

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF101818\
 Data File : BF109974.D
 Acq On : 18 Oct 2018 22:19
 Operator : JU/SJ
 Sample : J5573-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FIELD-BLANK

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-BF101518.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	2.310	32	38	54	rVB	2452904	3306592	71.24%	8.164%
2	5.204	526	530	540	rVB	105843	113657	2.45%	0.281%
3	5.592	592	596	599	rBV	2252813	1929305	41.56%	4.763%
4	6.504	748	751	758	rBV	173190	198990	4.29%	0.491%
5	6.587	761	765	769	rVV	1329518	1188800	25.61%	2.935%
6	6.745	787	792	795	rBV	3559610	3496175	75.32%	8.632%
7	6.975	827	831	834	rBV	1468040	1130594	24.36%	2.791%
8	7.134	854	858	861	rVB	3824587	3287206	70.82%	8.116%
9	7.539	921	927	930	rBV	2696662	2726285	58.73%	6.731%
10	7.751	959	963	964	rBV2	57768	58530	1.26%	0.145%
11	7.775	964	967	974	rVV	93253	96445	2.08%	0.238%
12	7.851	977	980	985	rBV	70370	57084	1.23%	0.141%
13	8.257	1045	1049	1052	rBV	1841141	1435196	30.92%	3.543%
14	8.533	1093	1096	1100	rBV	288471	237044	5.11%	0.585%
15	8.745	1128	1132	1135	rBV	226632	184198	3.97%	0.455%
16	8.875	1151	1154	1163	rVB	198269	151778	3.27%	0.375%
17	9.333	1227	1232	1235	rBV	5506817	4641681	100.00%	11.460%
18	10.016	1343	1348	1352	rBV2	1999481	1966722	42.37%	4.856%
19	10.804	1478	1482	1486	rBV	2743649	2742402	59.08%	6.771%
20	10.898	1495	1498	1503	rBV	77334	67554	1.46%	0.167%
21	11.510	1598	1602	1605	rBV	1701052	1509670	32.52%	3.727%
22	11.716	1633	1637	1641	rBV	74329	65766	1.42%	0.162%
23	12.533	1773	1776	1781	rBV	304627	243618	5.25%	0.601%
24	12.904	1835	1839	1842	rBV	103386	99617	2.15%	0.246%
25	13.092	1867	1871	1874	rBV	4399361	4033654	86.90%	9.959%
26	13.563	1944	1951	1952	rBV	123126	122357	2.64%	0.302%
27	13.580	1952	1954	1961	rVV2	136477	232063	5.00%	0.573%
28	13.639	1961	1964	1973	rVB	80952	91570	1.97%	0.226%
29	14.151	2047	2051	2055	rBV	1321783	1130626	24.36%	2.791%
30	14.604	2124	2128	2133	rVB	152153	146025	3.15%	0.361%
31	14.921	2176	2182	2187	rBV2	322081	509409	10.97%	1.258%
32	15.004	2193	2196	2200	rVB	58136	60347	1.30%	0.149%
33	15.280	2239	2243	2254	rVB	395484	455007	9.80%	1.123%
34	15.680	2305	2311	2318	rBV2	1209966	1710280	36.85%	4.222%

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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	16.145	2384	2390	2405	rBV	360281	597213	12.87%	1.474%
36	16.686	2476	2482	2495	rVB	194702	362270	7.80%	0.894%
37	17.321	2585	2590	2597	rVB2	60353	118433	2.55%	0.292%

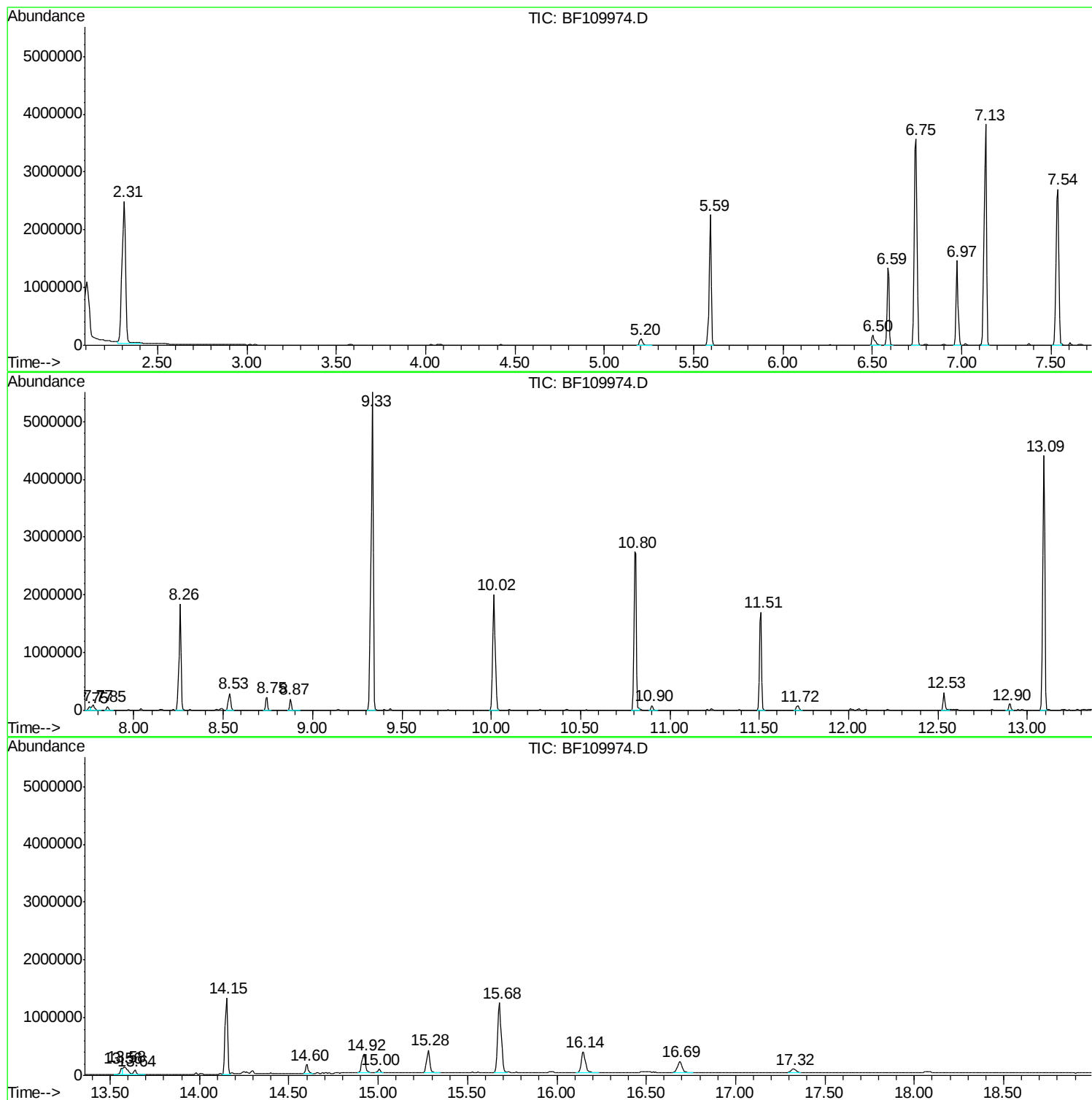
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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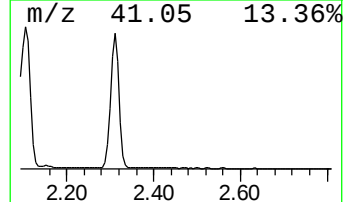
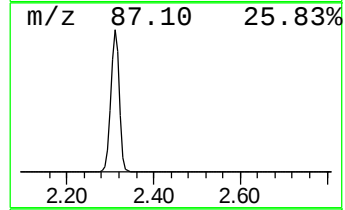
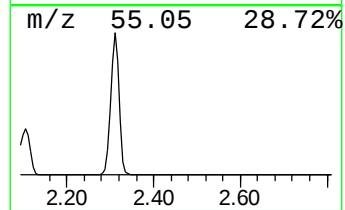
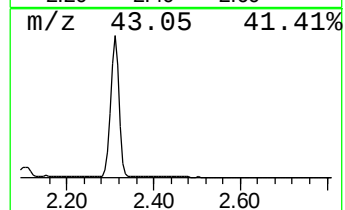
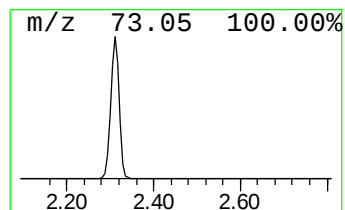
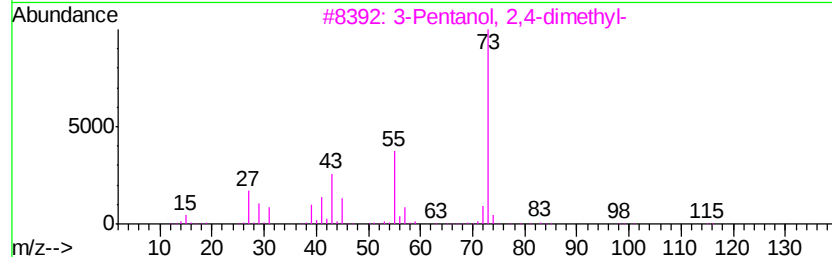
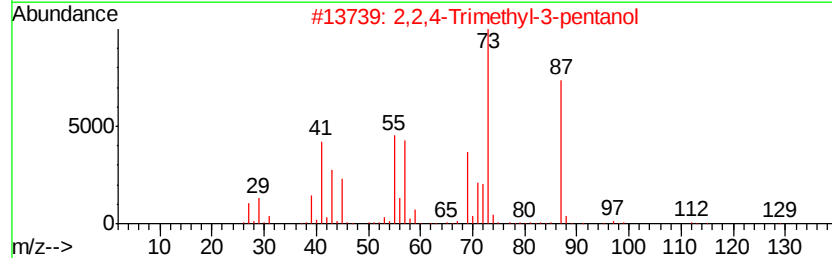
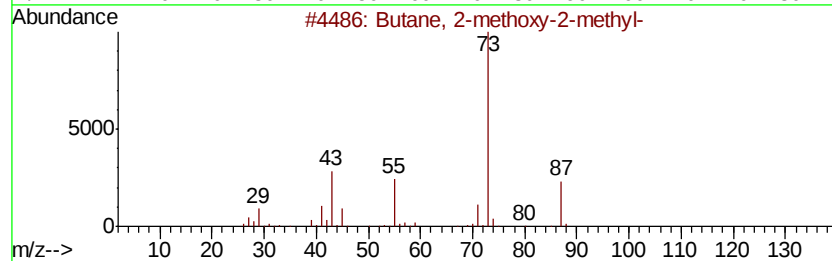
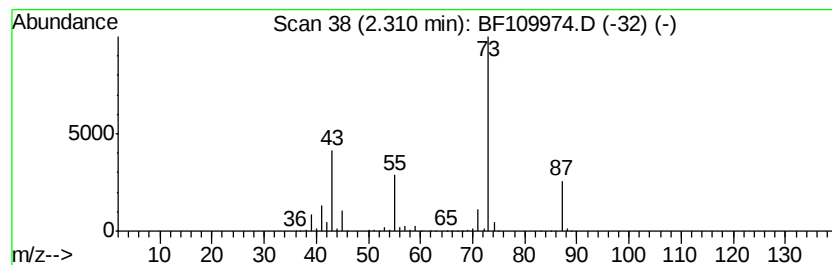
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.31	58.49 ng	3306590	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	27
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25



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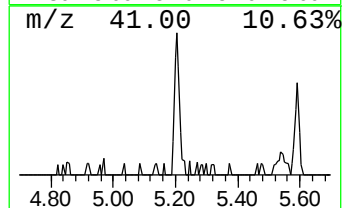
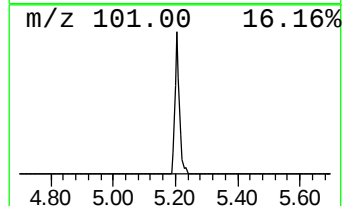
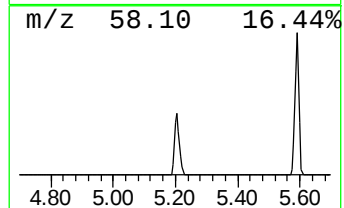
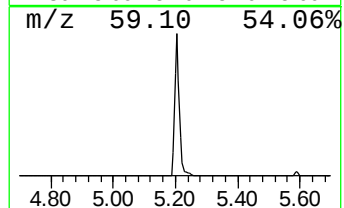
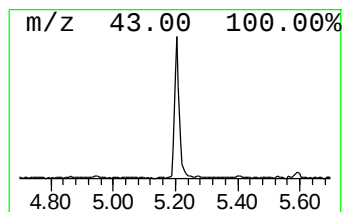
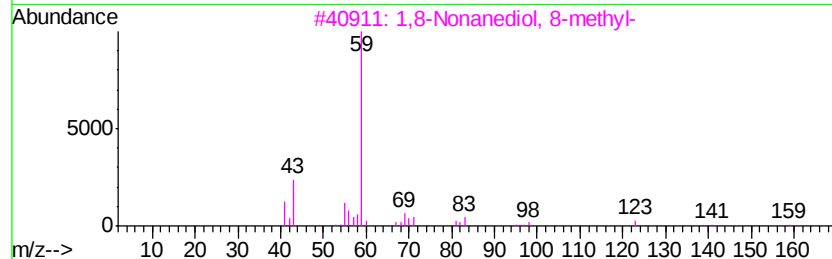
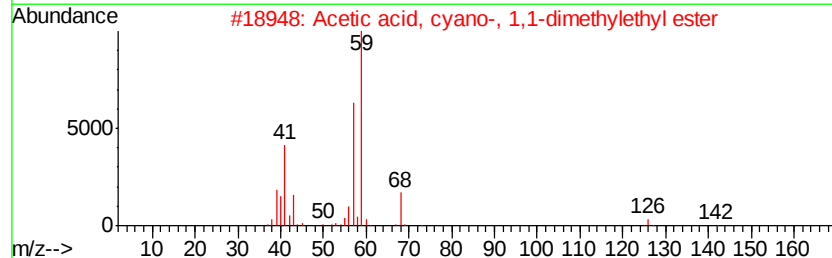
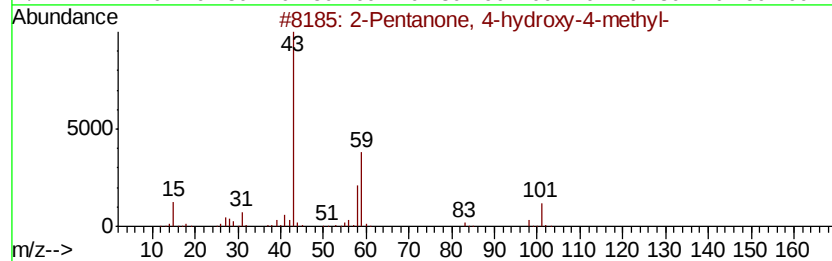
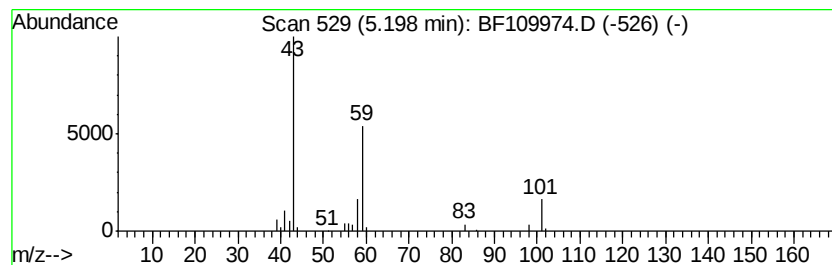
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	2.01 ng	113657	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	32
3			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	23
4			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	17
5			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	12



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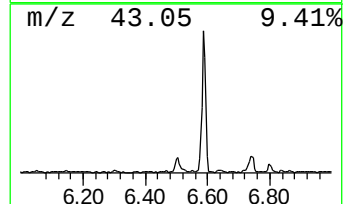
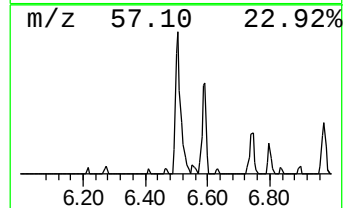
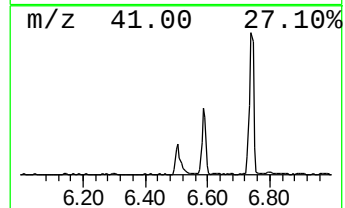
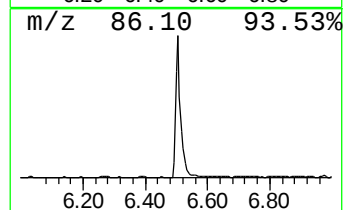
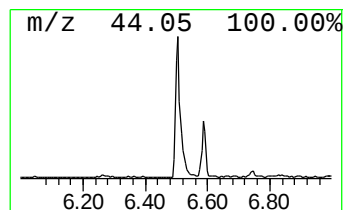
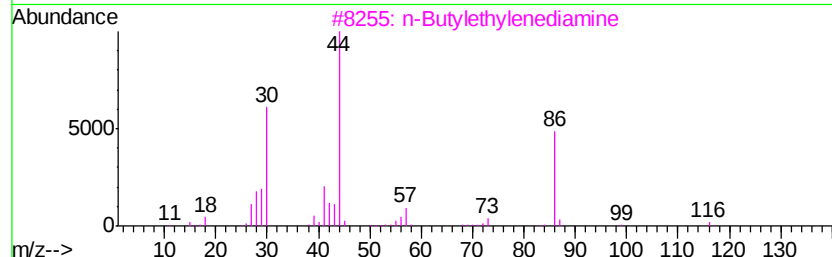
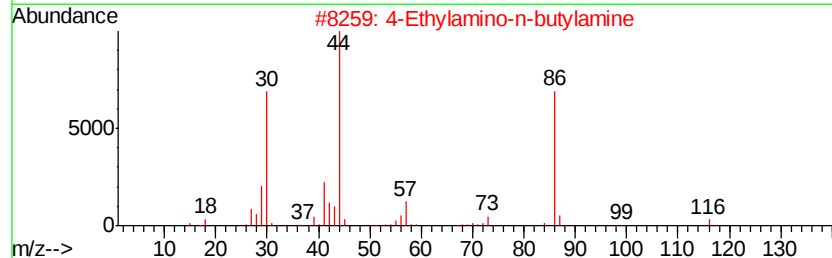
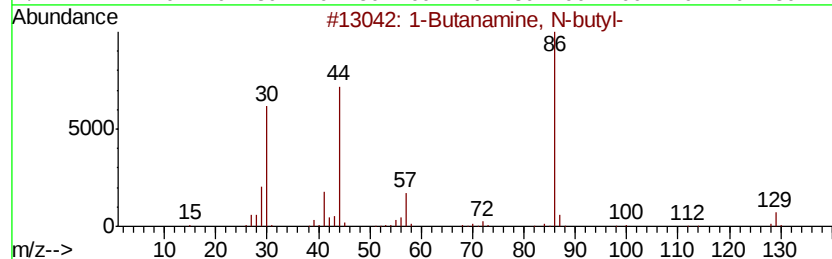
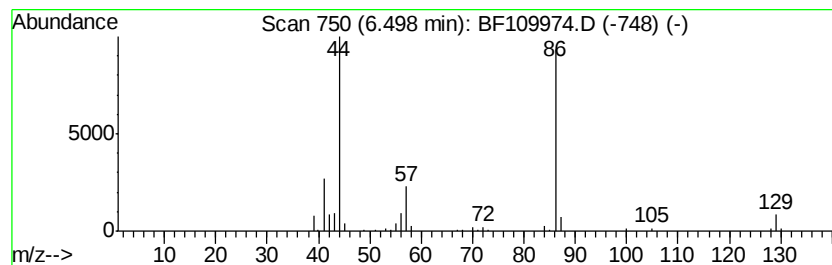
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1-Butanamine, N-butyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	3.52 ng	198990	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Butanamine, N-butyl-	129	C8H19N	000111-92-2	91
2			4-Ethylamino-n-butylamine	116	C6H16N2	064429-16-9	83
3			n-Butylethylenediamine	116	C6H16N2	019522-69-1	83
4			Prilocaine	220	C13H20N2O	000721-50-6	64
5			1-Propanamine, 2-methyl-N-(2-met...	129	C8H19N	000110-96-3	46



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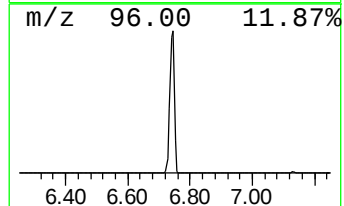
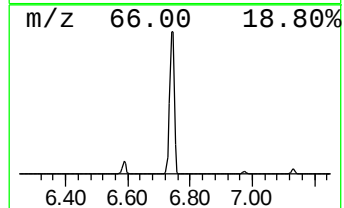
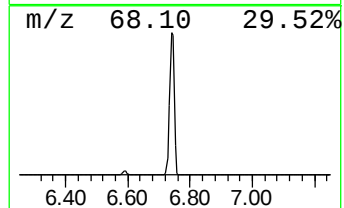
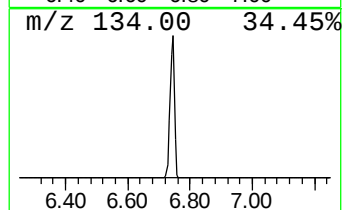
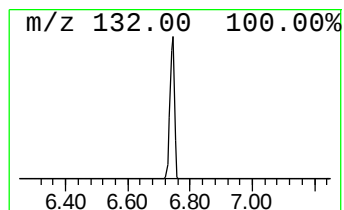
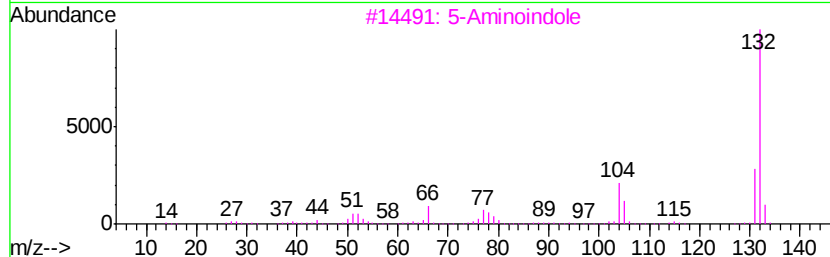
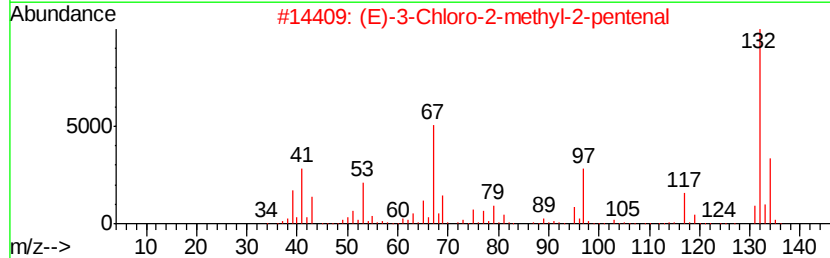
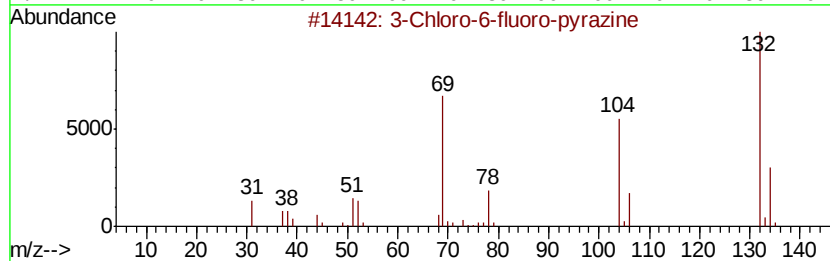
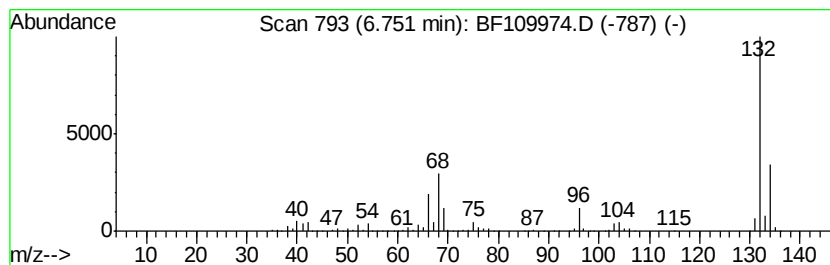
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown6.75 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	61.85 ng	3496180	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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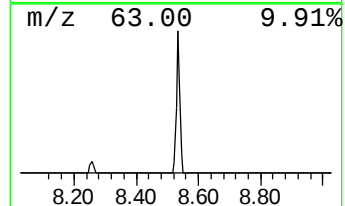
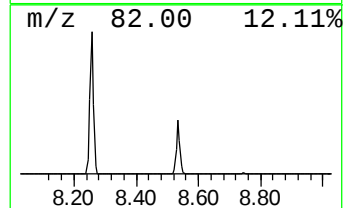
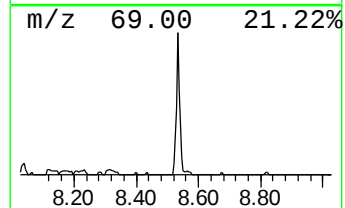
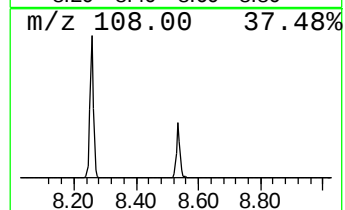
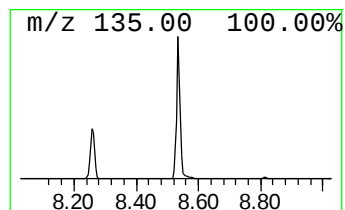
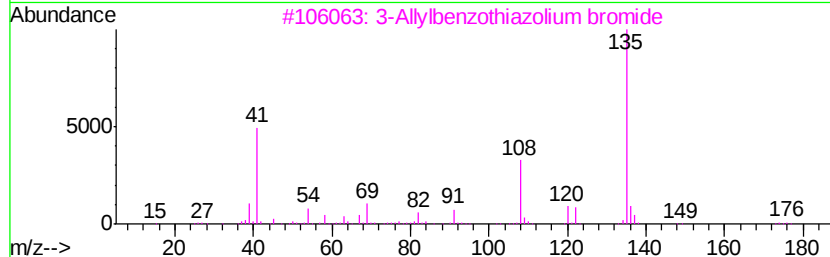
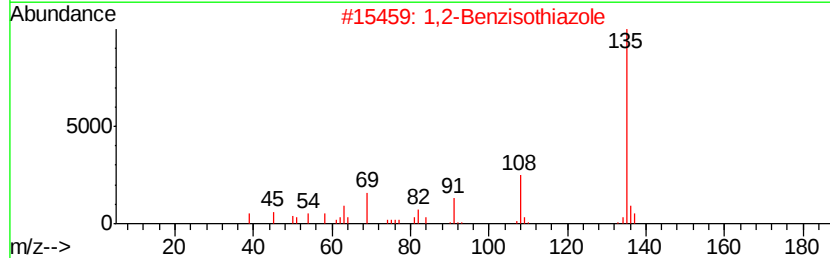
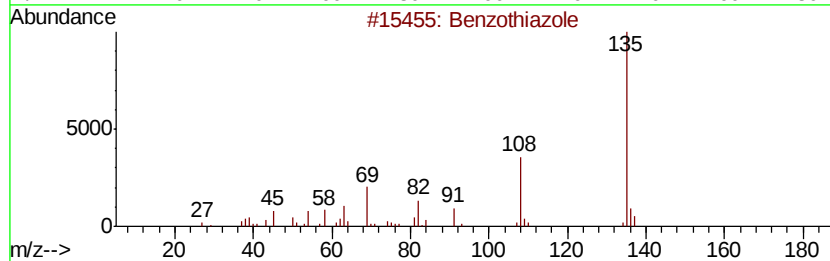
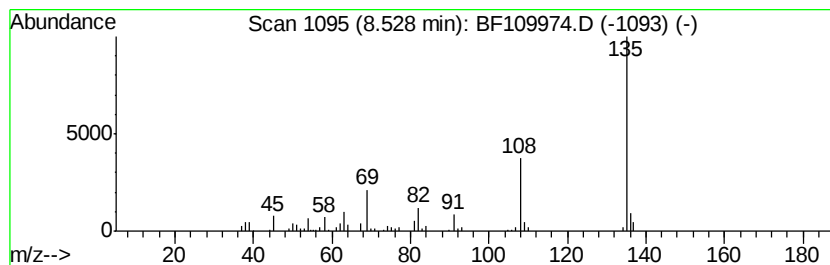
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzothiazole Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.53	3.30 ng	237044	Naphthalene-d8	8.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzothiazole	135	C7H5NS	000095-16-9	97
2			1,2-Benzisothiazole	135	C7H5NS	000272-16-2	91
3			3-Allylbenzothiazolium bromide	255	C10H10BrNS	016407-55-9	90
4			1H-Pyrazolo[3,4-d]pyrimidin-4-amine	135	C5H5N5	002380-63-4	64
5			Thieno[2,3-c]pyridine	135	C7H5NS	000272-12-8	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
Data File : BF109974.D
Acq On : 18 Oct 2018 22:19
Operator : JU/SJ
Sample : J5573-04
Misc :
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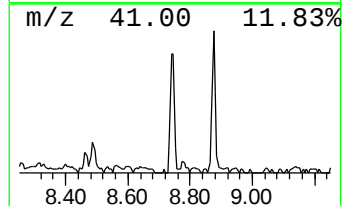
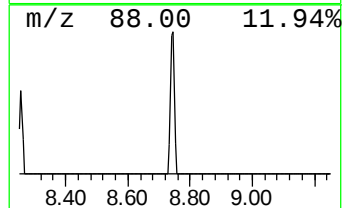
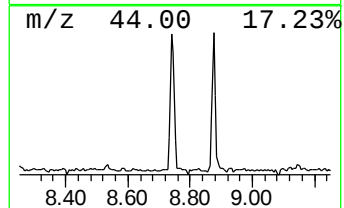
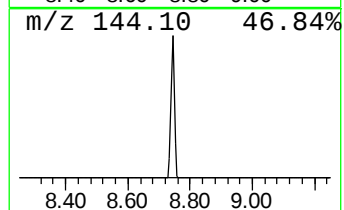
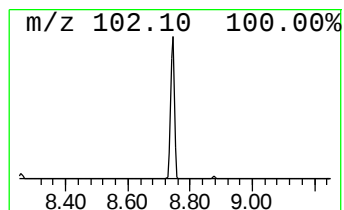
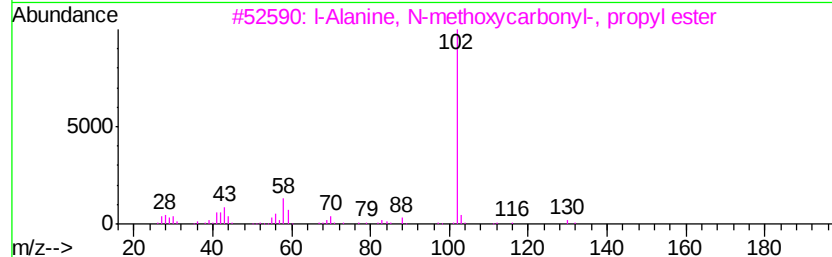
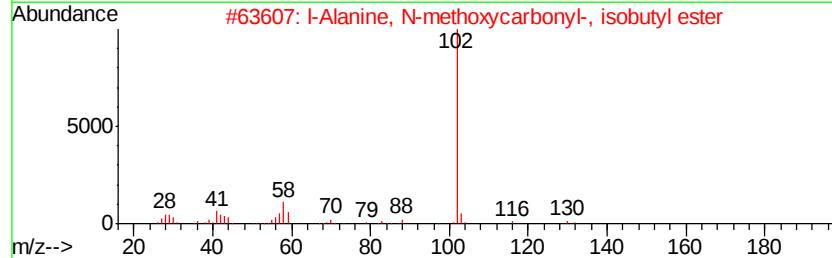
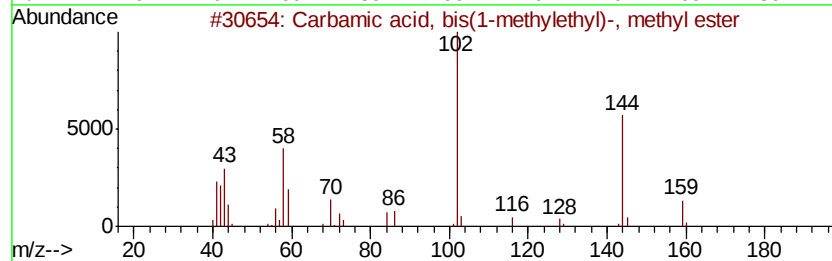
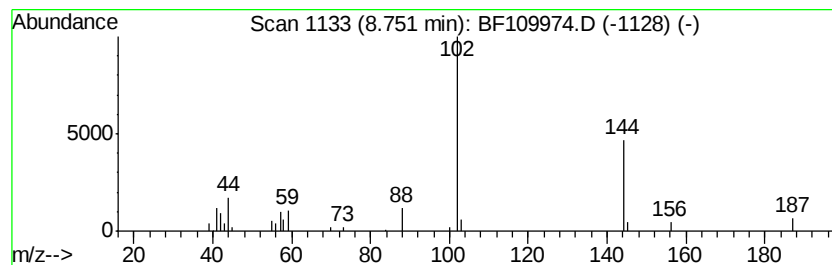
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown8.75 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.75	2.57 ng	184198	Naphthalene-d8	8.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Carbamic acid, bis(1-methylethyl...	159	C8H17N02	031603-49-3	42
2		l-Alanine, N-methoxycarbonyl-, i...	203	C9H17N04	1000313-37-2	40
3		l-Alanine, N-methoxycarbonyl-, p...	189	C8H15N04	1000313-37-1	28
4		l-Alanine, N-methoxycarbonyl-, h...	245	C12H23N04	1000313-37-7	28
5		Carbamic acid, butylmethyl-, met...	145	C7H15N02	054644-60-9	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
Data File : BF109974.D
Acq On : 18 Oct 2018 22:19
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Sample : J5573-04
Misc :
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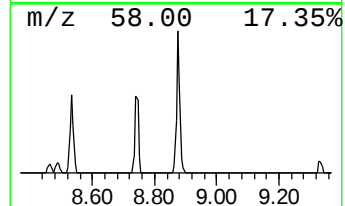
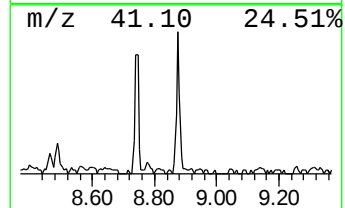
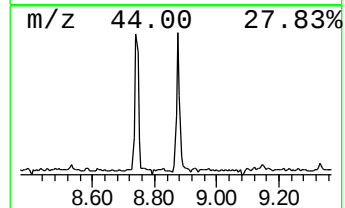
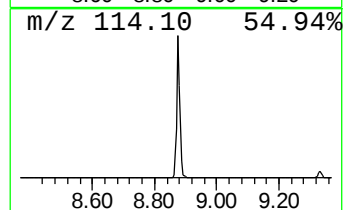
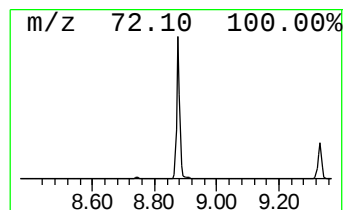
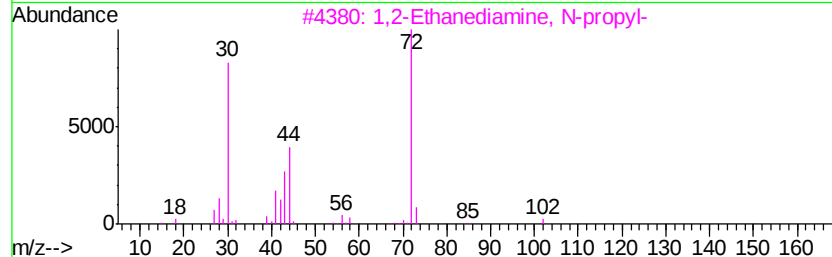
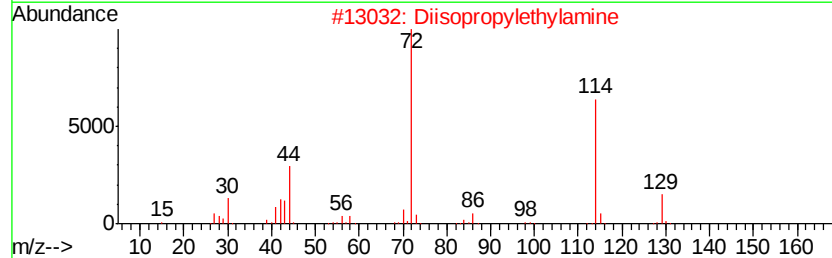
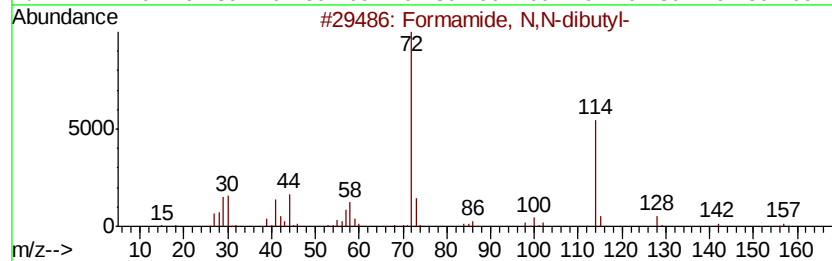
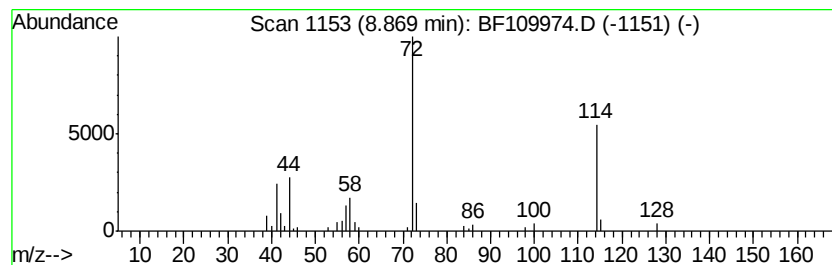
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Formamide, N,N-dibutyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.87	2.12 ng	151778	Naphthalene-d8	8.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Formamide, N,N-dibutyl-	157	C9H19NO	000761-65-9	94
2		Diisopropylethylamine	129	C8H19N	007087-68-5	59
3		1,2-Ethanediamine, N-propyl-	102	C5H14N2	000111-39-7	47
4		2-Pentanamine, N-ethyl-4-methyl-	129	C8H19N	042966-64-3	47
5		Dimethyl(2-octyl)amine	157	C10H23N	007378-97-4	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
Data File : BF109974.D
Acq On : 18 Oct 2018 22:19
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Sample : J5573-04
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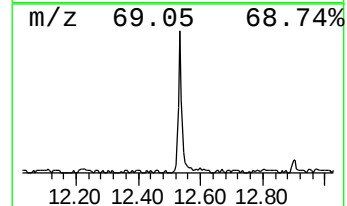
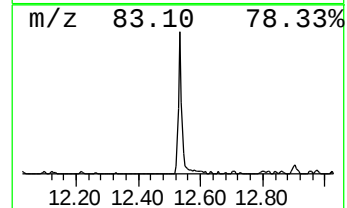
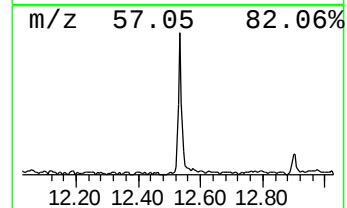
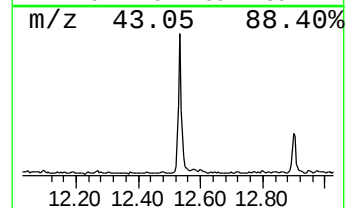
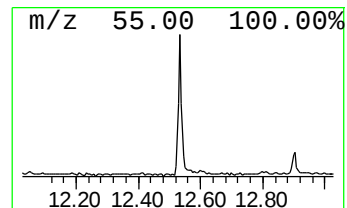
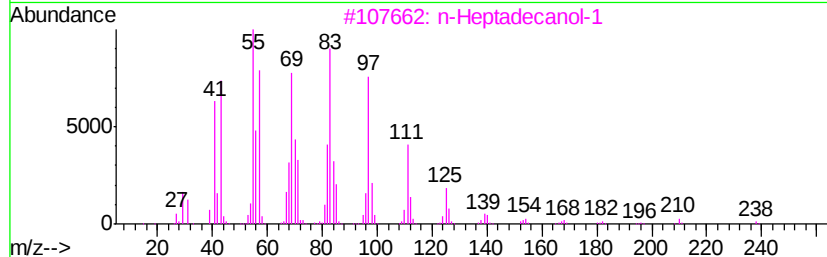
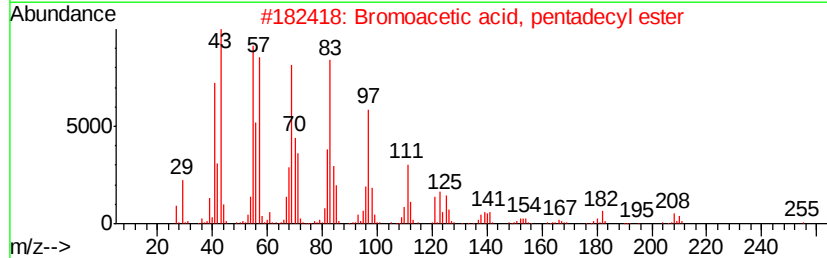
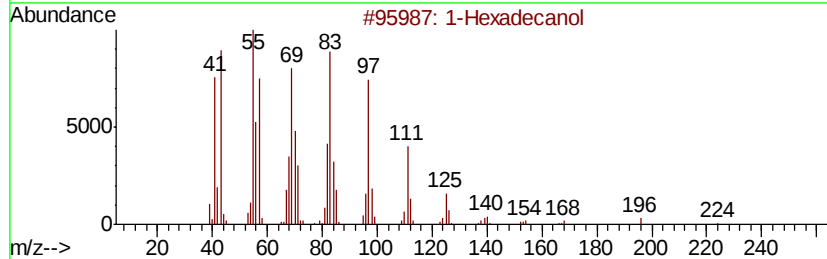
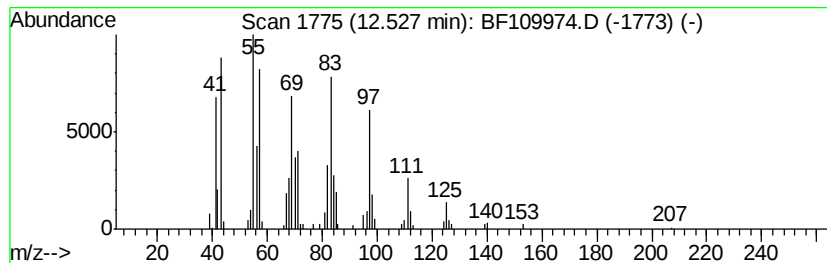
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Hexadecanol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.53	3.23 ng	243618	Phenanthrene-d10	11.51

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Hexadecanol	242	C16H34O	036653-82-4	95
2		Bromoacetic acid, pentadecyl ester	348	C17H33BrO2	131143-01-6	94
3		n-Heptadecanol-1	256	C17H36O	001454-85-9	94
4		1-Nonadecene	266	C19H38	018435-45-5	91
5		5-Octadecene, (E)-	252	C18H36	007206-21-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
Data File : BF109974.D
Acq On : 18 Oct 2018 22:19
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Sample : J5573-04
Misc :
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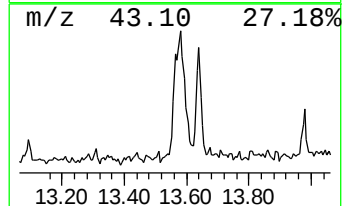
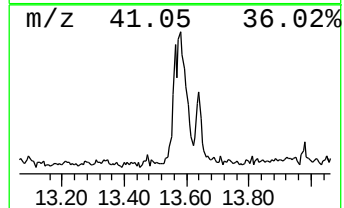
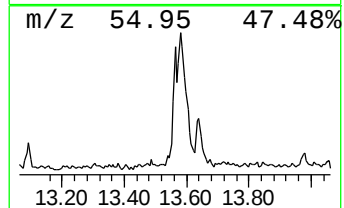
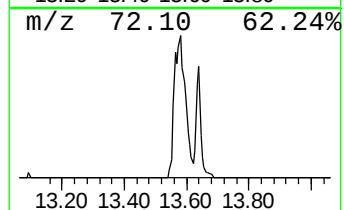
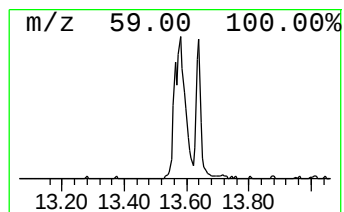
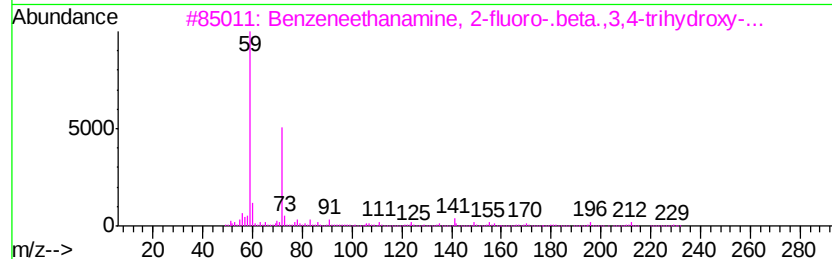
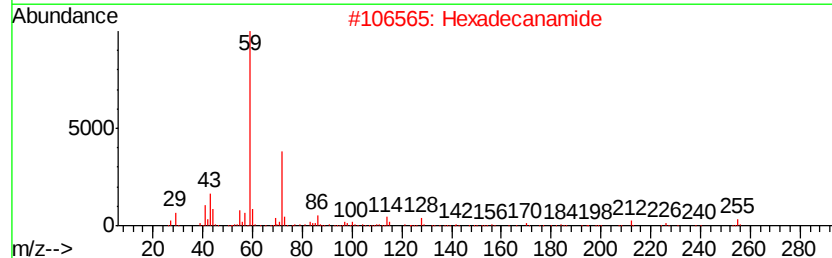
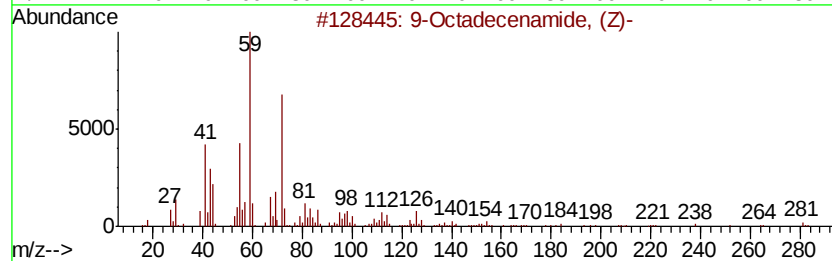
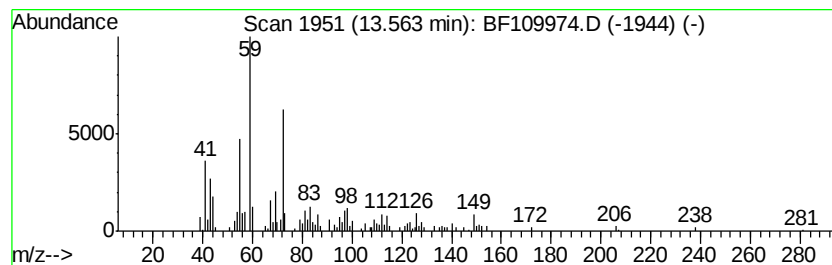
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 9-Octadecenamide, (Z)- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.56	2.16 ng	122357	Chrysene-d12	14.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	91
2		Hexadecanamide	255	C16H33NO	000629-54-9	64
3		Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FN03	061338-98-5	59
4		Tetradecanamide	227	C14H29NO	000638-58-4	53
5		Dodecanamide	199	C12H25NO	001120-16-7	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
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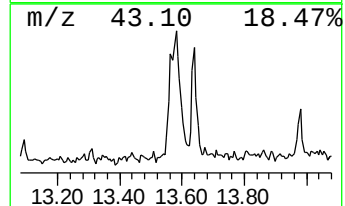
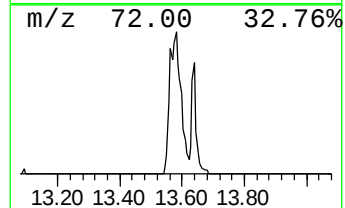
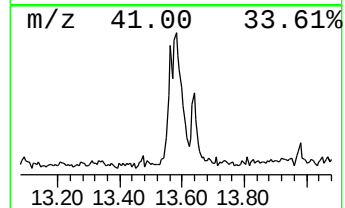
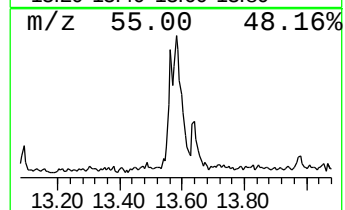
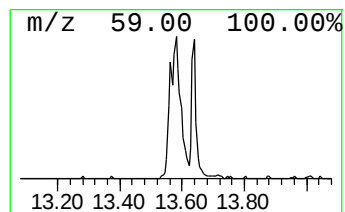
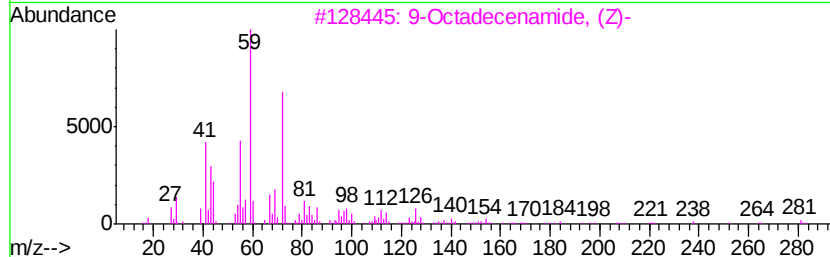
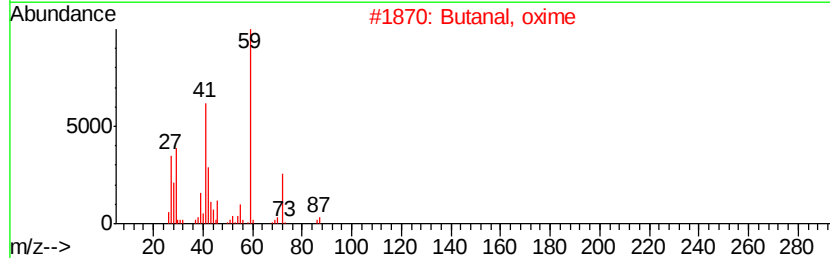
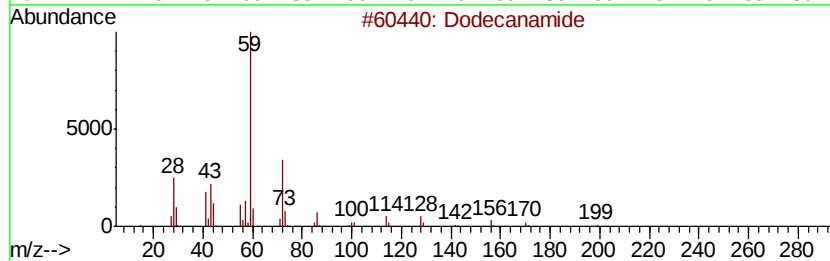
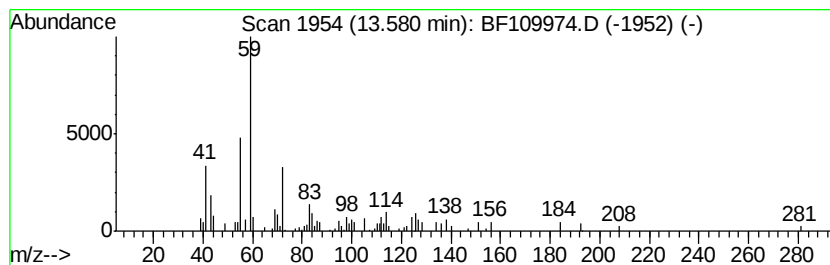
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Dodecanamide Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.58	4.11 ng	232063	Chrysene-d12	14.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecanamide	199	C12H25NO	001120-16-7	53
2			Butanal, oxime	87	C4H9NO	000110-69-0	50
3			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	45
4			Succinic acid, 2-methoxyethyl oc...	288	C15H28O5	1000325-80-8	43
5			Tetradecanamide	227	C14H29NO	000638-58-4	42



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
Data File : BF109974.D
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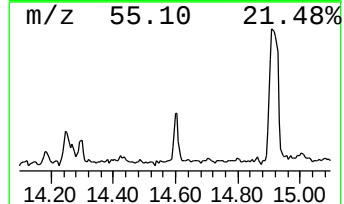
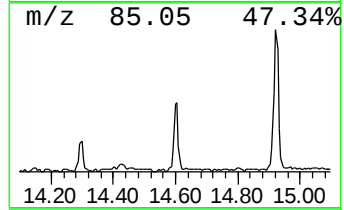
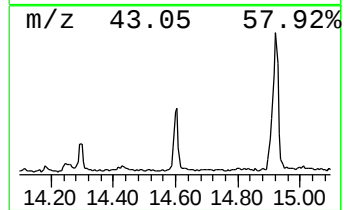
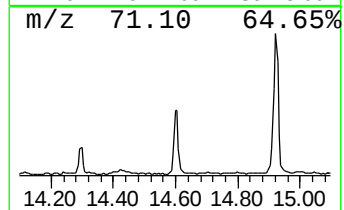
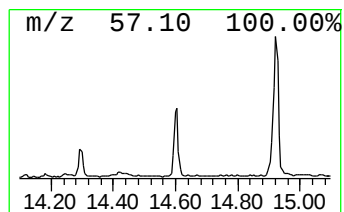
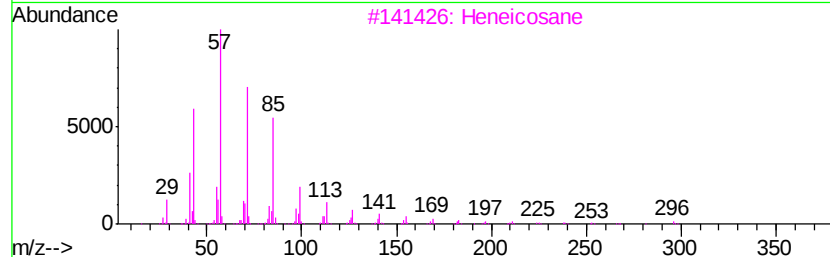
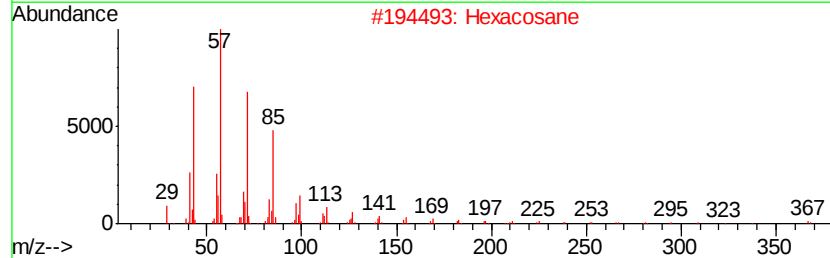
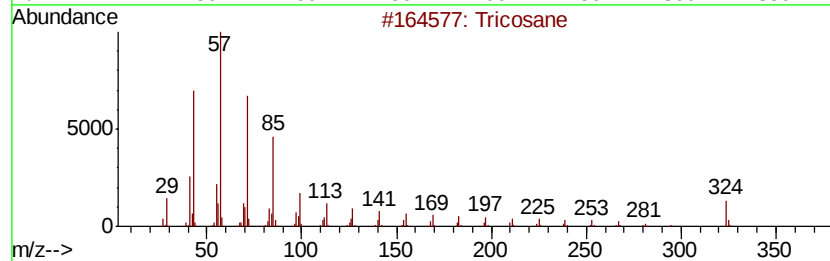
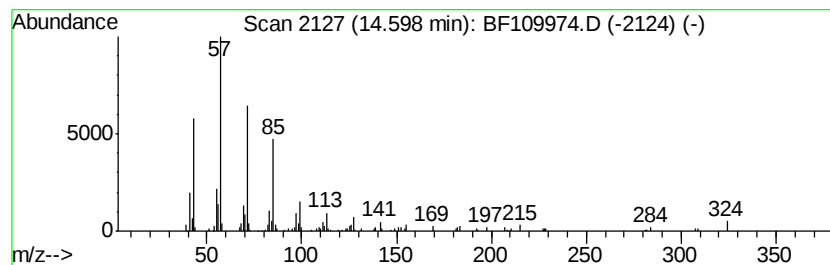
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tricosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	2.58 ng	146025	Chrysene-d12	14.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tricosane	324	C23H48	000638-67-5	93
2			Hexacosane	366	C26H54	000630-01-3	90
3			Heneicosane	296	C21H44	000629-94-7	90
4			Octacosane	394	C28H58	000630-02-4	90
5			Tetracosane	338	C24H50	000646-31-1	90



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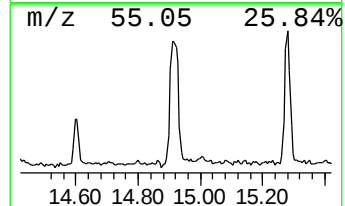
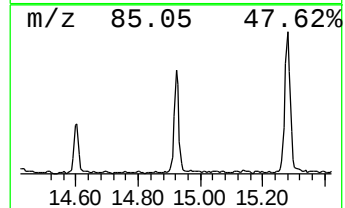
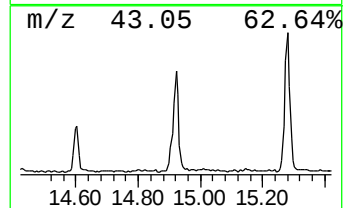
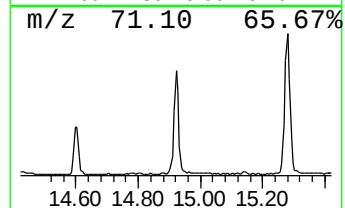
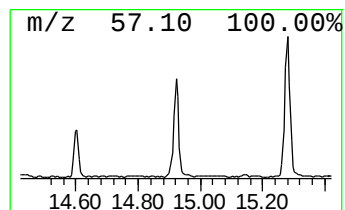
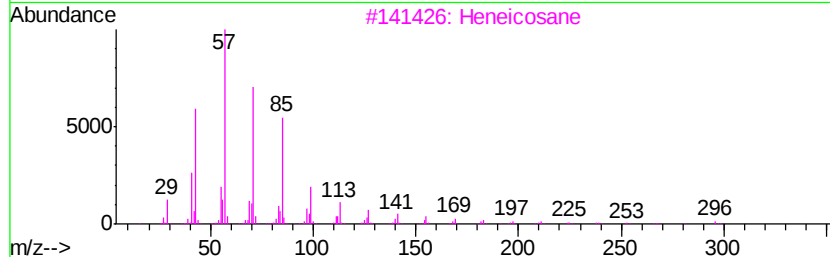
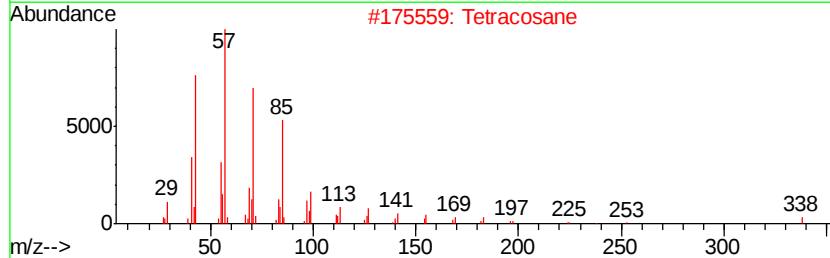
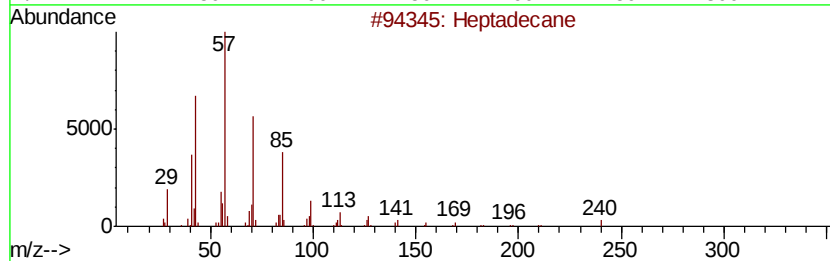
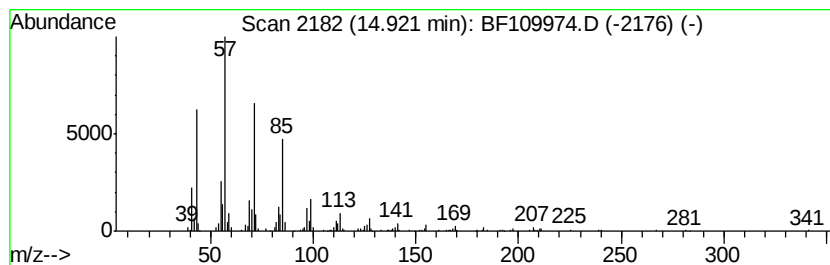
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heptadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.92	5.96 ng	509409	Perylene-d12	15.68

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	96
2	Tetracosane		338	C24H50	000646-31-1	91
3	Heneicosane		296	C21H44	000629-94-7	91
4	Octacosane		394	C28H58	000630-02-4	91
5	Eicosane		282	C20H42	000112-95-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
Data File : BF109974.D
Acq On : 18 Oct 2018 22:19
Operator : JU/SJ
Sample : J5573-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

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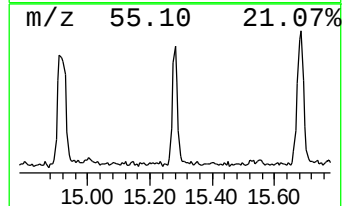
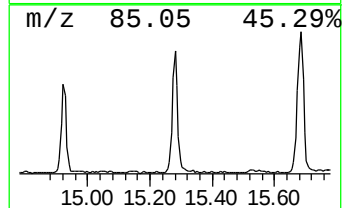
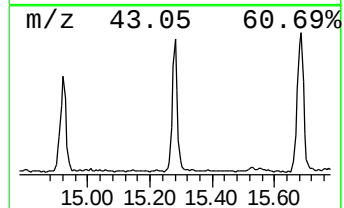
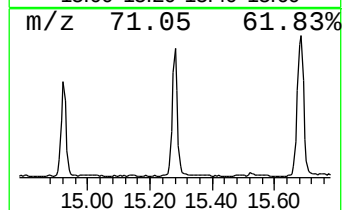
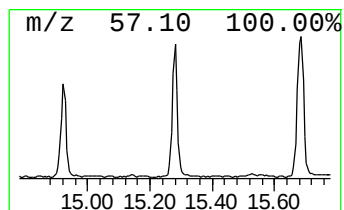
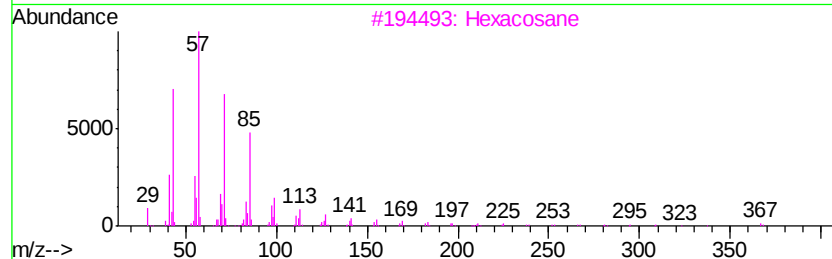
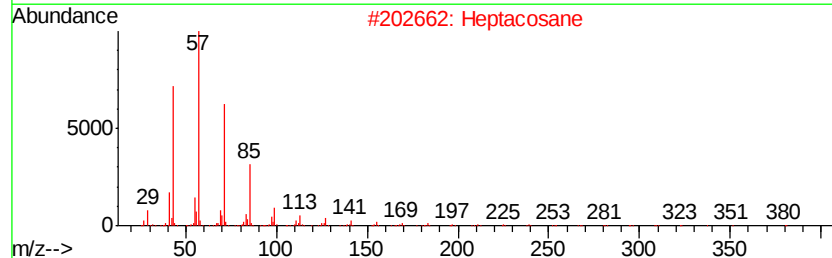
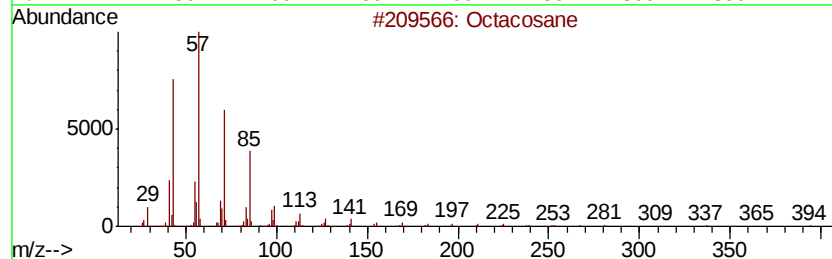
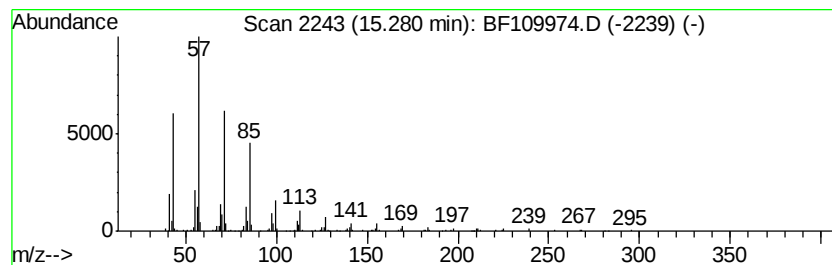
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Octacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.28	5.32 ng	455007	Perylene-d12	15.68

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	91
2			Heptacosane	380	C27H56	000593-49-7	91
3			Hexacosane	366	C26H54	000630-01-3	91
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5			2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
Data File : BF109974.D
Acq On : 18 Oct 2018 22:19
Operator : JU/SJ
Sample : J5573-04
Misc :
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Instrument :
BNA_F
ClientSampleId :
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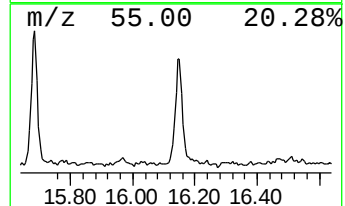
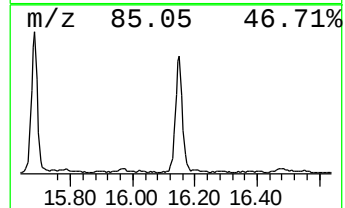
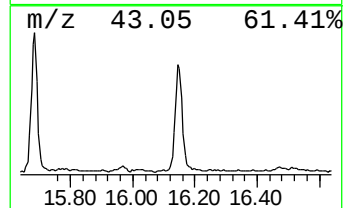
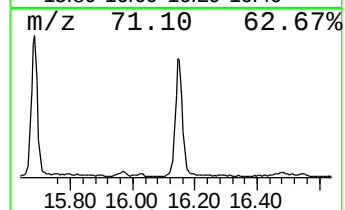
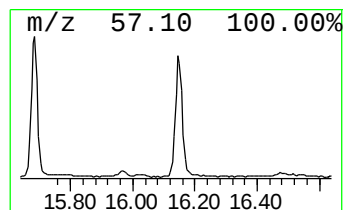
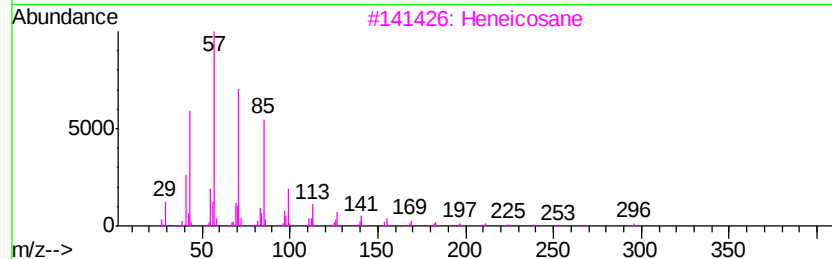
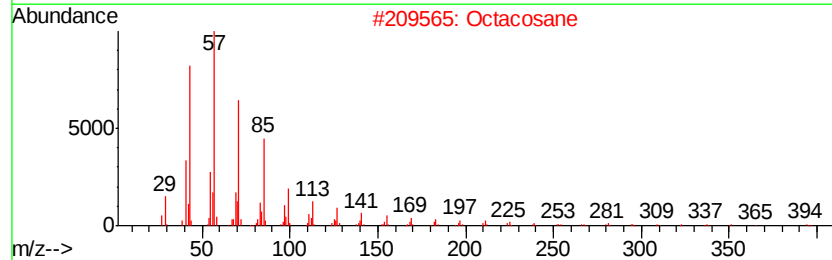
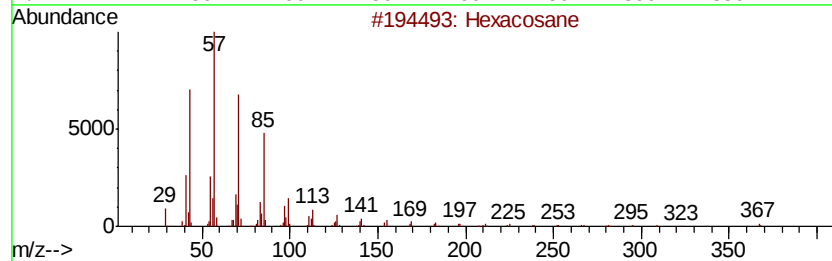
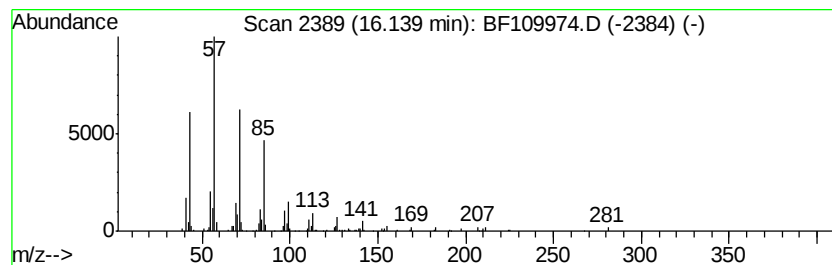
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Hexacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.14	6.98 ng	597213	Perylene-d12	15.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Octacosane	394	C28H58	000630-02-4	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Tetratetracontane	619	C44H90	007098-22-8	91
5		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
Data File : BF109974.D
Acq On : 18 Oct 2018 22:19
Operator : JU/SJ
Sample : J5573-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

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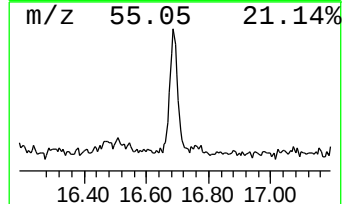
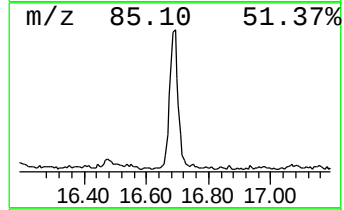
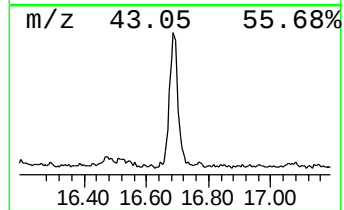
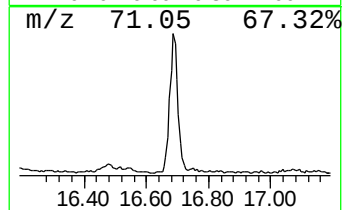
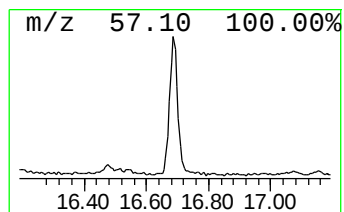
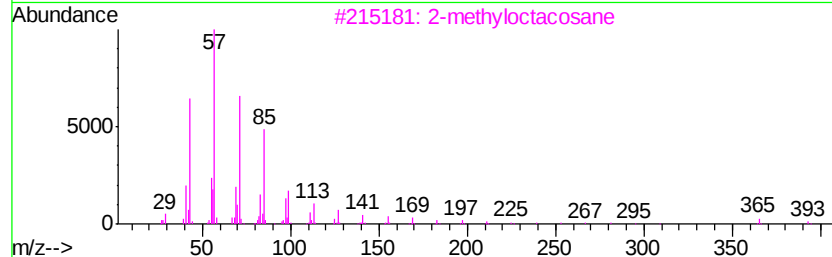
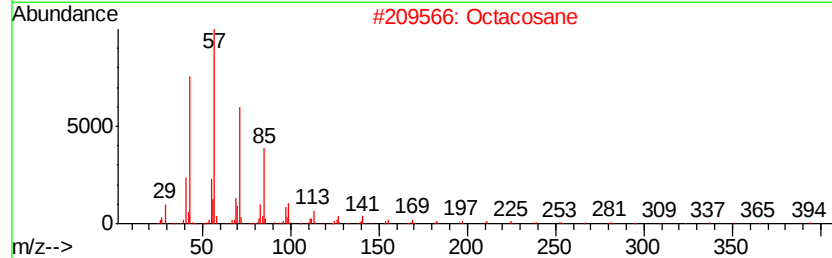
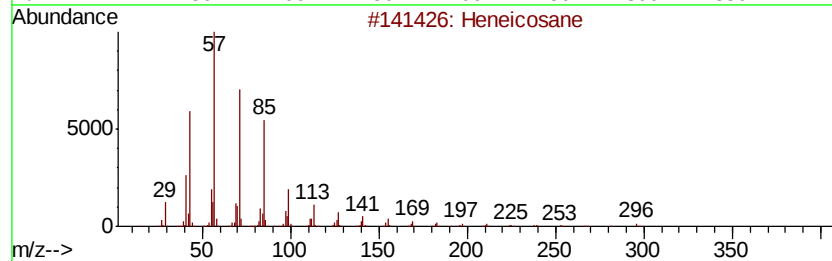
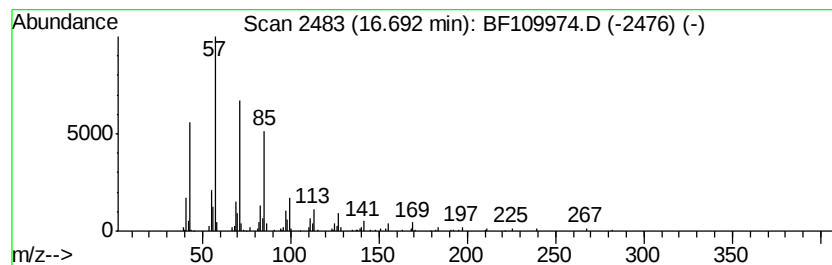
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Heneicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.69	4.24 ng	362270	Perylene-d12	15.68

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Octacosane	394	C28H58	000630-02-4	91
3			2-methyloctacosane	408	C29H60	1000376-72-8	91
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5			Hentriacontane	437	C31H64	000630-04-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF101818\
 Data File : BF109974.D
 Acq On : 18 Oct 2018 22:19
 Operator : JU/SJ
 Sample : J5573-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101518.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
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2-Pentanone, 4-hy...	5.20	2.0	ng	113657	1	6.97	1130590	20.0
1-Butanamine, N-b...	6.50	3.5	ng	198990	1	6.97	1130590	20.0
unknown6.75	6.75	61.9	ng	3496180	1	6.97	1130590	20.0
Benzothiazole	8.53	3.3	ng	237044	2	8.26	1435200	20.0
unknown8.75	8.75	2.6	ng	184198	2	8.26	1435200	20.0
Formamide, N,N-di...	8.87	2.1	ng	151778	2	8.26	1435200	20.0
1-Hexadecanol	12.53	3.2	ng	243618	4	11.51	1509670	20.0
9-Octadecenamide, ...	13.56	2.2	ng	122357	5	14.15	1130630	20.0
Dodecanamide	13.58	4.1	ng	232063	5	14.15	1130630	20.0
Tricosane	14.60	2.6	ng	146025	5	14.15	1130630	20.0
Heptadecane	14.92	6.0	ng	509409	6	15.68	1710280	20.0
Octacosane	15.28	5.3	ng	455007	6	15.68	1710280	20.0
Hexacosane	16.14	7.0	ng	597213	6	15.68	1710280	20.0
Heneicosane	16.69	4.2	ng	362270	6	15.68	1710280	20.0