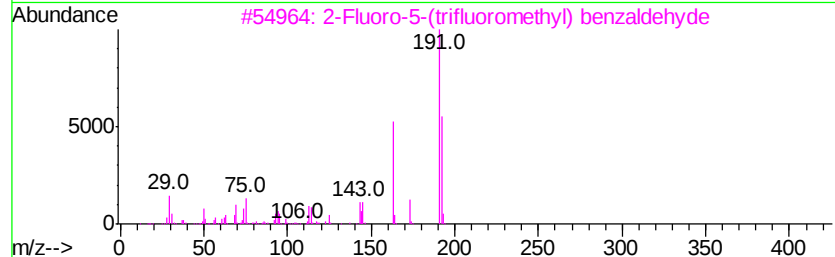
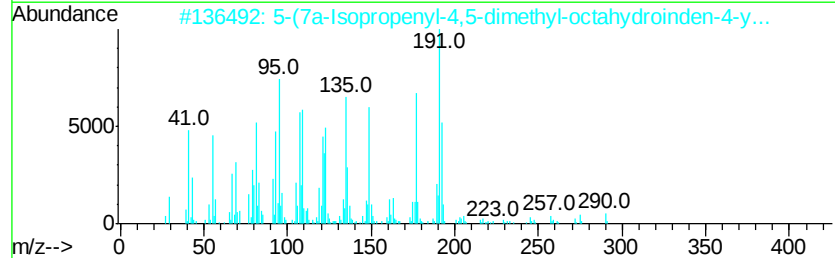
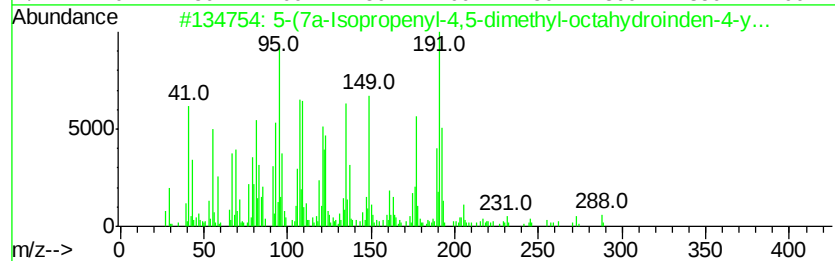
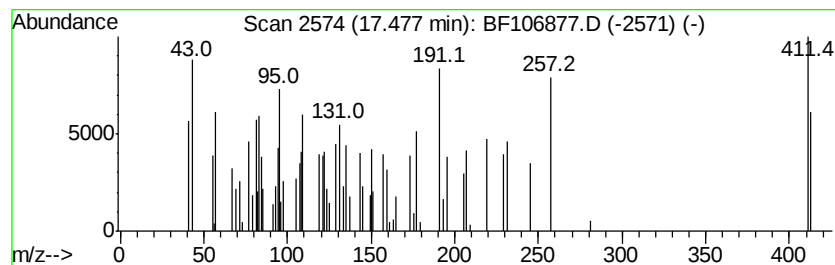
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown17.48 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.48	2.45 ng	48975	Perylene-d12	15.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-(7a-Isopropenyl-4,5-dimethyl-o...	288	C20H32O	1000193-54-1	25
2		5-(7a-Isopropenyl-4,5-dimethyl-o...	290	C20H34O	1000193-54-0	16
3		2-Fluoro-5-(trifluoromethyl) ben...	192	C8H4F4O	1000342-45-2	11
4		7-Hydroxy-2,5,8-quinoline trione	191	C9H5NO4	015544-54-4	11
5		Propenamide, 3-(3-methoxyphenyl)...	191	C11H13NO2	1000259-85-9	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF062718\
Data File : BF106877.D
Acq On : 27 Jun 2018 17:35
Operator : JU/SJ
Sample : J3570-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
20180374-AMENDED-COMPOSITE-

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.23	178.3	ng	2268390	1	6.95	254436	20.0
unknown6.73	6.73	2.7	ng	34353	1	6.95	254436	20.0
Triethylene glycol	8.38	4.8	ng	78071	2	8.24	321693	20.0
1H-Indene, octahy...	9.38	2.5	ng	43100	3	10.00	343664	20.0
1-Trimethylsilylp...	10.40	2.3	ng	39063	3	10.00	343664	20.0
Tetradecane	10.88	2.8	ng	49259	4	11.49	349959	20.0
unknown11.35	11.35	2.4	ng	42130	4	11.49	349959	20.0
Pentafluoropropio...	13.95	4.4	ng	63843	5	14.14	292513	20.0
unknown16.87	16.87	2.6	ng	51667	6	15.68	399473	20.0
unknown17.48	17.48	2.5	ng	48975	6	15.68	399473	20.0