

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061418\
 Data File : BF106440.D
 Acq On : 14 Jun 2018 14:23
 Operator : JU/SJ
 Sample : J3239-11
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-06-061318-B

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-BF061118.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	5.201	482	487	495	rVB	208085	246582	2.58%	0.540%
2	5.607	553	556	569	rBV	2237588	2508504	26.22%	5.493%
3	6.607	723	726	734	rBV	2482034	2536344	26.51%	5.553%
4	6.748	746	750	753	rVB	2599316	2457048	25.68%	5.380%
5	6.966	783	787	790	rBV	920443	739945	7.73%	1.620%
6	7.119	809	813	817	rVB	2143051	1990265	20.80%	4.358%
7	7.525	878	882	885	rBV	1741882	1686756	17.63%	3.693%
8	8.248	1001	1005	1008	rBV	1129956	933753	9.76%	2.045%
9	9.319	1183	1187	1191	rBV	3097406	2851431	29.80%	6.243%
10	9.713	1250	1254	1258	rBV	342069	299240	3.13%	0.655%
11	10.007	1300	1304	1307	rBV	1146434	1033251	10.80%	2.262%
12	10.419	1371	1374	1377	rVB	134999	108019	1.13%	0.237%
13	10.795	1434	1438	1442	rBV	1777006	1748237	18.27%	3.828%
14	10.895	1451	1455	1463	rBV	139322	164878	1.72%	0.361%
15	11.001	1469	1473	1476	rVB	294251	221918	2.32%	0.486%
16	11.495	1553	1557	1560	rBV	1201685	1011521	10.57%	2.215%
17	11.795	1606	1608	1613	rVB	405133	334794	3.50%	0.733%
18	11.877	1618	1622	1625	rBV	4574146	4906640	51.28%	10.743%
19	12.018	1640	1646	1652	rBV4	140902	200693	2.10%	0.439%
20	12.195	1674	1676	1679	rVB2	115061	102455	1.07%	0.224%
21	12.277	1687	1690	1695	rVB3	148028	132506	1.38%	0.290%
22	12.583	1731	1742	1743	rBV	5124214	9567789	100.00%	20.949%
23	12.671	1754	1757	1761	rVV	3025754	2889674	30.20%	6.327%
24	12.718	1761	1765	1769	rVV	299700	316386	3.31%	0.693%
25	12.901	1793	1796	1800	rBV2	92699	115473	1.21%	0.253%
26	12.954	1801	1805	1811	rVB	271578	351312	3.67%	0.769%
27	13.095	1824	1829	1832	rBV	2701405	2438040	25.48%	5.338%
28	13.336	1864	1870	1876	rVB2	473391	601585	6.29%	1.317%
29	13.413	1880	1883	1886	rVB	359749	314094	3.28%	0.688%
30	13.860	1955	1959	1967	rVB6	56961	115099	1.20%	0.252%
31	14.024	1984	1987	1995	rVB2	211978	252371	2.64%	0.553%
32	14.130	2002	2005	2011	rVB	262615	243620	2.55%	0.533%
33	14.212	2014	2019	2022	rVV2	1012837	1046958	10.94%	2.292%
34	15.359	2210	2214	2218	rBV	123174	204846	2.14%	0.449%

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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	15.754	2277	2281	2288	rvB	97001	134000	1.40%	0.293%
36	15.824	2288	2293	2297	rvB	663888	865413	9.05%	1.895%

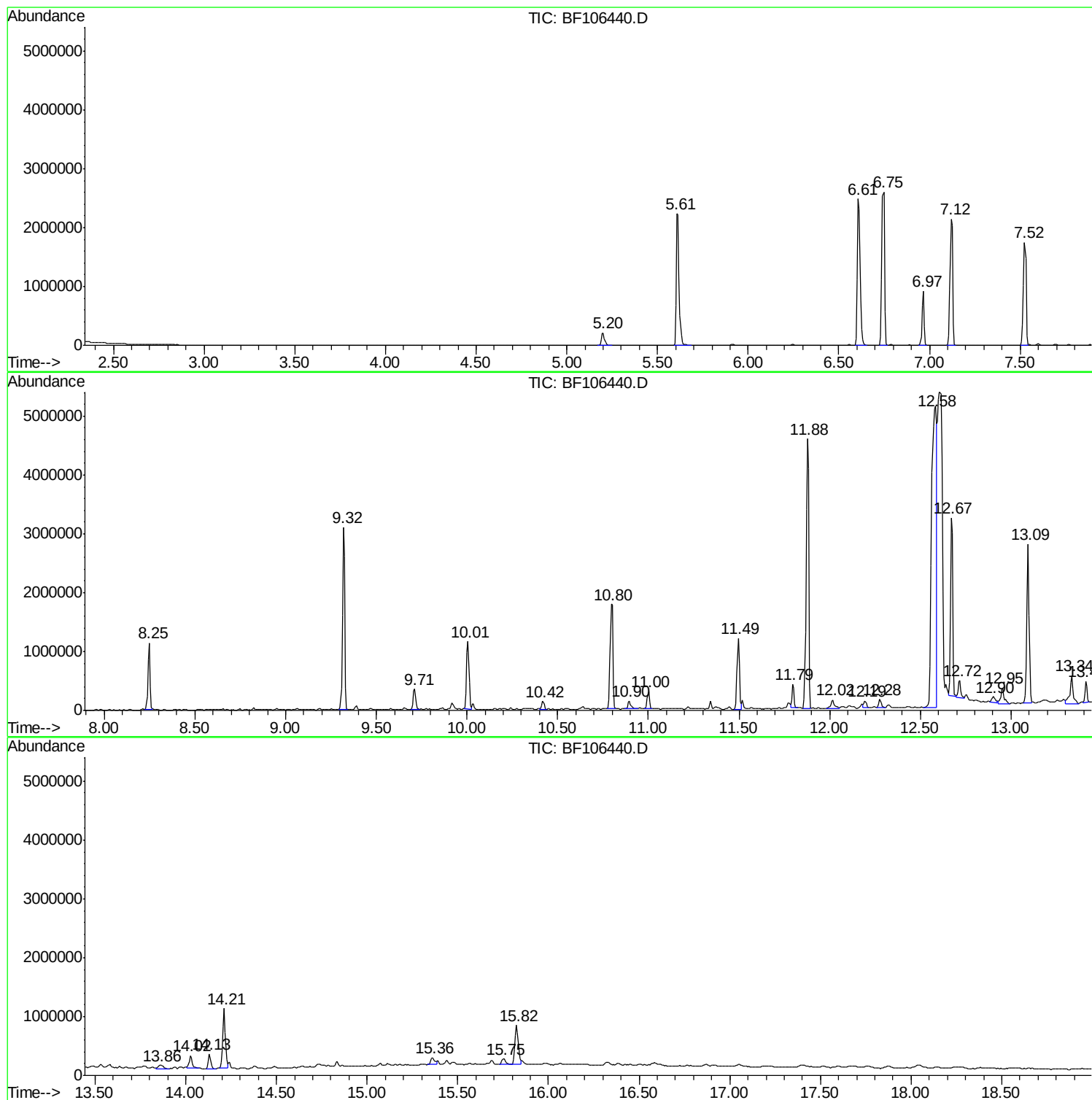
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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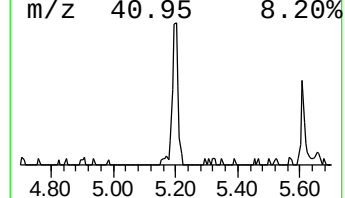
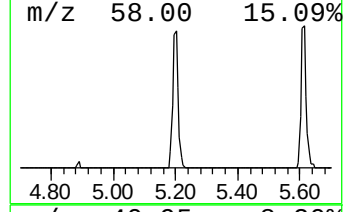
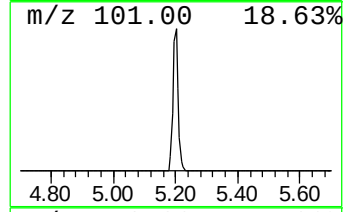
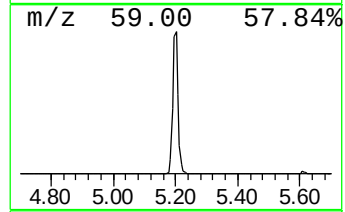
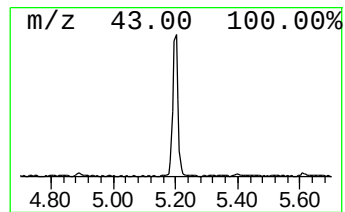
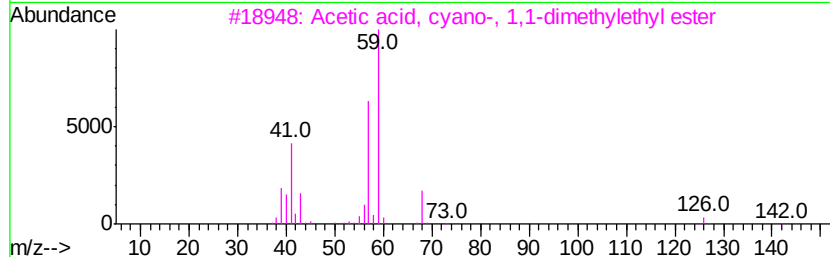
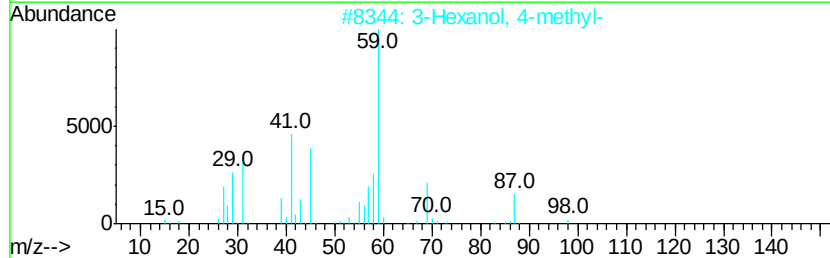
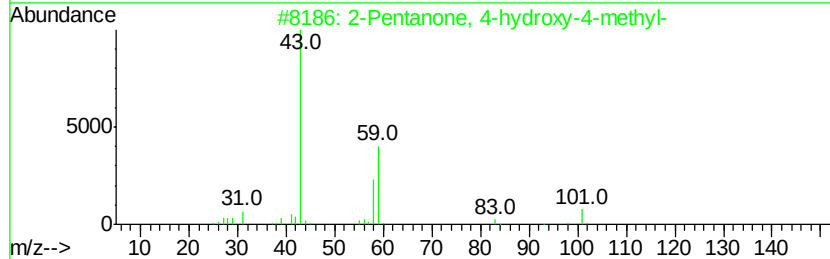
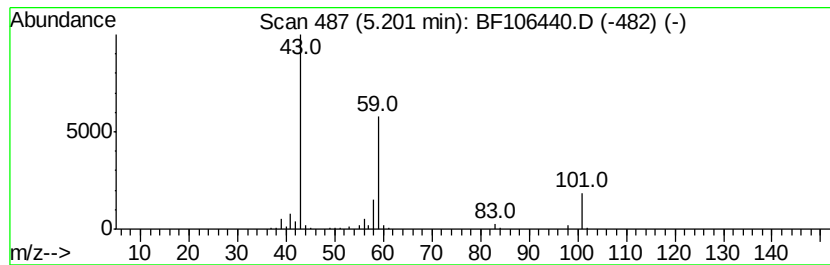
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown5.20 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	6.66 ng	246582	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	45
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5			2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9



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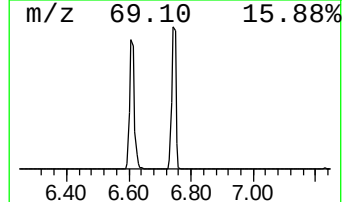
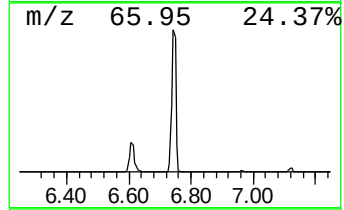
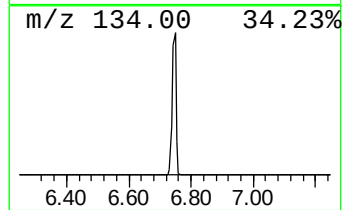
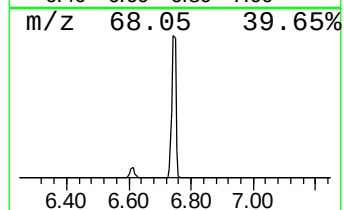
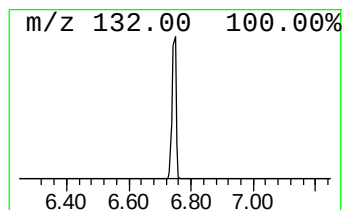
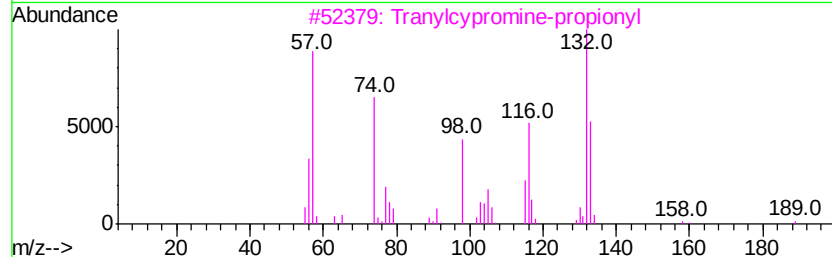
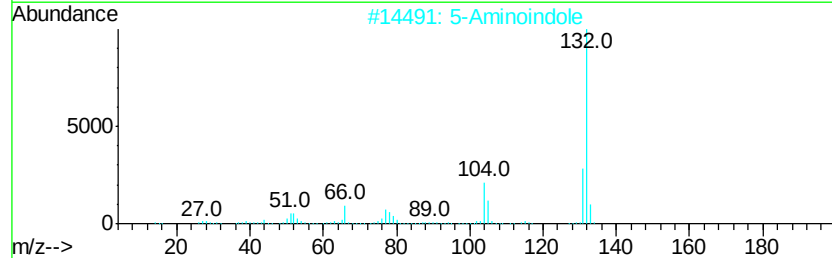
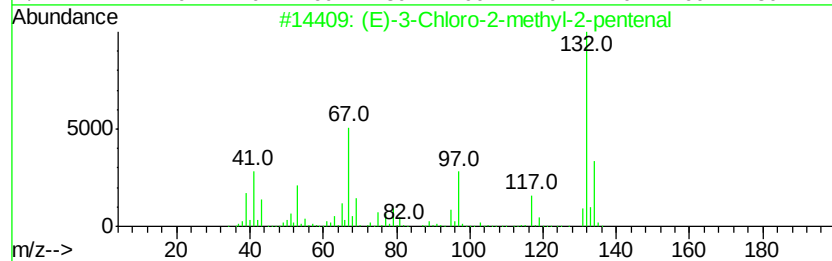
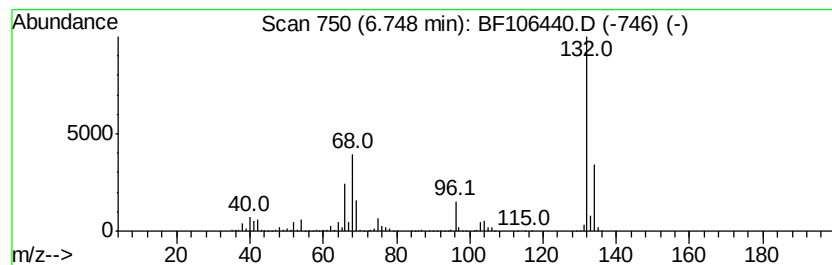
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown6.75 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	66.41 ng	2457050	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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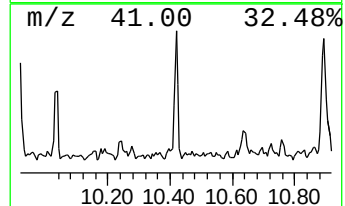
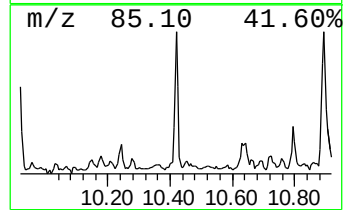
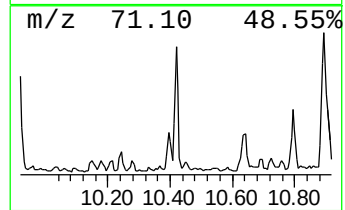
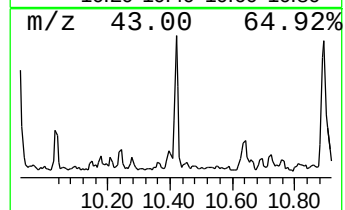
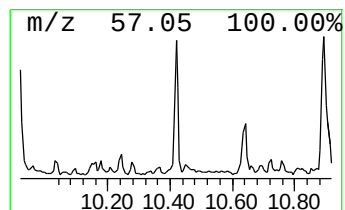
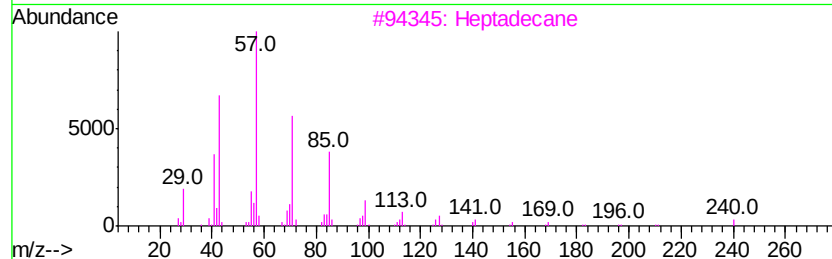
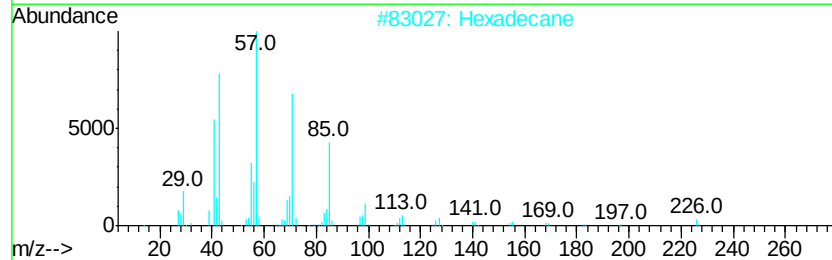
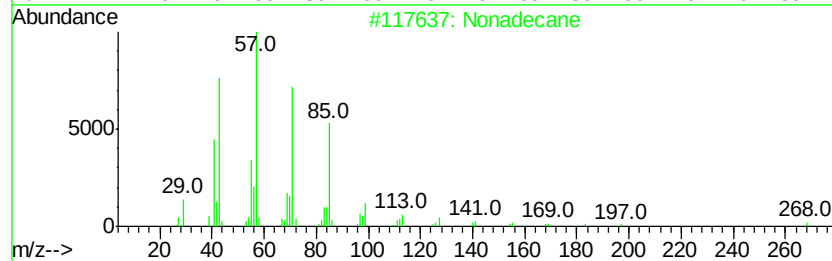
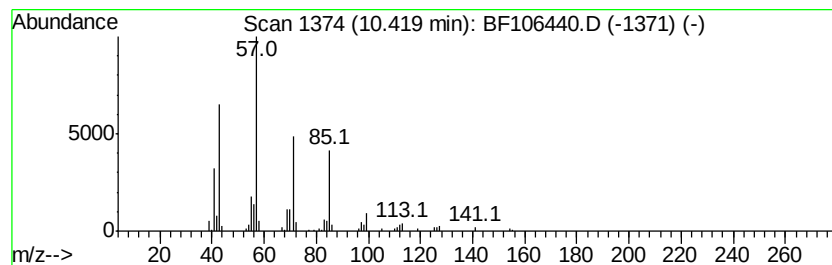
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Nonadecane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.42	2.09 ng	108019	Acenaphthene-d10		10.01
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	87
2	Hexadecane	226	C16H34	000544-76-3	64
3	Heptadecane	240	C17H36	000629-78-7	64
4	Pentadecane	212	C15H32	000629-62-9	59
5	Nonane, 5-butyl-	184	C13H28	017312-63-9	59



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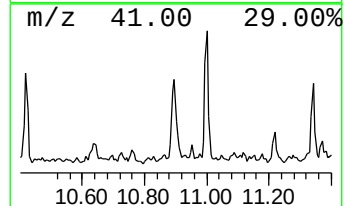
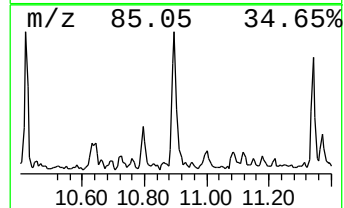
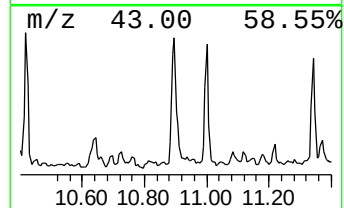
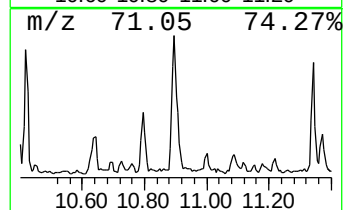
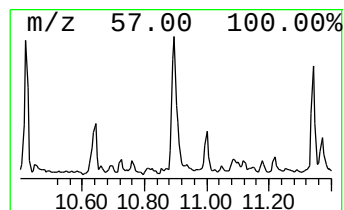
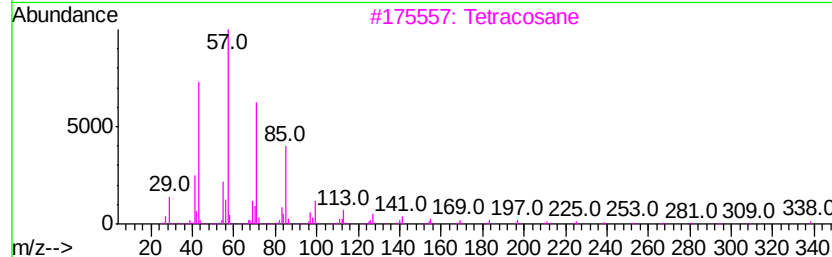
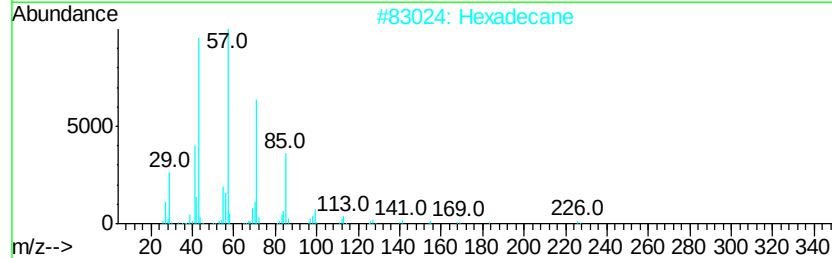
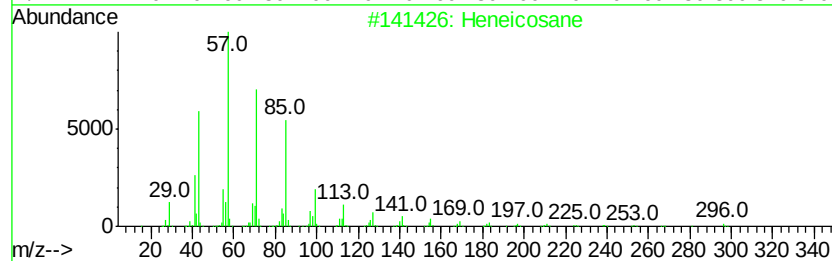
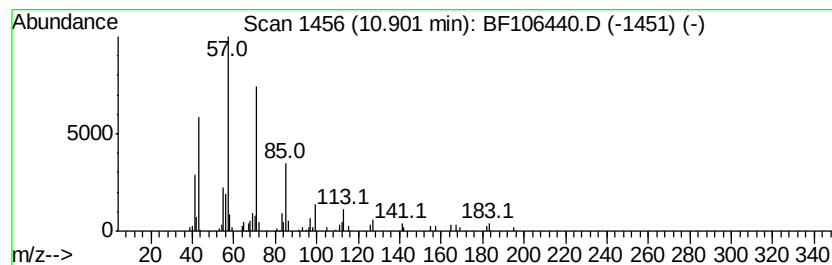
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Heneicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.90	3.26 ng	164878	Phenanthrene-d10	11.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	90
2	Hexadecane		226	C16H34	000544-76-3	90
3	Tetracosane		338	C24H50	000646-31-1	90
4	Nonadecane		268	C19H40	000629-92-5	86
5	Heptadecane, 8-methyl-		254	C18H38	013287-23-5	86



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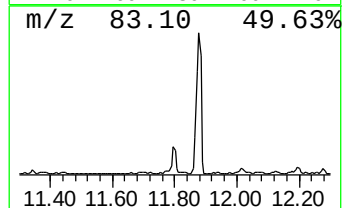
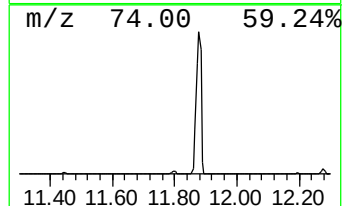
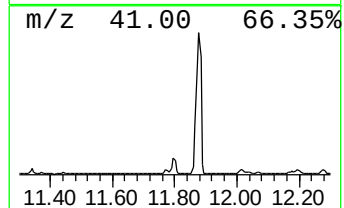
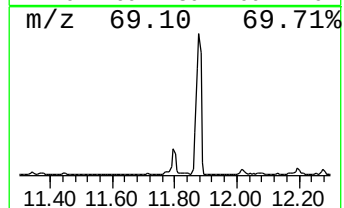
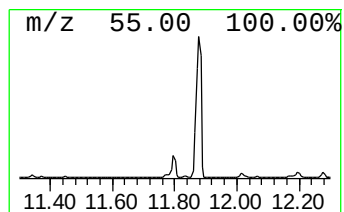
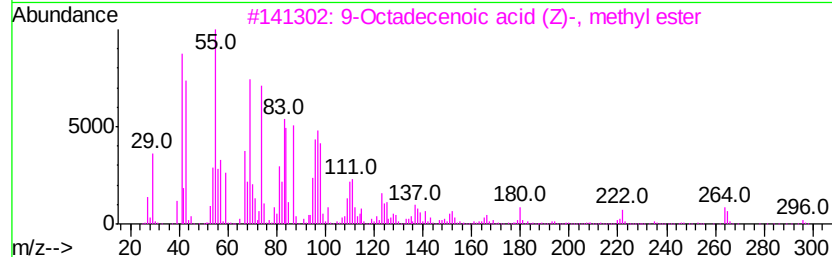
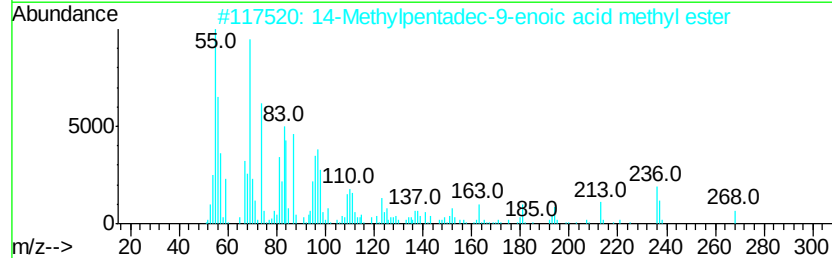
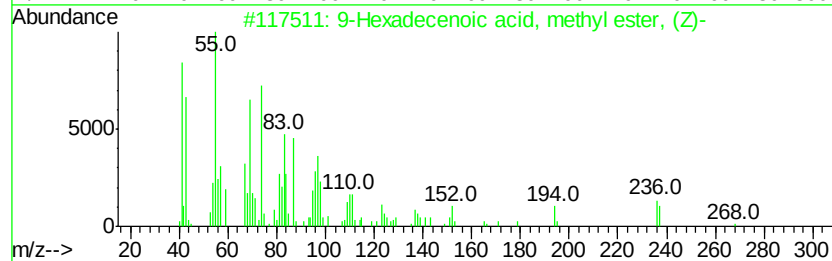
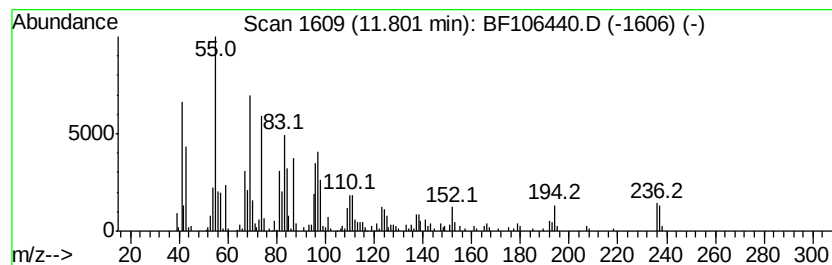
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 9-Hexadecenoic acid, methyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.80	6.62 ng	334794	Phenanthrene-d10	11.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Hexadecenoic acid, methyl este...	268	C17H32O2	001120-25-8	98
2		14-Methylpentadec-9-enoic acid m...	268	C17H32O2	1000365-89-7	91
3		9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	90
4		Methyl hexadec-9-enoate	268	C17H32O2	010030-74-7	90
5		Cyclopropaneoctanoic acid, 2-hex...	282	C18H34O2	010152-61-1	87



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Operator : JU/SJ
Sample : J3239-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-061318-B

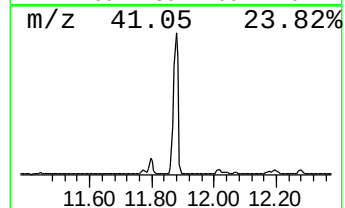
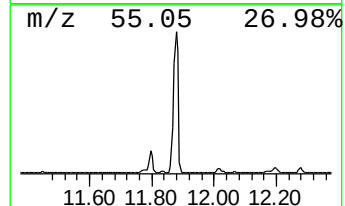
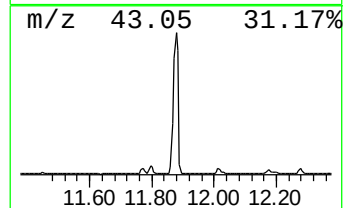
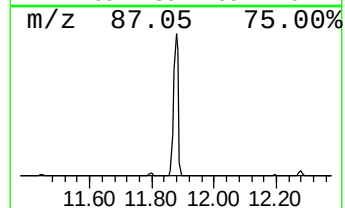
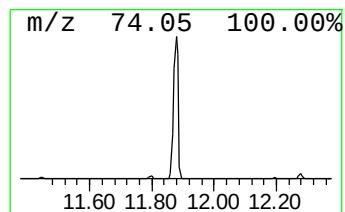
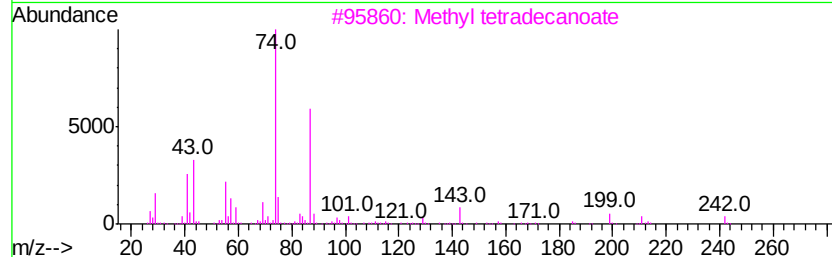
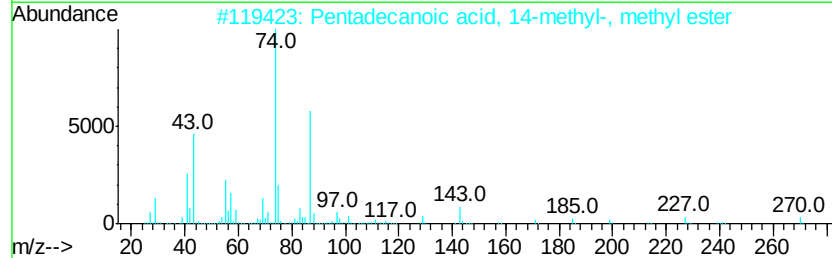
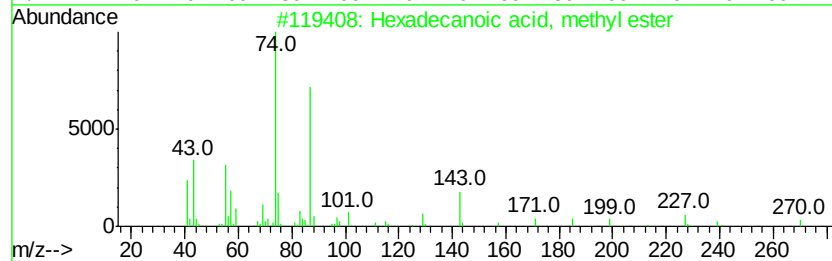
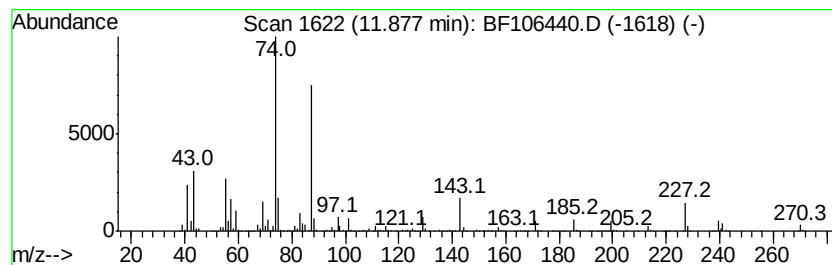
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexadecanoic acid, methyl e... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.88	97.02 ng	4906640	Phenanthrene-d10	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	95
3		Methyl tetradecanoate	242	C15H30O2	000124-10-7	91
4		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	91
5		Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061418\
Data File : BF106440.D
Acq On : 14 Jun 2018 14:23
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Sample : J3239-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

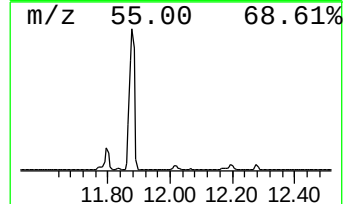
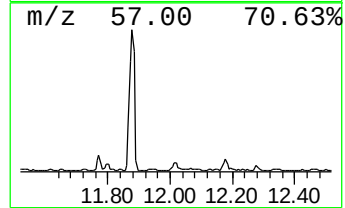
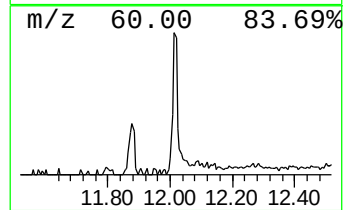
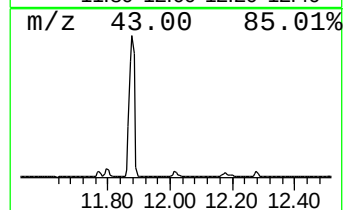
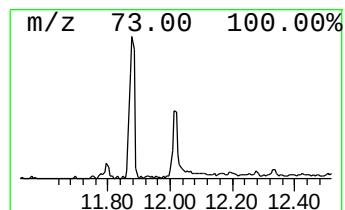
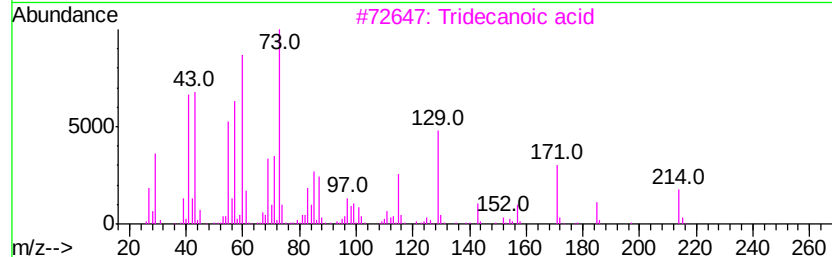
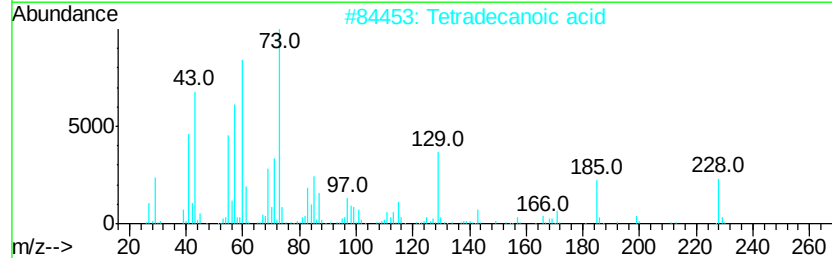
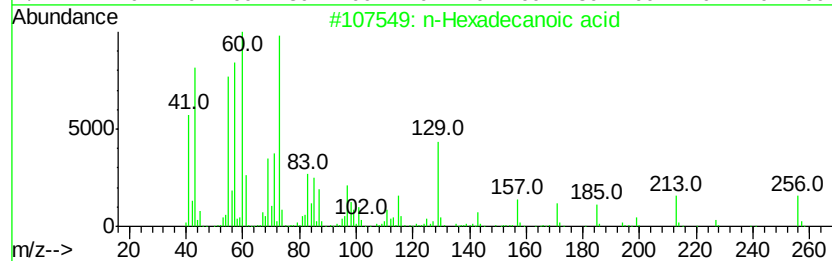
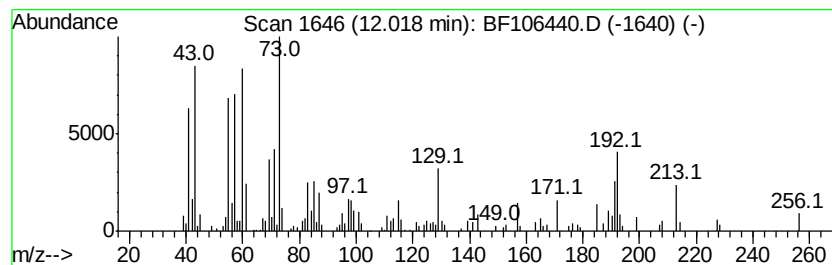
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.02	3.97 ng	200693	Phenanthrene-d10		11.49
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	93
3	Tridecanoic acid	214	C13H26O2	000638-53-9	91
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	58
5	Dodecanoic acid	200	C12H24O2	000143-07-7	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061418\
Data File : BF106440.D
Acq On : 14 Jun 2018 14:23
Operator : JU/SJ
Sample : J3239-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

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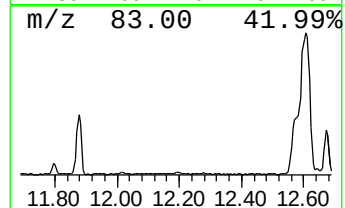
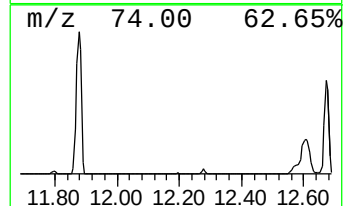
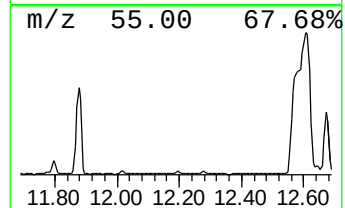
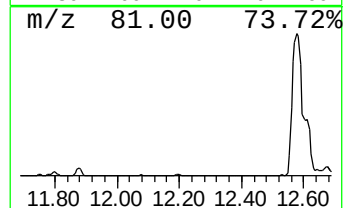
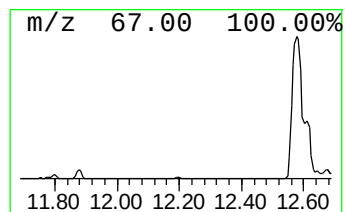
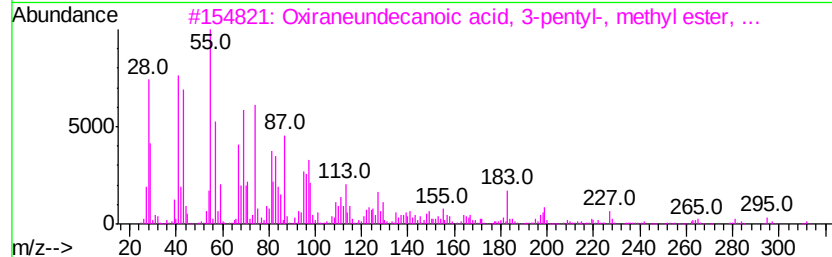
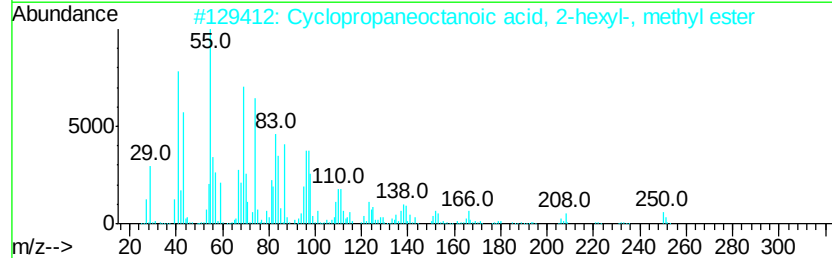
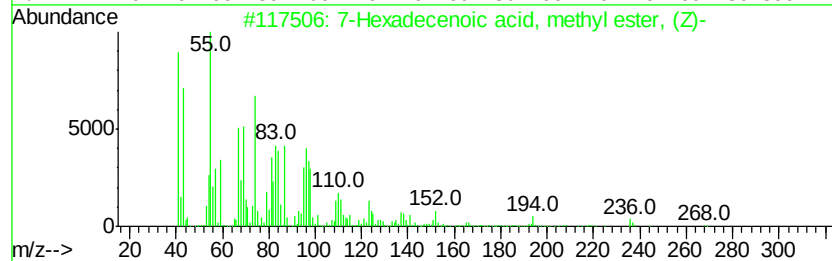
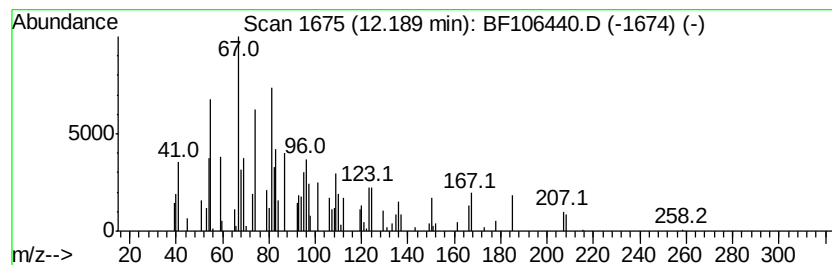
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 7-Hexadecenoic acid, methyl... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.19	2.03 ng	102455	Phenanthrene-d10	11.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	7-Hexadecenoic acid, methyl este...	268	C17H32O2	056875-67-3	83
2		Cyclopropaneoctanoic acid, 2-hex...	282	C18H34O2	010152-61-1	58
3		Oxiraneundecanoic acid, 3-pentyl...	312	C19H36O3	038520-31-9	58
4		9-Hexadecenoic acid, methyl este...	268	C17H32O2	001120-25-8	52
5		Methyl Z-11-tetradecenoate	240	C15H28O2	1000130-82-8	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061418\
Data File : BF106440.D
Acq On : 14 Jun 2018 14:23
Operator : JU/SJ
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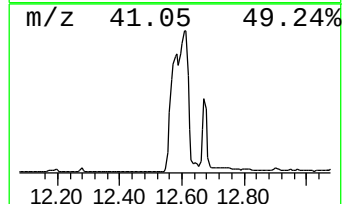
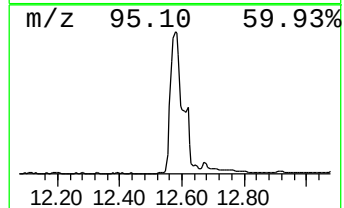
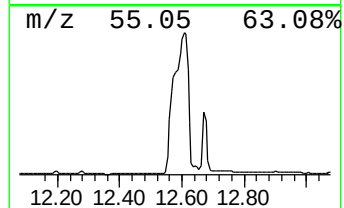
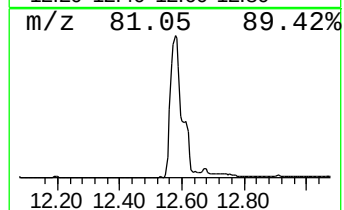
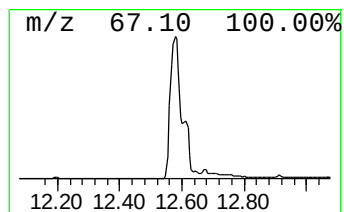
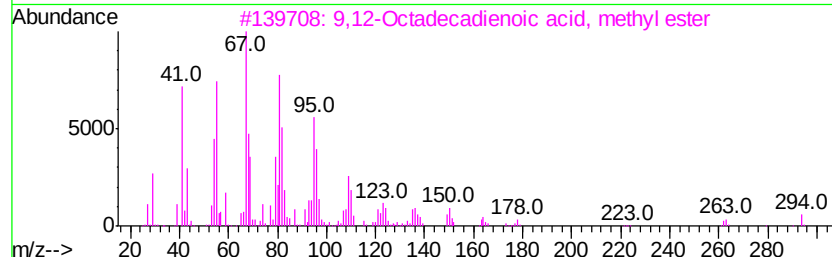
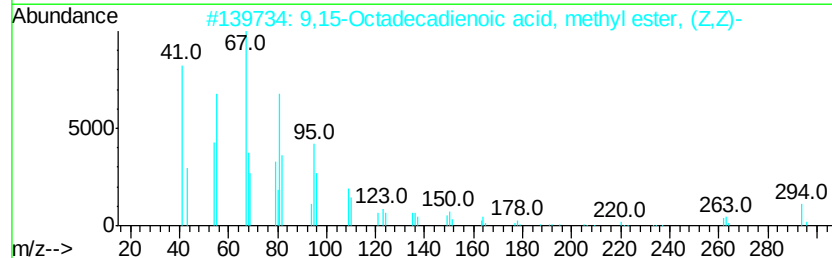
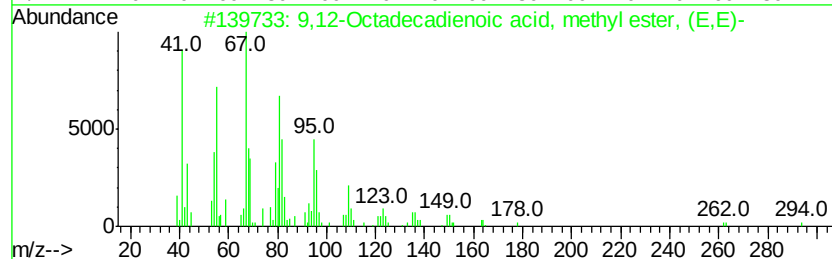
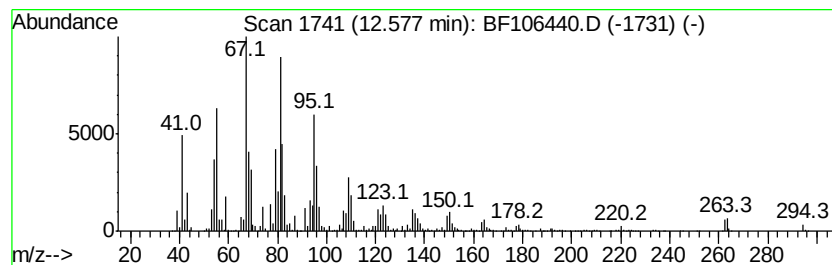
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.58	189.18 ng	9567790	Phenanthrene-d10	11.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
3	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
4	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	98
5	11,14-Octadecadienoic acid, meth...	294	C19H34O2	056554-61-1	98



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061418\
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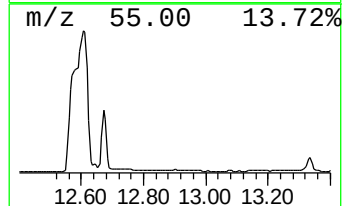
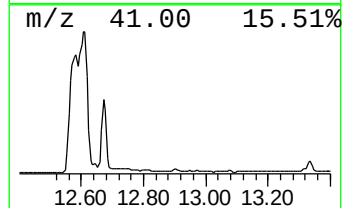
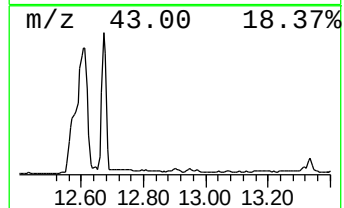
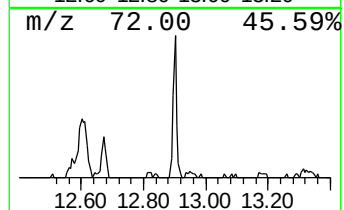
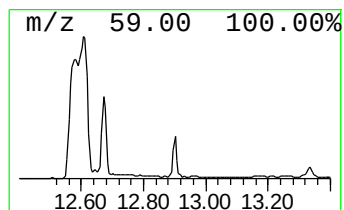
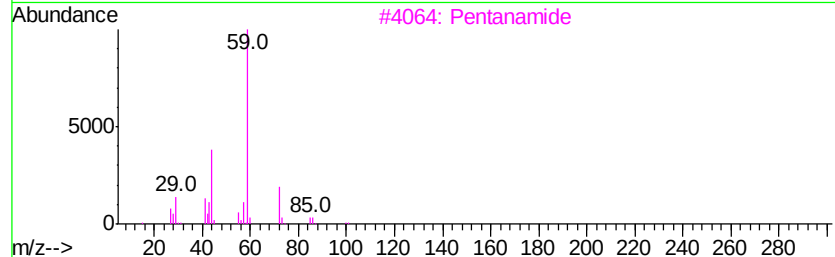
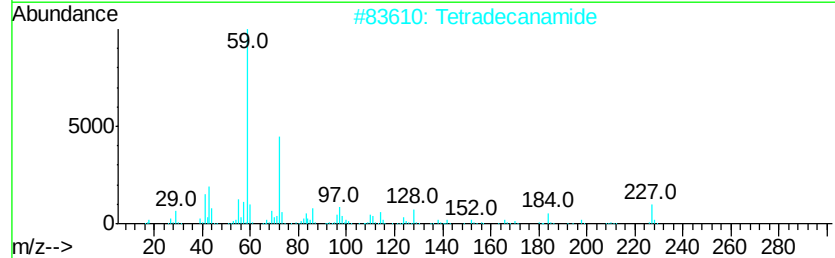
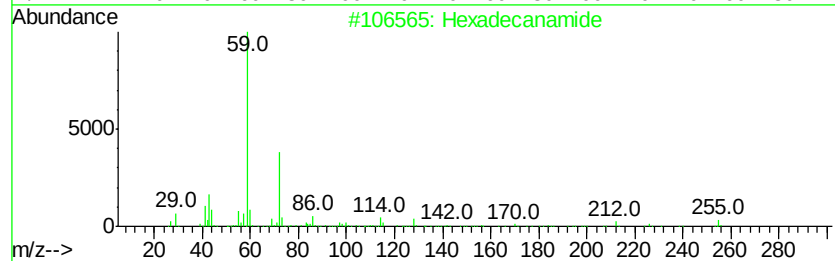
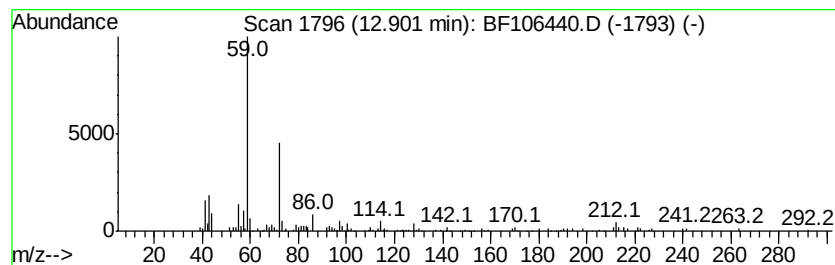
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexadecanamide Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.90	2.21 ng	115473	Chrysene-d12	14.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanamide	255	C16H33NO	000629-54-9	86
2		Tetradecanamide	227	C14H29NO	000638-58-4	83
3		Pentanamide	101	C5H11NO	000626-97-1	64
4		Hexanamide	115	C6H13NO	000628-02-4	64
5		Heptanamide, 4-ethyl-5-methyl-	171	C10H21NO	054789-40-1	56



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061418\
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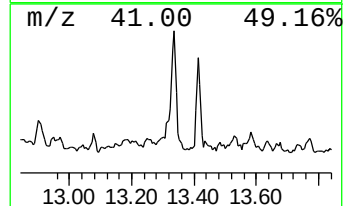
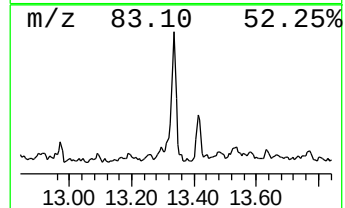
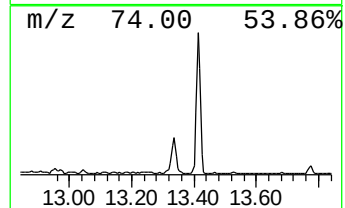
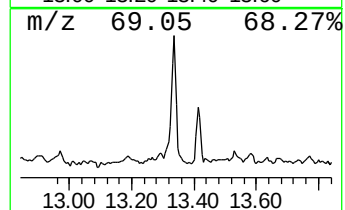
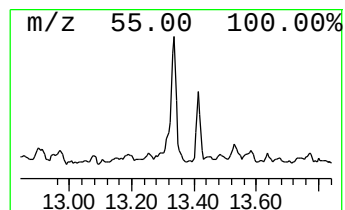
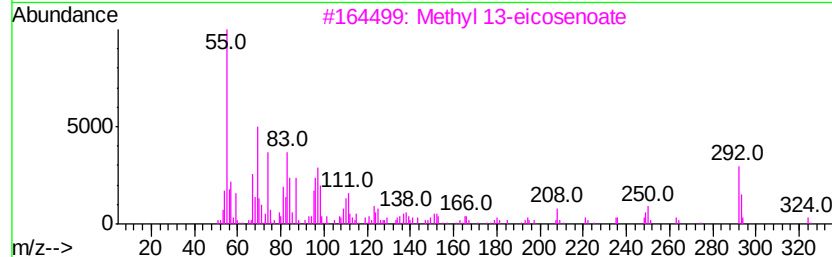
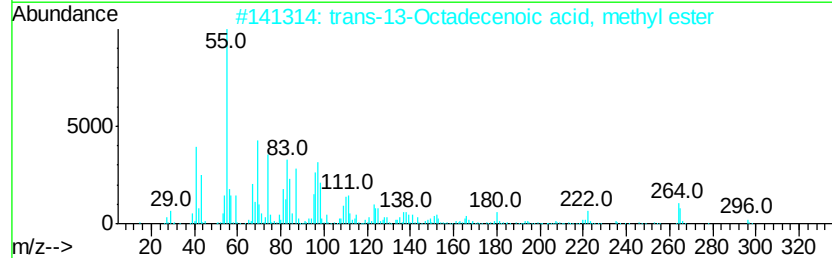
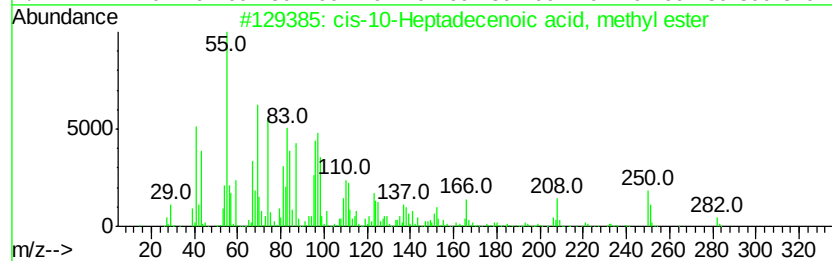
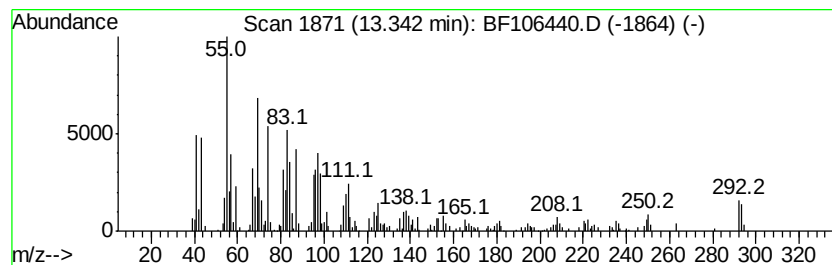
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 cis-10-Heptadecenoic acid, ... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	11.49 ng	601585	Chrysene-d12	14.21

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-10-Heptadecenoic acid, methy...	282	C18H34O2	1000333-62-1	91
2			trans-13-Octadecenoic acid, meth...	296	C19H36O2	1000333-61-3	91
3			Methyl 13-eicosenoate	324	C21H40O2	1000336-48-4	91
4			10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	81
5			Cyclopropanoic acid, 2-hex...	282	C18H34O2	010152-61-1	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061418\
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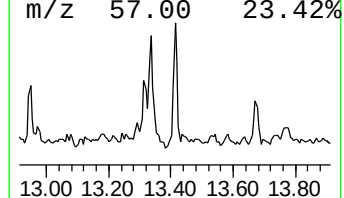
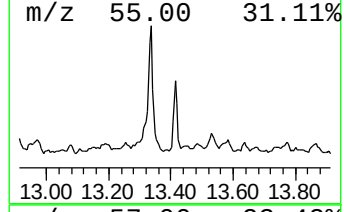
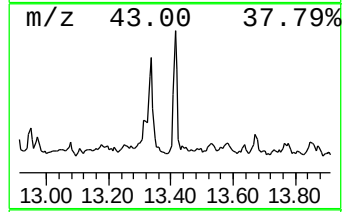
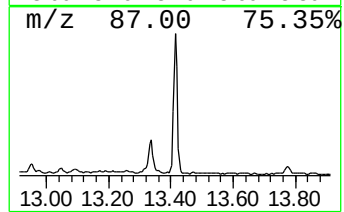
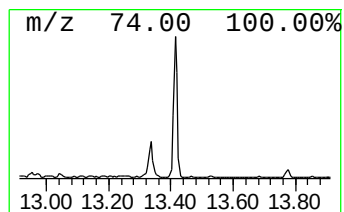
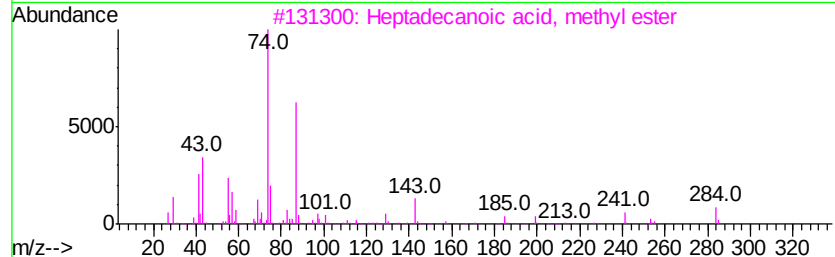
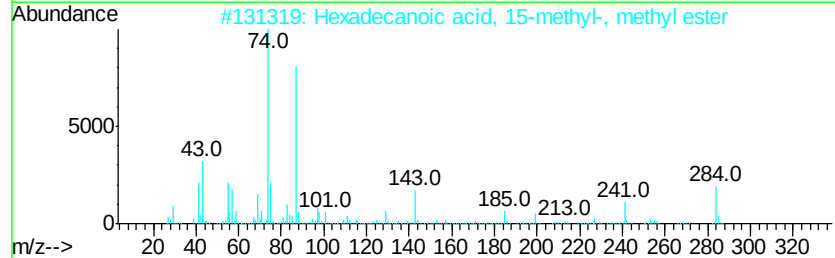
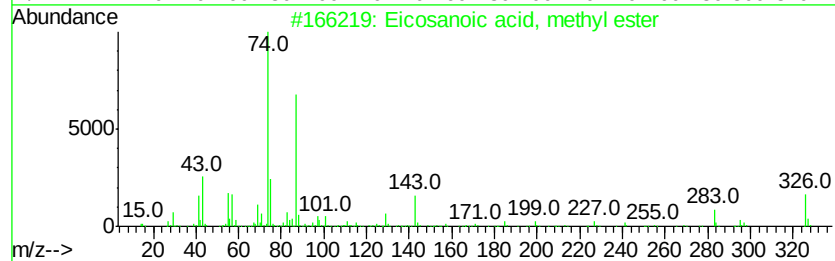
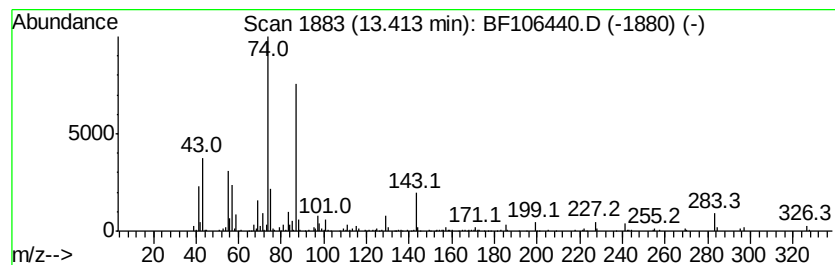
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Eicosanoic acid, methyl ester Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.41	6.00 ng	314094	Chrysene-d12	14.21	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosanoic acid, methyl ester	326	C21H42O2	001120-28-1	98
2	Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	94
3	Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	93
4	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	93
5	Methyl stearate	298	C19H38O2	000112-61-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061418\
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Acq On : 14 Jun 2018 14:23
Operator : JU/SJ
Sample : J3239-11
Misc :
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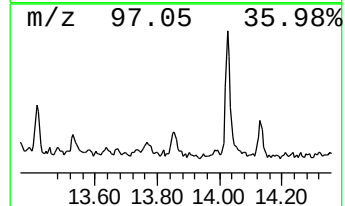
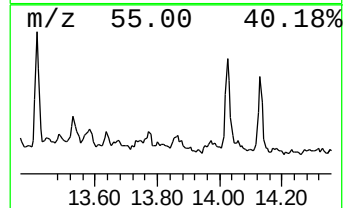
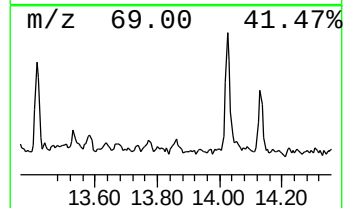
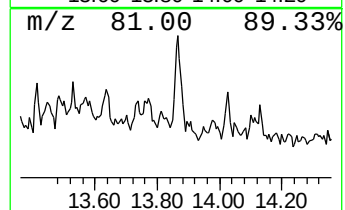
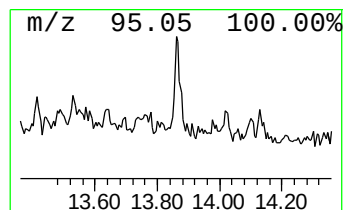
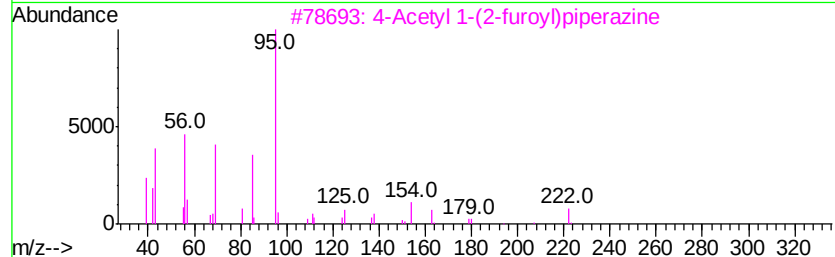
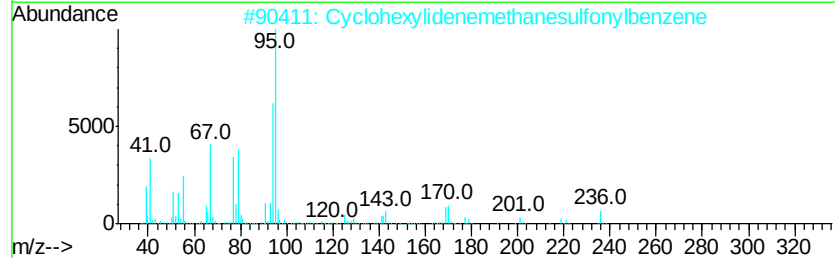
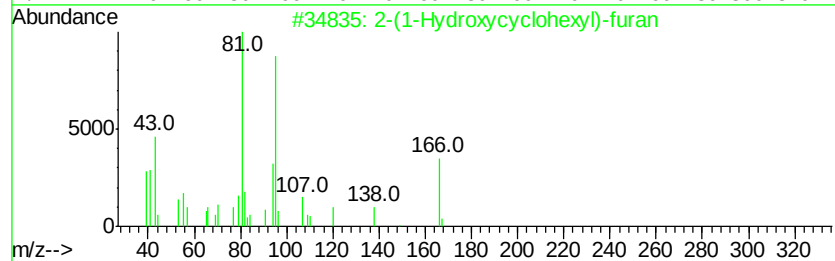
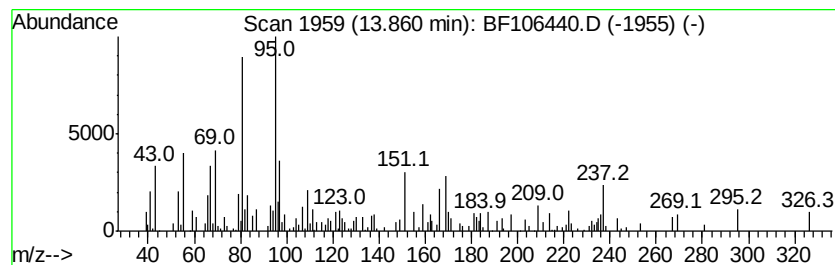
Instrument :
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ClientSampleId :
AU-06-061318-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown13.86 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	2.20 ng	115099	Chrysene-d12	14.21
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2-(1-Hydroxycyclohexyl)-furan	166 C10H14O2	036169-67-2	35
2	Cyclohexylidenemethanesulfonylbe...	236 C13H16O2S	090020-44-3	25
3	4-Acetyl 1-(2-furoyl)piperazine	222 C11H14N2O3	1000122-86-9	22
4	1H-Benzimidazol-1-ol, 2,4,5,6,7,...	184 C9H16N2O2	027992-37-6	22
5	6-Methyl-2-heptyne	110 C8H14	051065-64-6	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061418\
Data File : BF106440.D
Acq On : 14 Jun 2018 14:23
Operator : JU/SJ
Sample : J3239-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-061318-B

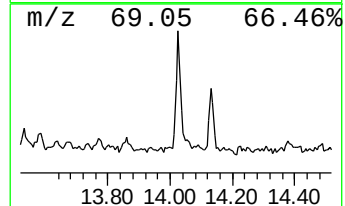
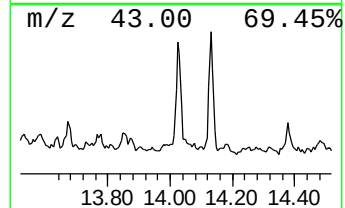
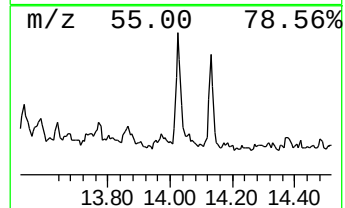
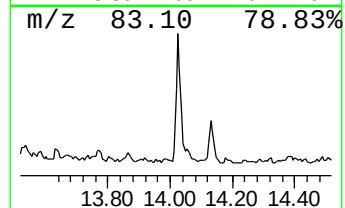
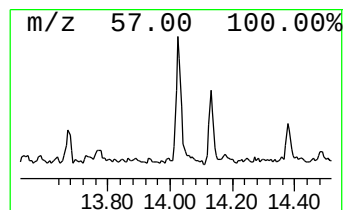
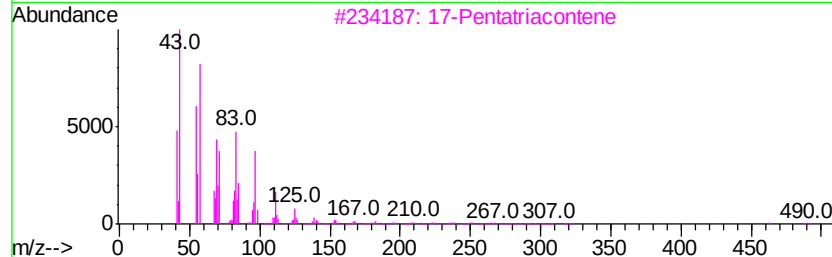
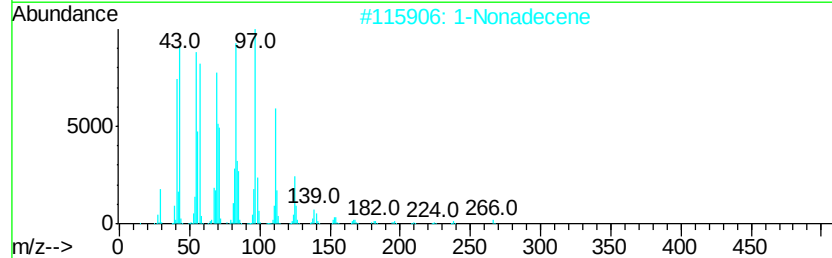
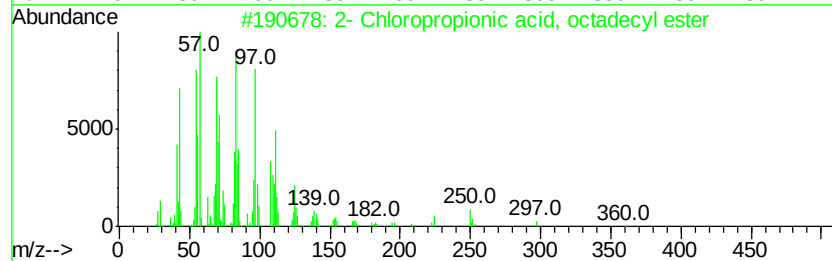
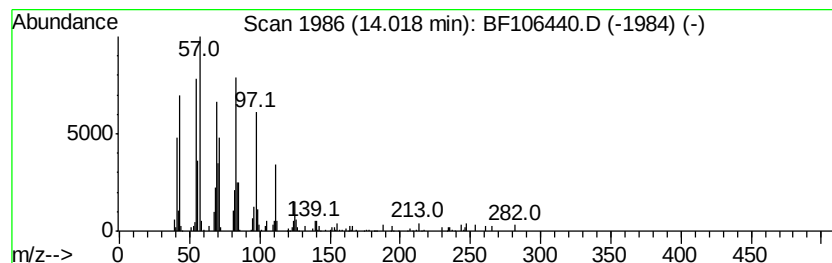
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 2- Chloropropionic acid, oc... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.02	4.82 ng	252371	Chrysene-d12	14.21

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	91
2			1-Nonadecene	266	C19H38	018435-45-5	90
3			17-Pentatriacontene	491	C35H70	006971-40-0	87
4			3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	87
5			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061418\
Data File : BF106440.D
Acq On : 14 Jun 2018 14:23
Operator : JU/SJ
Sample : J3239-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-06-061318-B

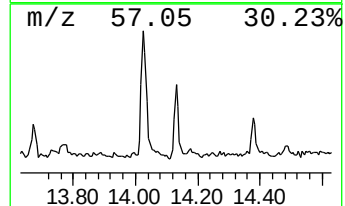
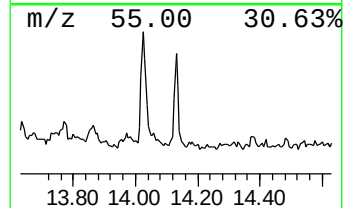
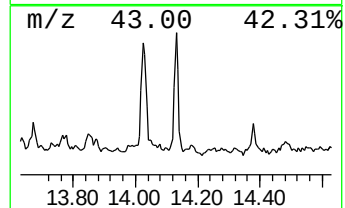
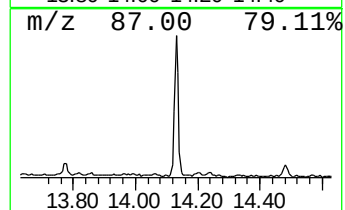
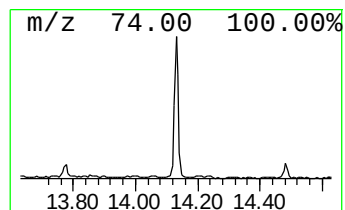
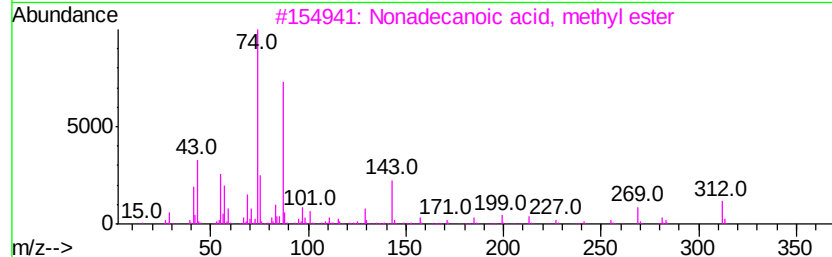
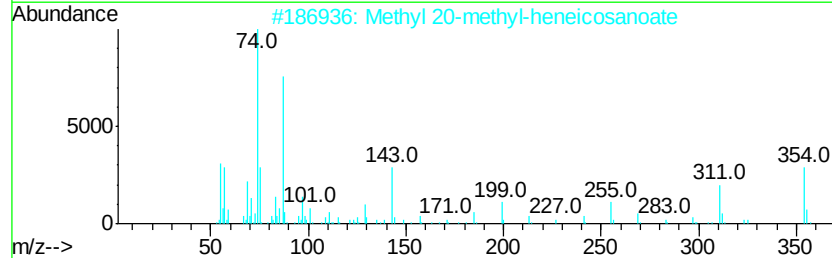
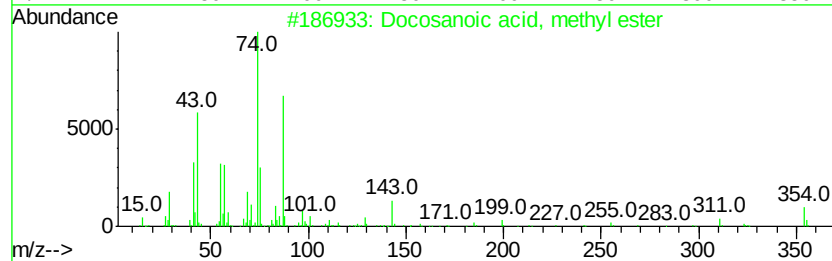
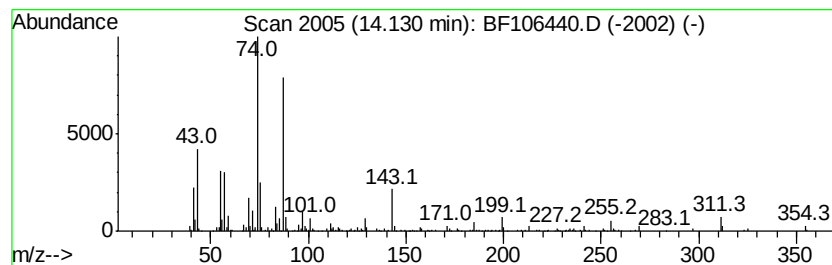
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Docosanoic acid, methyl ester Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.13	4.65 ng	243620	Chrysene-d12	14.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosanoic acid, methyl ester	354	C23H46O2	000929-77-1	98
2		Methyl 20-methyl-heneicosanoate	354	C23H46O2	1000336-47-4	98
3		Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	94
4		Methyl stearate	298	C19H38O2	000112-61-8	87
5		Octadecanoic acid, 10-methyl-, m...	312	C20H40O2	002490-19-9	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF061418\
 Data File : BF106440.D
 Acq On : 14 Jun 2018 14:23
 Operator : JU/SJ
 Sample : J3239-11
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-06-061318-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061118.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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unknown6.75	6.75	66.4	ng	2457050	1	6.97	739945	20.0
Nonadecane	10.42	2.1	ng	108019	3	10.01	1033250	20.0
Heneicosane	10.90	3.3	ng	164878	4	11.49	1011520	20.0
9-Hexadecenoic ac...	11.80	6.6	ng	334794	4	11.49	1011520	20.0
Hexadecanoic acid...	11.88	97.0	ng	4906640	4	11.49	1011520	20.0
n-Hexadecanoic acid	12.02	4.0	ng	200693	4	11.49	1011520	20.0
7-Hexadecenoic ac...	12.19	2.0	ng	102455	4	11.49	1011520	20.0
9,12-Octadecadien...	12.58	189.2	ng	9567790	4	11.49	1011520	20.0
Hexadecanamide	12.90	2.2	ng	115473	5	14.21	1046960	20.0
cis-10-Heptadecen...	13.34	11.5	ng	601585	5	14.21	1046960	20.0
Eicosanoic acid, ...	13.41	6.0	ng	314094	5	14.21	1046960	20.0
unknown13.86	13.86	2.2	ng	115099	5	14.21	1046960	20.0
2-Chloropropioni...	14.02	4.8	ng	252371	5	14.21	1046960	20.0
Docosanoic acid, ...	14.13	4.7	ng	243620	5	14.21	1046960	20.0