

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122717\
 Data File : BF101769.D
 Acq On : 27 Dec 2017 18:22
 Operator : SJ/JU
 Sample : I6685-09 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-05-122617-C

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:\HPCHEM1\BNA_F\METHODS\8270-BF121417.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	5.004	493	496	514	rBV	125465	221830	1.44%	0.447%
2	5.410	562	565	599	rBV	1314447	2135121	13.90%	4.307%
3	6.434	736	739	757	rBV	1440025	2114333	13.76%	4.265%
4	6.563	757	761	767	rVV	2008762	1990627	12.96%	4.015%
5	6.786	796	799	814	rBV	1043863	992990	6.46%	2.003%
6	6.939	822	825	837	rBV	1514090	1546223	10.06%	3.119%
7	7.357	893	896	912	rBV	1095416	1385205	9.02%	2.794%
8	8.069	1014	1017	1023	rBV	1283311	1217598	7.92%	2.456%
9	8.645	1110	1115	1119	rBV	231532	202658	1.32%	0.409%
10	9.145	1196	1200	1208	rBV	2011442	1999231	13.01%	4.033%
11	9.210	1208	1211	1215	rVB	324525	256514	1.67%	0.517%
12	9.545	1266	1268	1274	rVB	171514	184829	1.20%	0.373%
13	9.739	1298	1301	1305	rVB	374598	266422	1.73%	0.537%
14	9.827	1312	1316	1323	rVB	1285098	1192687	7.76%	2.406%
15	10.239	1380	1386	1389	rBV	369003	303767	1.98%	0.613%
16	10.621	1448	1451	1463	rBV	960036	1044737	6.80%	2.107%
17	10.710	1463	1466	1476	rVB	319350	405442	2.64%	0.818%
18	11.157	1539	1542	1545	rBV	223763	191477	1.25%	0.386%
19	11.321	1566	1570	1578	rVB	1009184	1070789	6.97%	2.160%
20	11.698	1629	1634	1637	rBV	5598700	6829448	44.45%	13.775%
21	11.833	1654	1657	1667	rBV2	129199	181244	1.18%	0.366%
22	12.404	1745	1754	1756	rBV2	5892922	15364078	100.00%	30.990%
23	12.486	1765	1768	1772	rVB	3671942	3538289	23.03%	7.137%
24	12.721	1804	1808	1811	rVB	271251	261639	1.70%	0.528%
25	12.904	1833	1839	1842	rBV	1456047	1388517	9.04%	2.801%
26	13.127	1875	1877	1884	rVB	293998	318395	2.07%	0.642%
27	13.204	1887	1890	1893	rBV	462454	410502	2.67%	0.828%
28	13.639	1957	1964	1969	rBV4	299247	602207	3.92%	1.215%
29	13.874	2000	2004	2009	rBV	405020	373282	2.43%	0.753%
30	13.974	2017	2021	2025	rBV	881165	882351	5.74%	1.780%
31	15.445	2267	2271	2278	rBV	524042	704927	4.59%	1.422%

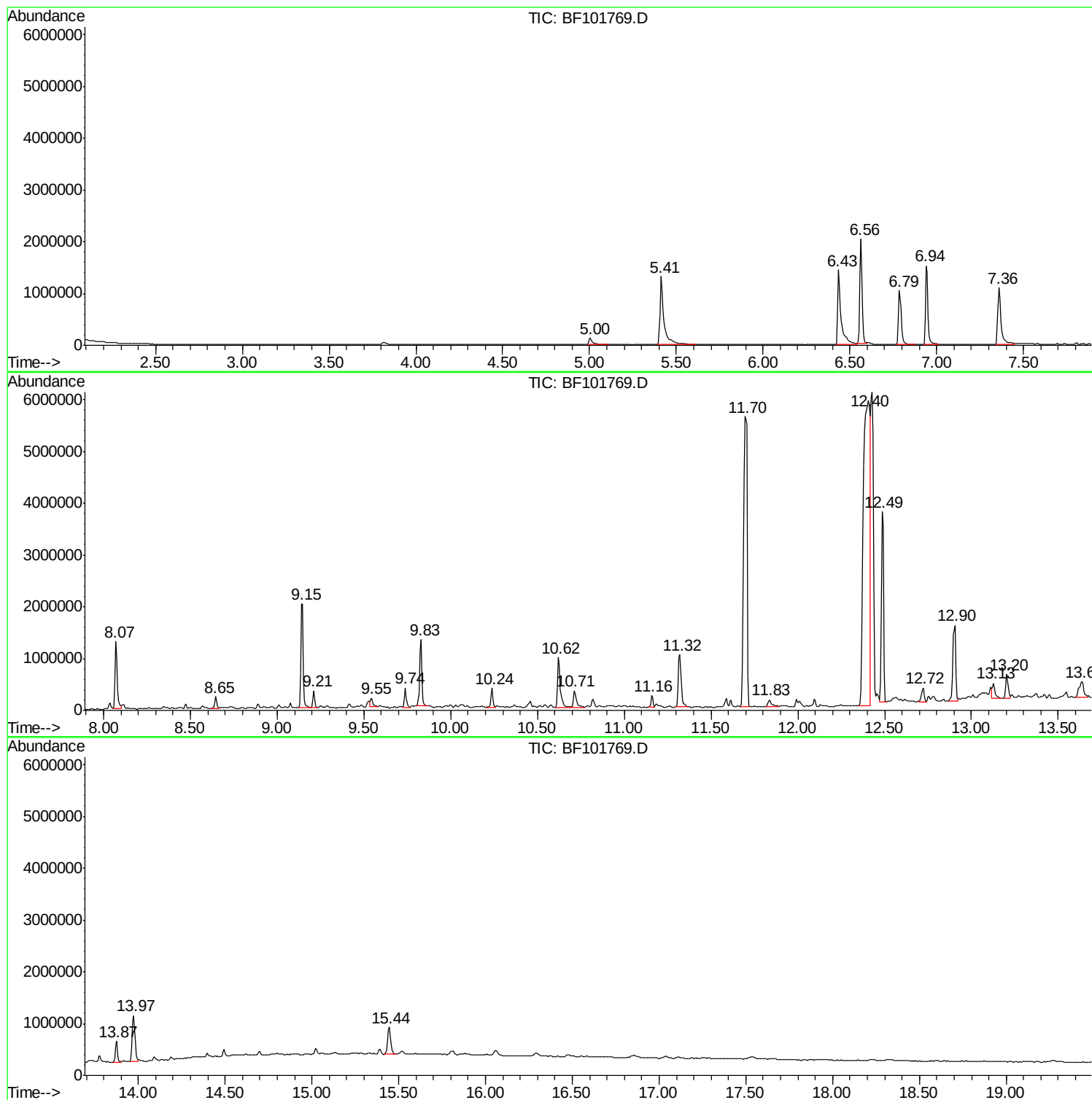
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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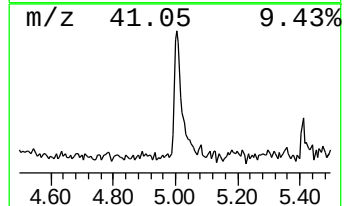
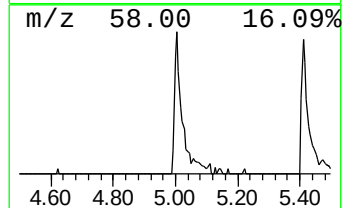
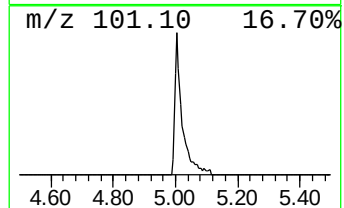
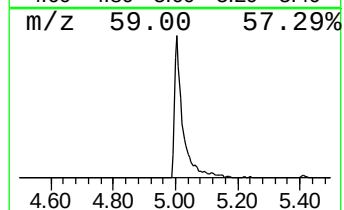
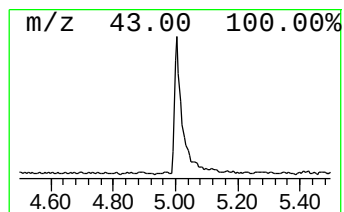
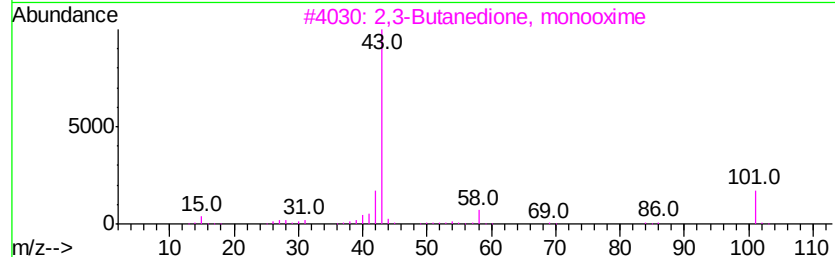
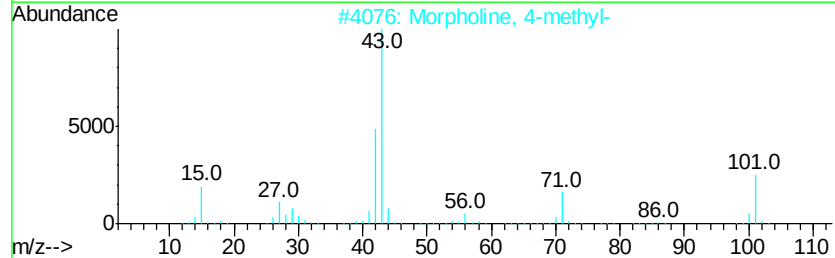
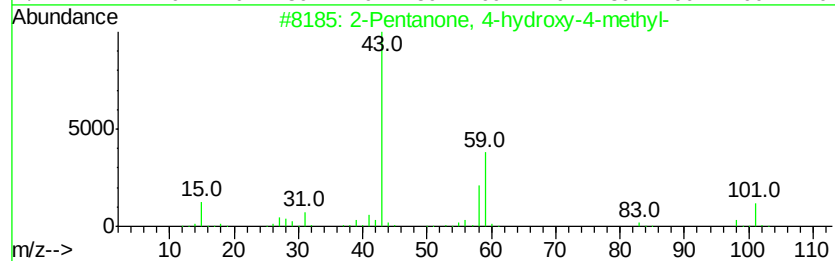
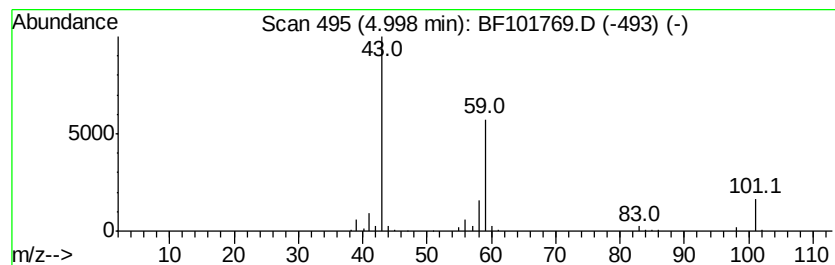
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	4.47 ng	221830	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			5-Oxohexanenitrile	111	C6H9NO	010412-98-3	9
5			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	9



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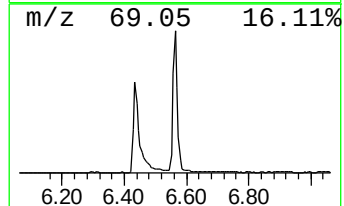
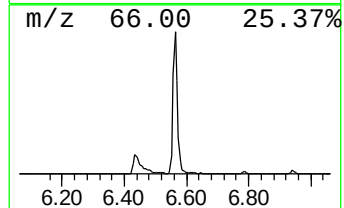
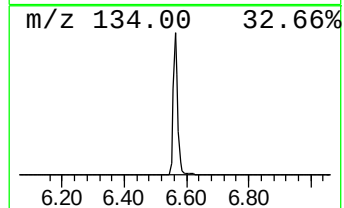
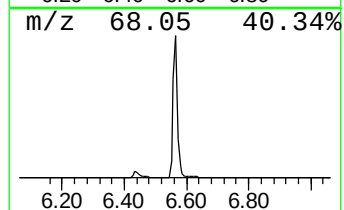
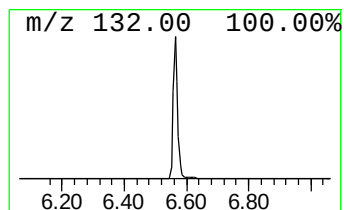
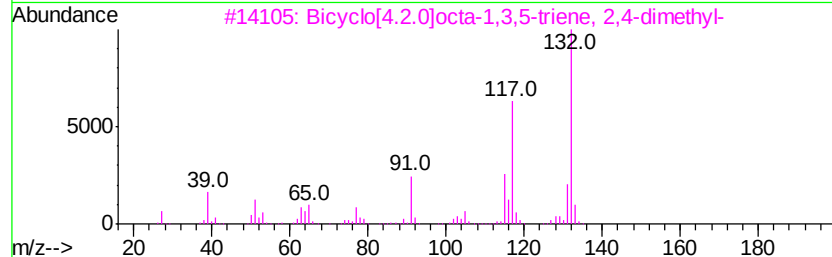
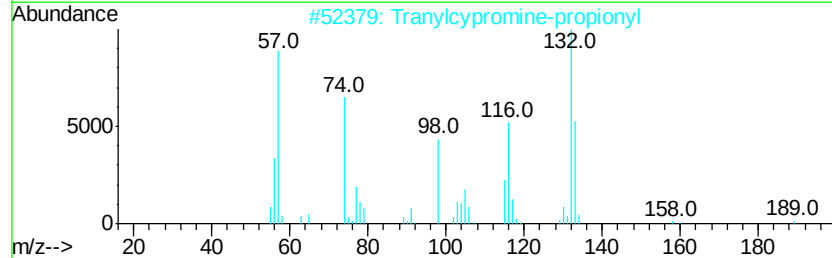
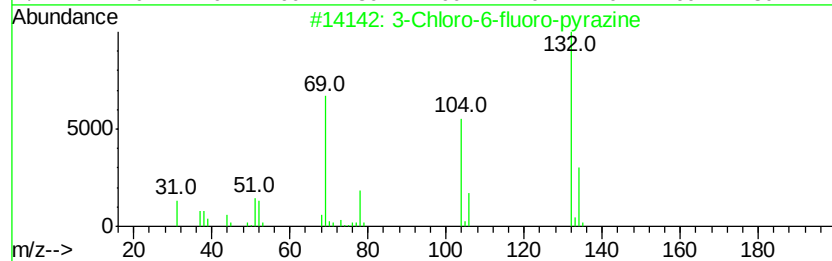
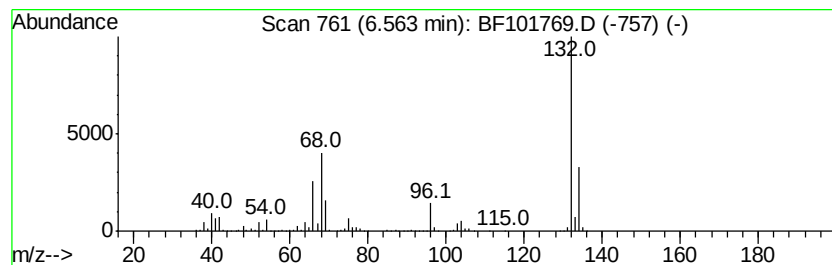
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown6.56 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	40.09 ng	1990630	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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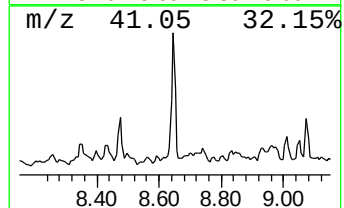
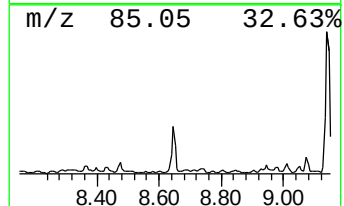
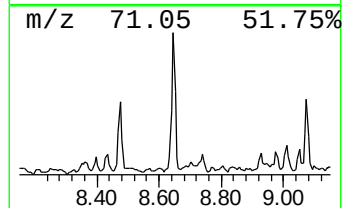
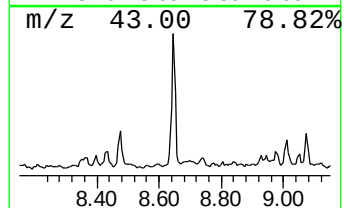
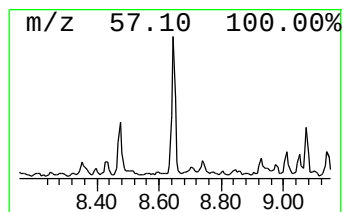
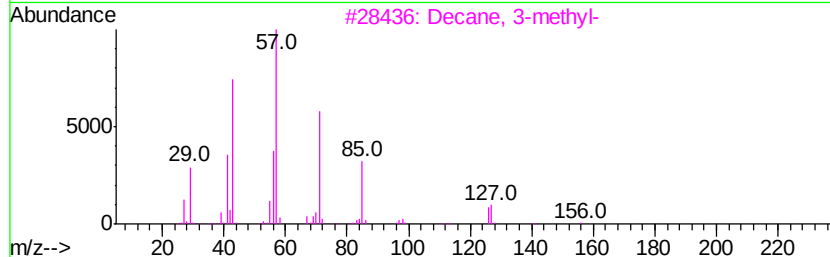
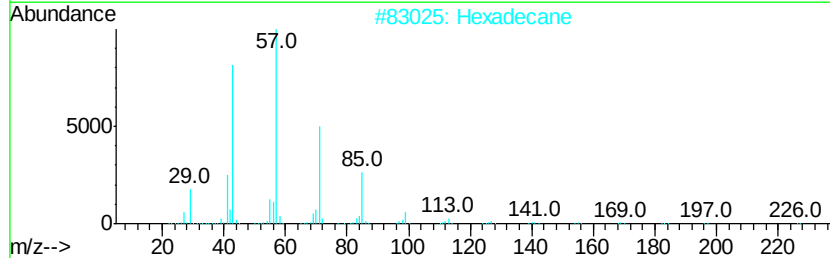
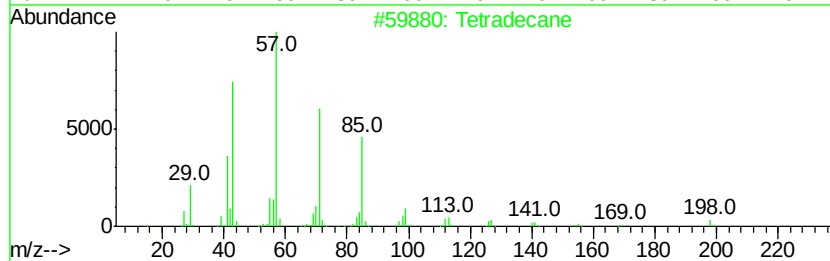
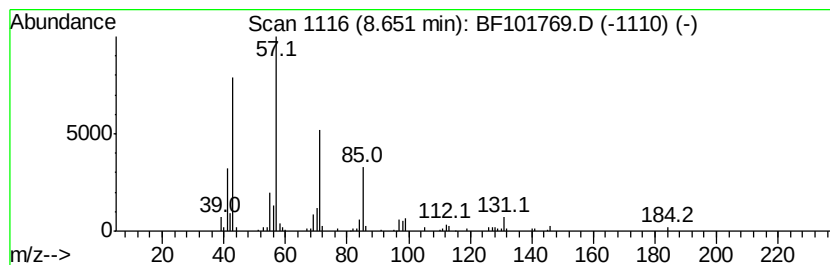
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Peak Number 4 Tetradecane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.65	3.33 ng	202658	Napthalene-d8	8.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	90
2		Hexadecane	226	C16H34	000544-76-3	86
3		Decane, 3-methyl-	156	C11H24	013151-34-3	64
4		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	64
5		Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	62



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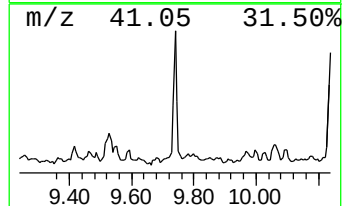
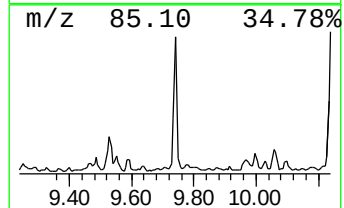
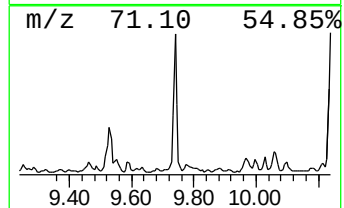
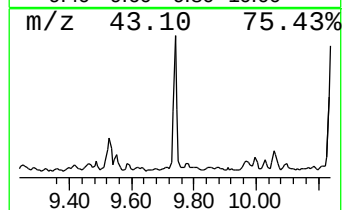
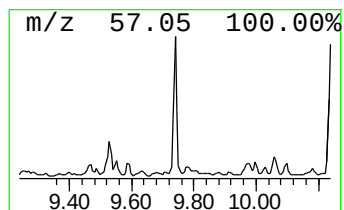
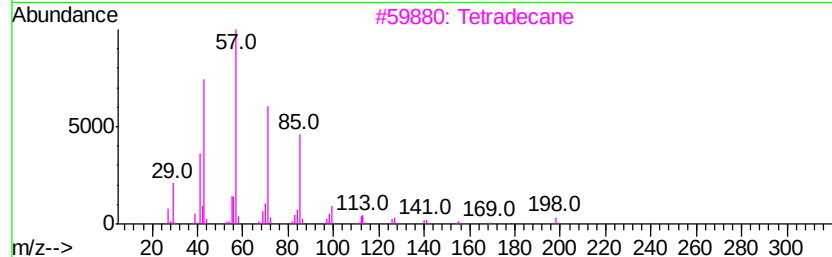
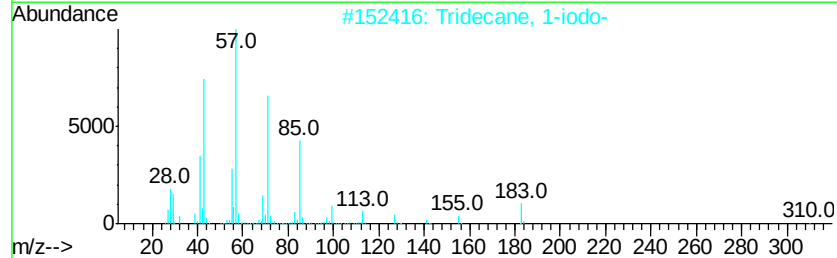
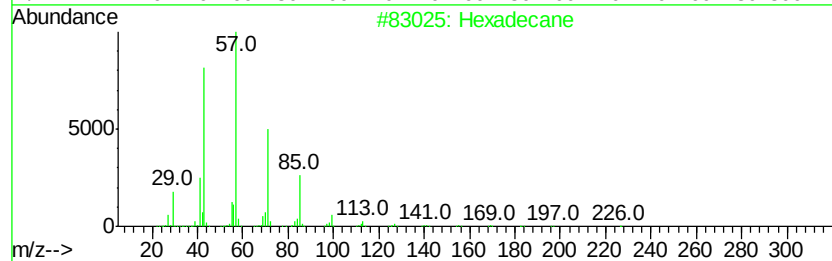
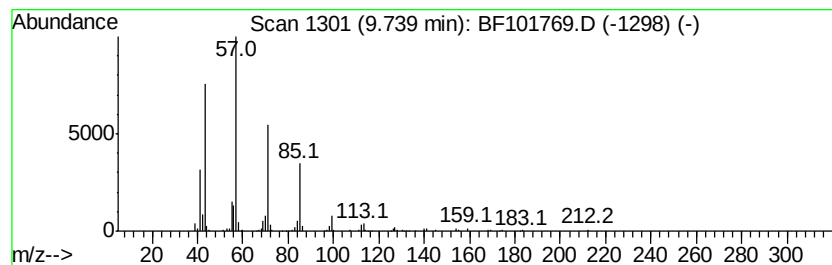
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 Hexadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.74	4.47 ng	266422	Acenaphthene-d10	9.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	91
2		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90
3		Tetradecane	198	C14H30	000629-59-4	90
4		Bacchotricuneatin c	342	C20H22O5	066563-30-2	74
5		Heptadecane	240	C17H36	000629-78-7	72



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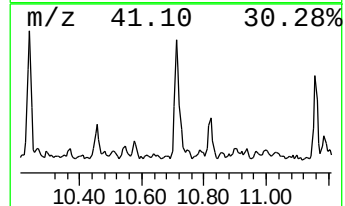
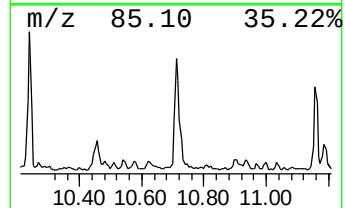
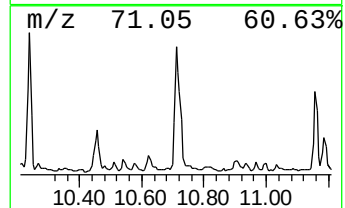
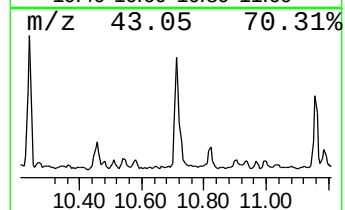
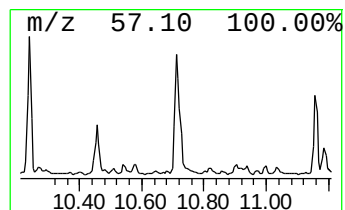
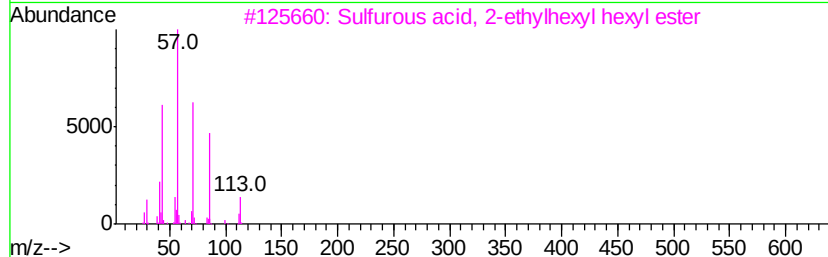
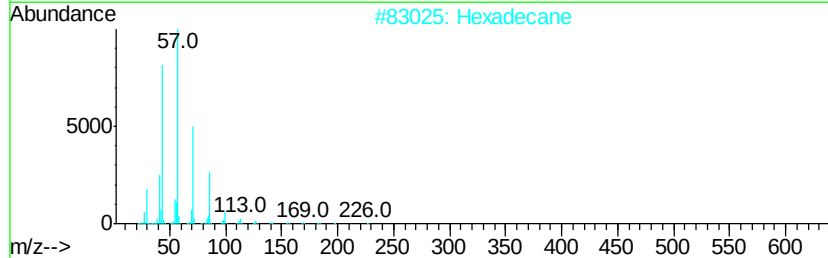
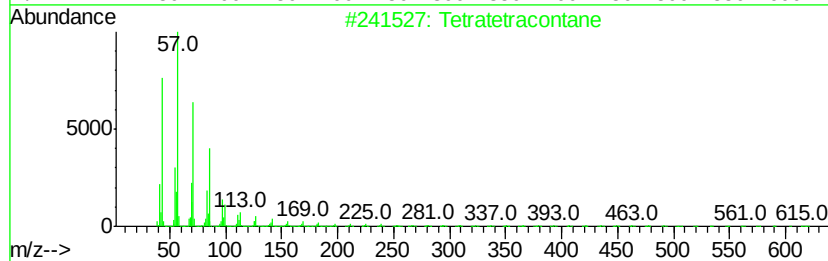
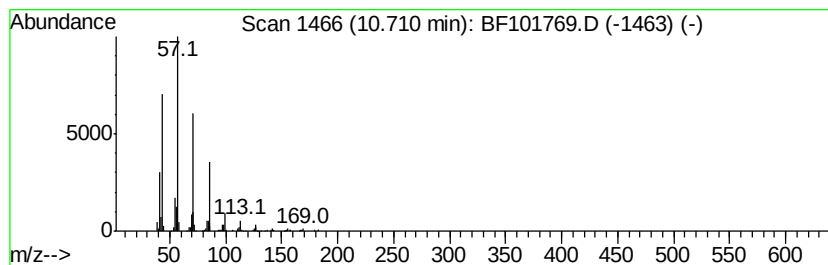
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 8 Tetratetracontane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.71	7.57 ng	405442	Phenanthrene-d10	11.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetratetracontane	619	C44H90	007098-22-8	90
2		Hexadecane	226	C16H34	000544-76-3	90
3		Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	86
4		Heneicosane	296	C21H44	000629-94-7	83
5		Tridecane	184	C13H28	000629-50-5	83



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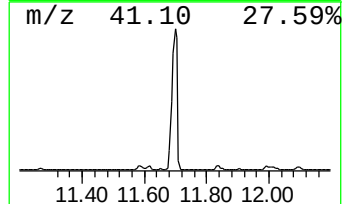
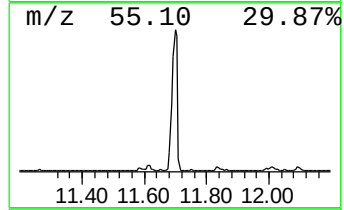
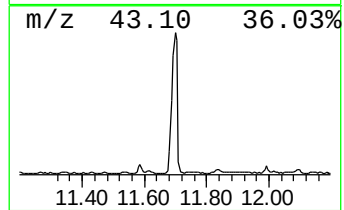
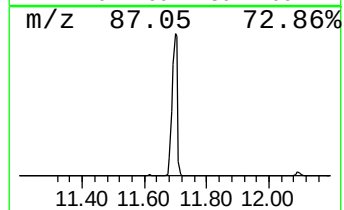
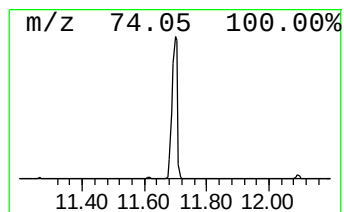
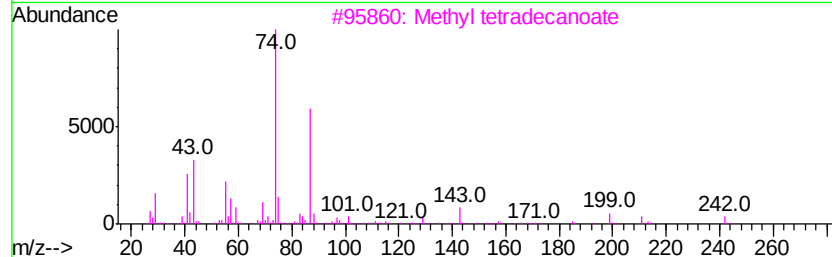
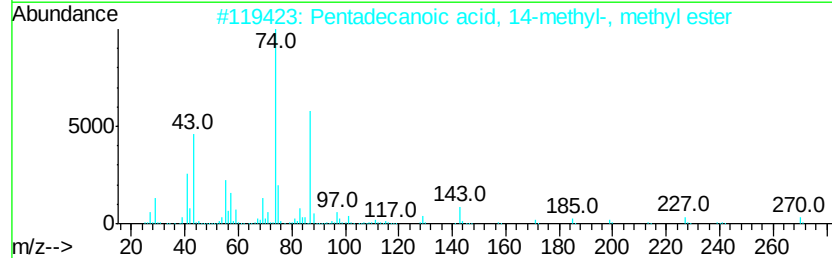
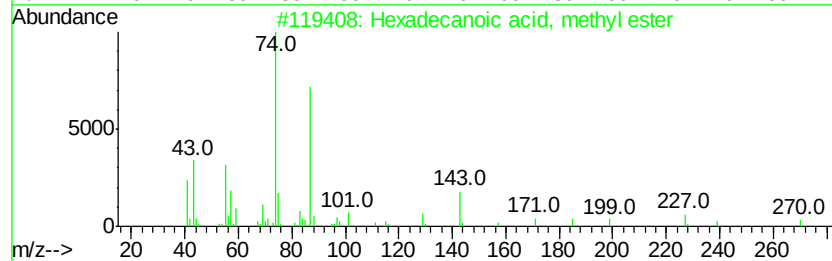
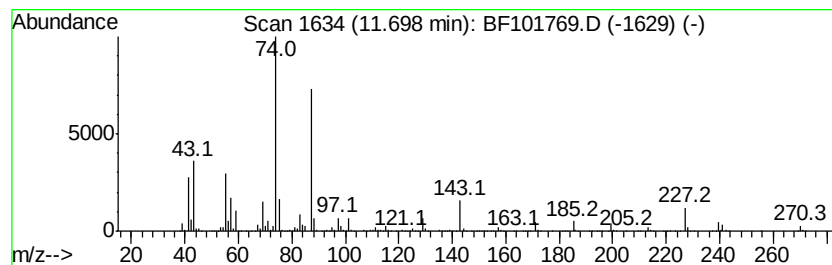
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecanoic acid, methyl e... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.70	127.56 ng	6829450	Phenanthrene-d10	11.32

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	96
3		Methyl tetradecanoate	242	C15H30O2	000124-10-7	91
4		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	90
5		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122717\
Data File : BF101769.D
Acq On : 27 Dec 2017 18:22
Operator : SJ/JU
Sample : I6685-09 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
AU-05-122617-C

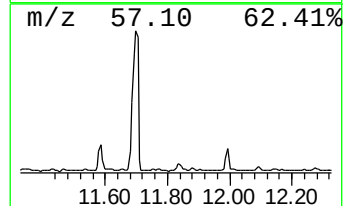
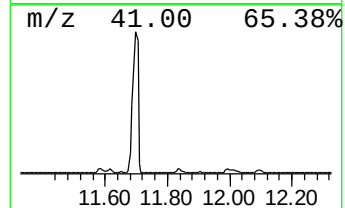
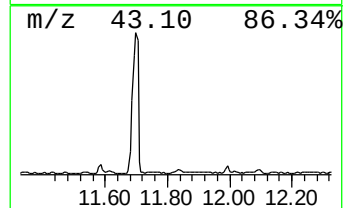
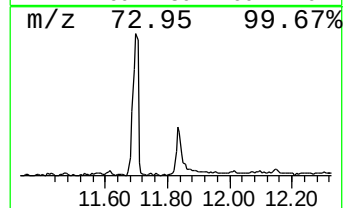
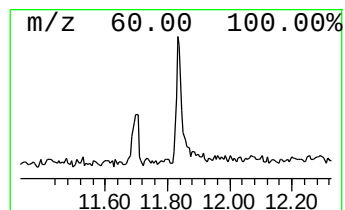
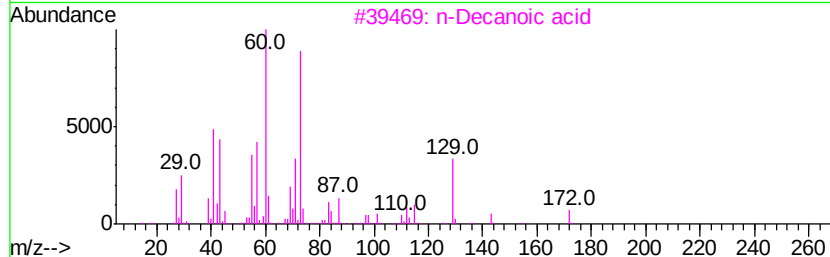
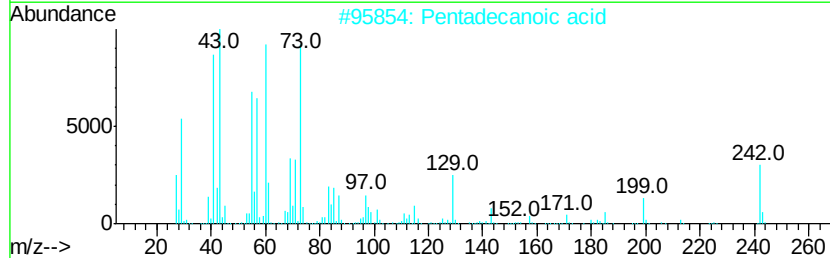
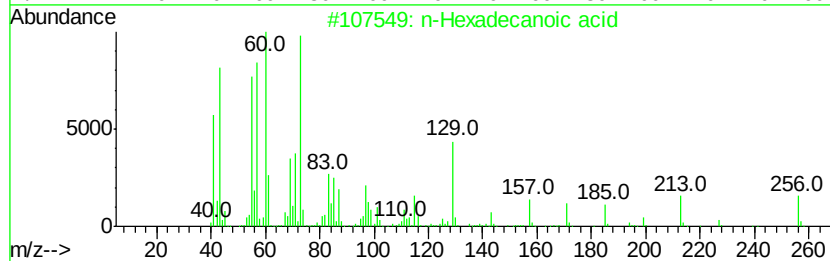
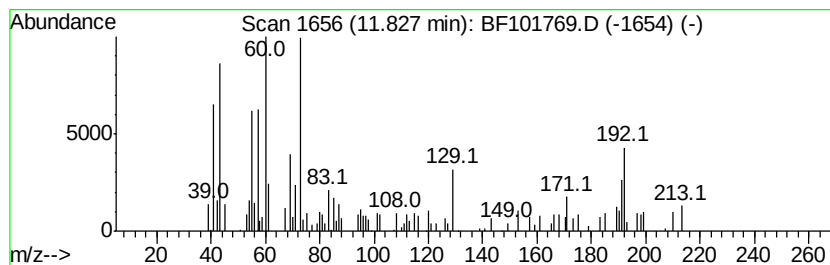
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 n-Hexadecanoic acid Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.83	3.39 ng	181244	Phenanthrene-d10	11.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	86
3			n-Decanoic acid	172	C10H20O2	000334-48-5	64
4			Dodecanoic acid	200	C12H24O2	000143-07-7	53
5			Oxalic acid, dodecyl propyl ester	300	C17H32O4	1000309-26-5	30



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122717\
Data File : BF101769.D
Acq On : 27 Dec 2017 18:22
Operator : SJ/JU
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Misc :
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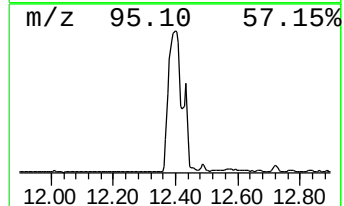
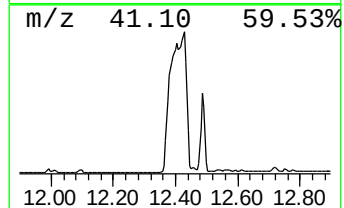
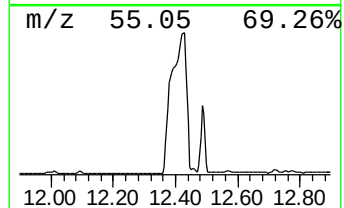
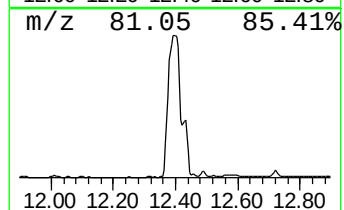
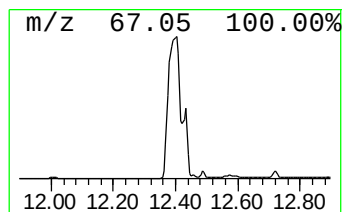
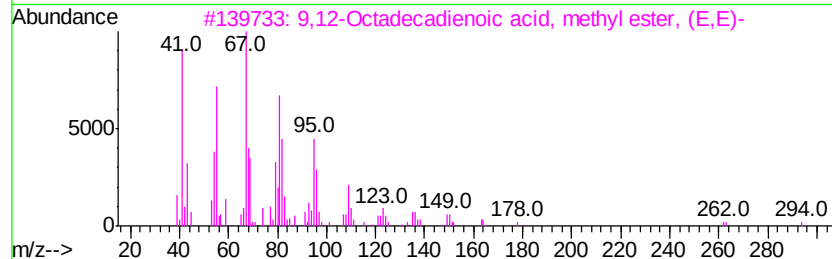
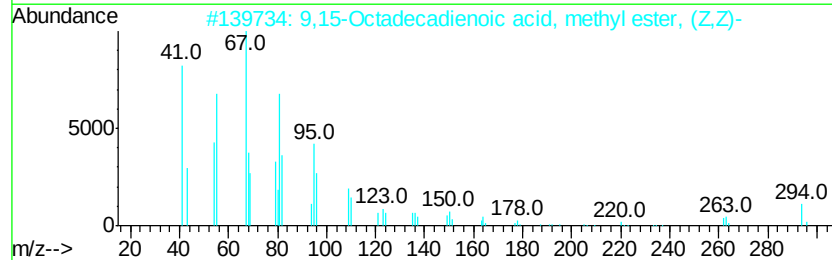
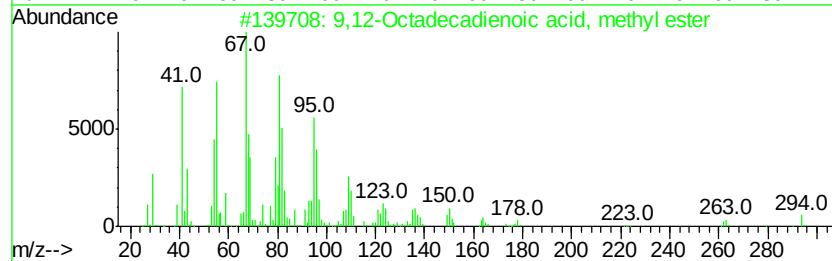
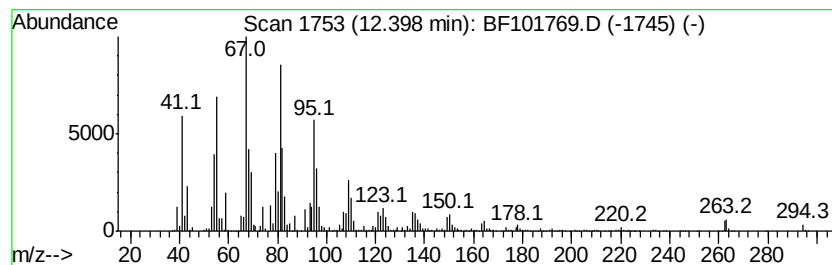
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.40	286.97 ng	15364100	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
2		9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
3		9,12-Octadecadienoic acid, methv...	294	C19H34O2	002566-97-4	99
4		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
5		8,11-Octadecadienoic acid, methy...	294	C19H34O2	056599-58-7	98



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122717\
Data File : BF101769.D
Acq On : 27 Dec 2017 18:22
Operator : SJ/JU
Sample : I6685-09 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

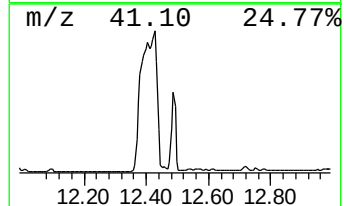
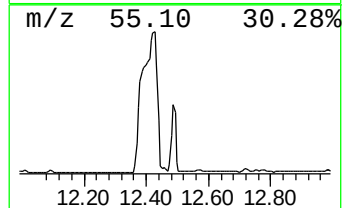
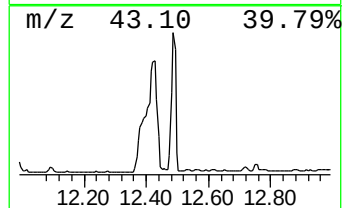
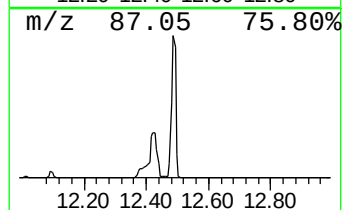
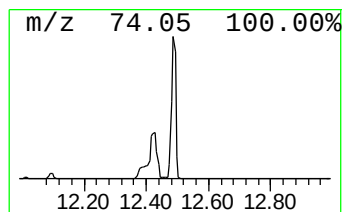
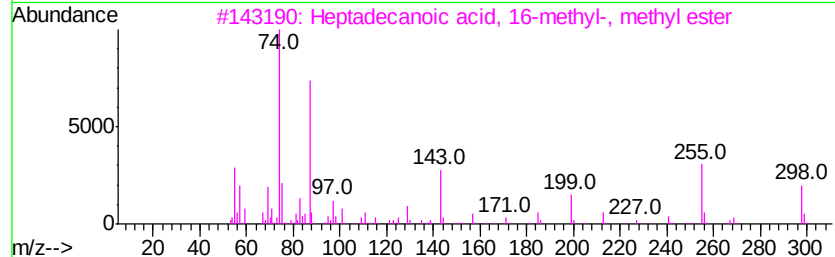
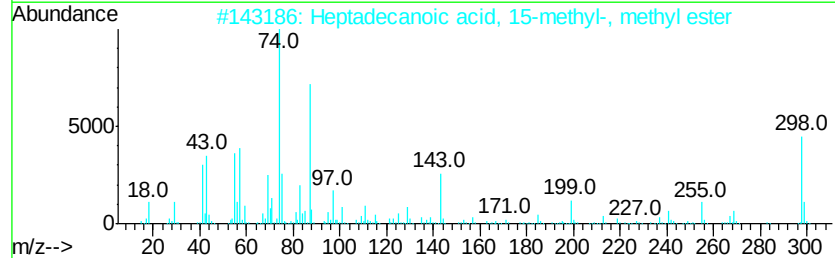
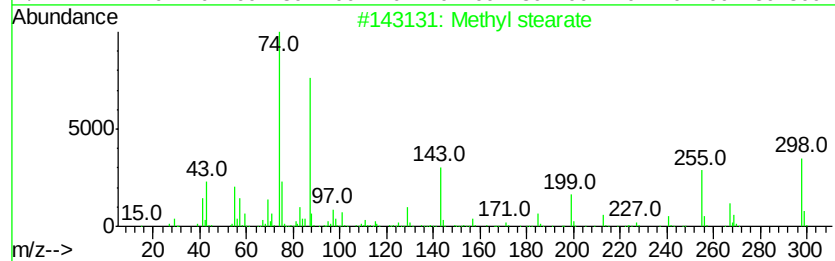
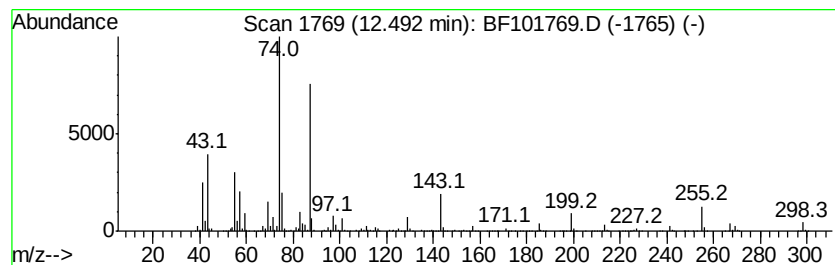
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Methyl stearate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.49	66.09 ng	3538290	Phenanthrene-d10	11.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Methyl stearate	298 C19H38O2	000112-61-8 98
2		Heptadecanoic acid, 15-methyl-, ...	298 C19H38O2	054833-55-5 93
3		Heptadecanoic acid, 16-methyl-, ...	298 C19H38O2	005129-61-3 91
4		Heptadecanoic acid, methyl ester	284 C18H36O2	001731-92-6 91
5		Pentadecanoic acid, methyl ester	256 C16H32O2	007132-64-1 90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122717\
Data File : BF101769.D
Acq On : 27 Dec 2017 18:22
Operator : SJ/JU
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ALS Vial : 10 Sample Multiplier: 1

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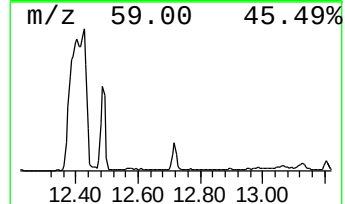
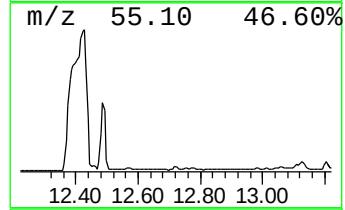
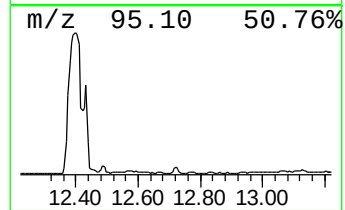
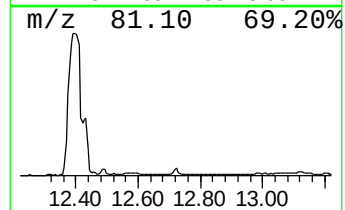
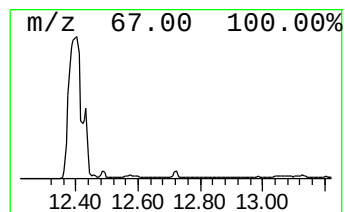
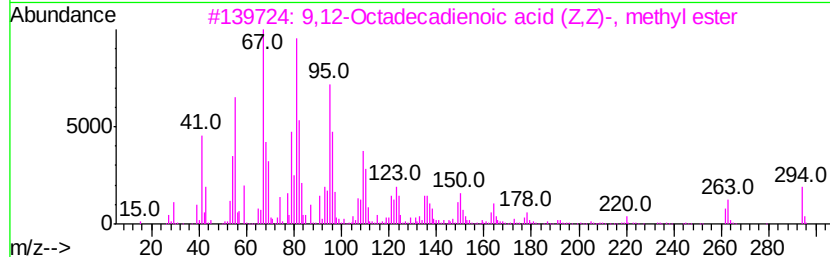
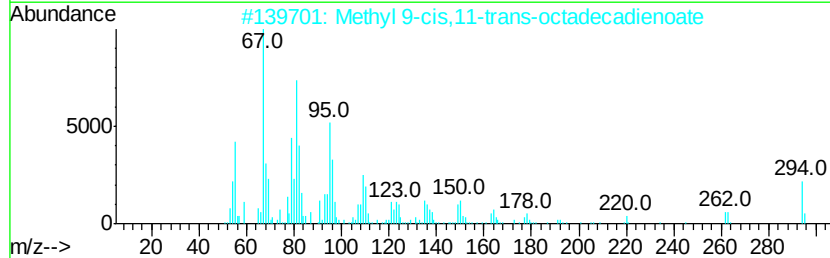
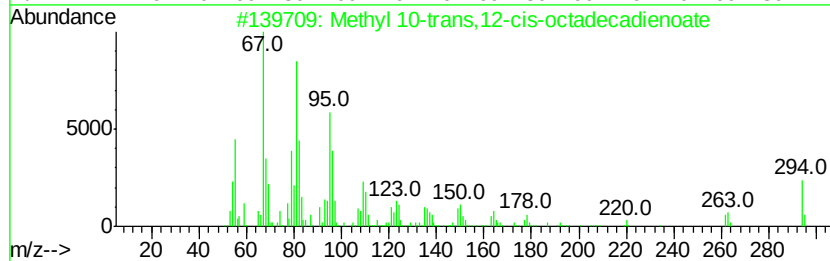
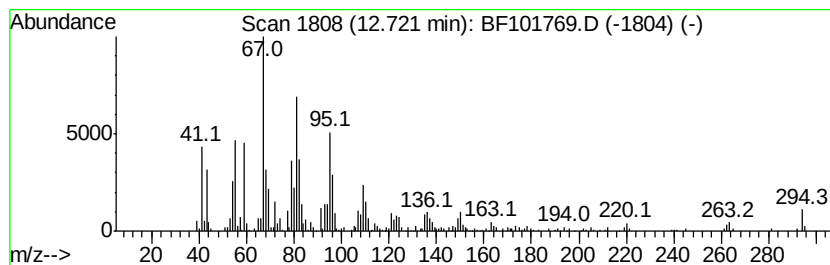
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Methyl 10-trans,12-cis-octa... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.72	5.93 ng	261639	Chrysene-d12	13.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	99
2			Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	1000336-44-0	99
3			9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
4			9,11-Octadecadienoic acid, methy...	294	C19H34O2	013038-47-6	99
5			9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	96



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122717\
Data File : BF101769.D
Acq On : 27 Dec 2017 18:22
Operator : SJ/JU
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Misc :
ALS Vial : 10 Sample Multiplier: 1

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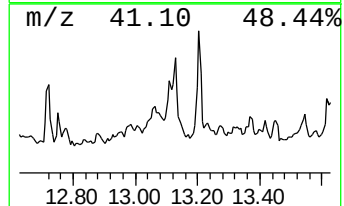
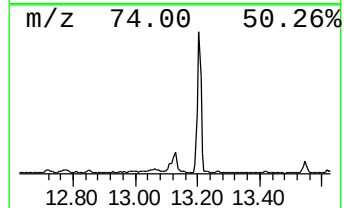
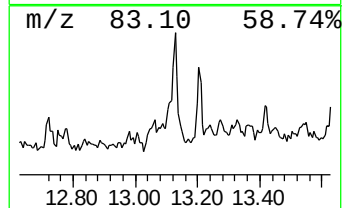
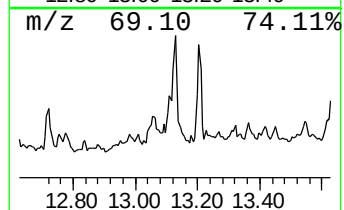
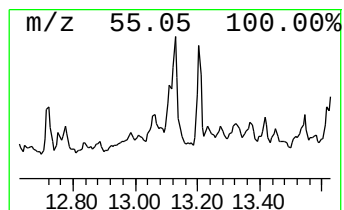
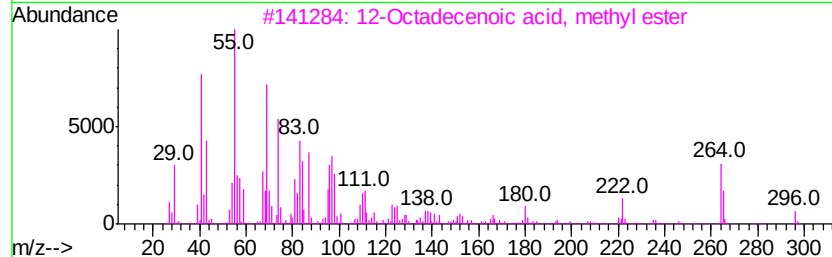
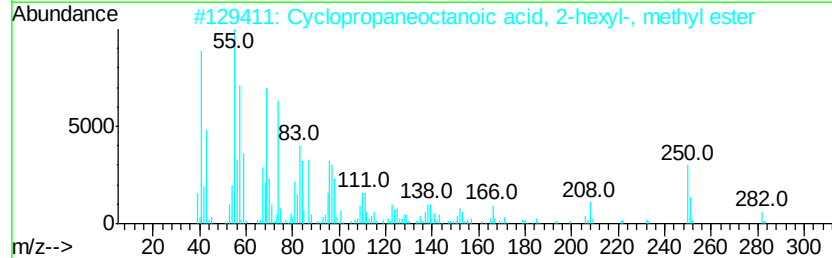
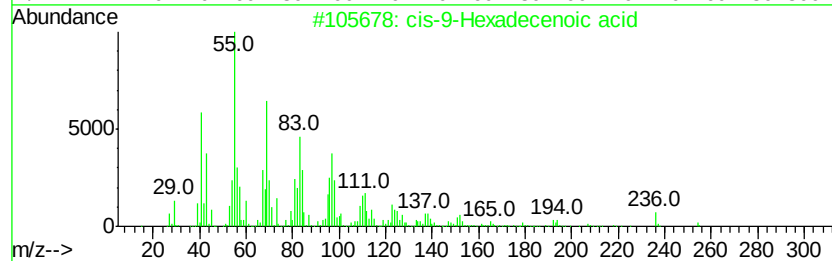
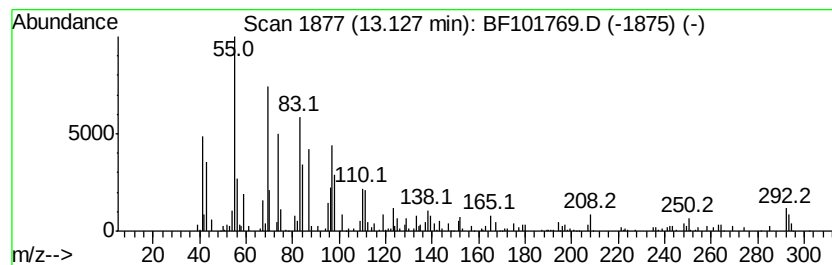
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 cis-9-Hexadecenoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.13	7.22 ng	318395	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	cis-9-Hexadecenoic acid	254	C16H30O2	1000333-19-5	87
2		Cyclopropaneoctanoic acid, 2-hex...	282	C18H34O2	010152-61-1	52
3		12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	50
4		9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	50
5		14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	49



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122717\
Data File : BF101769.D
Acq On : 27 Dec 2017 18:22
Operator : SJ/JU
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Misc :
ALS Vial : 10 Sample Multiplier: 1

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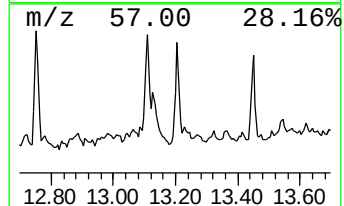
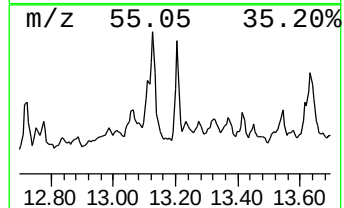
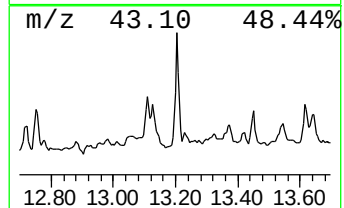
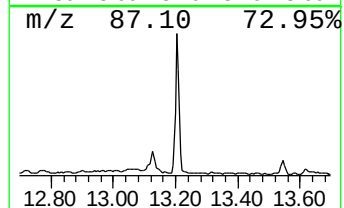
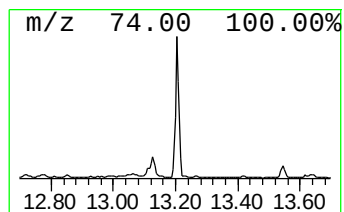
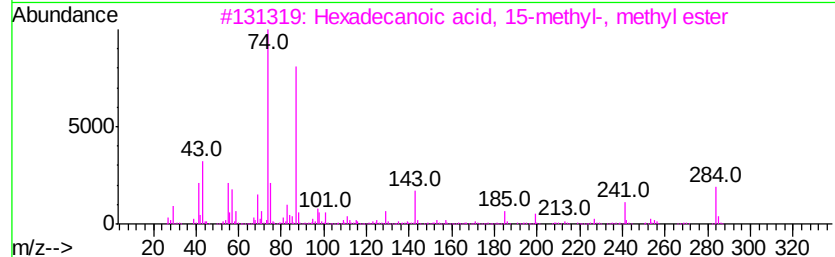
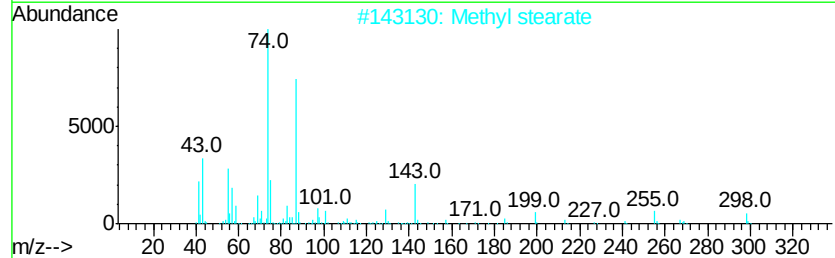
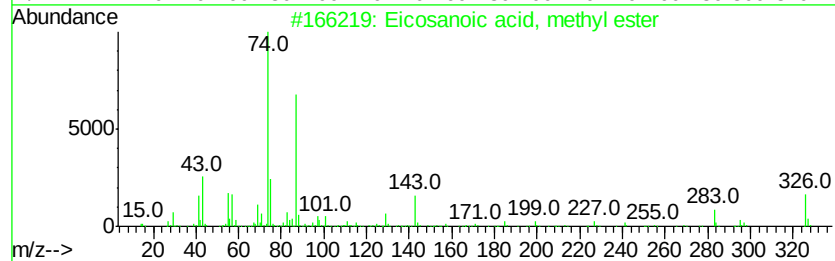
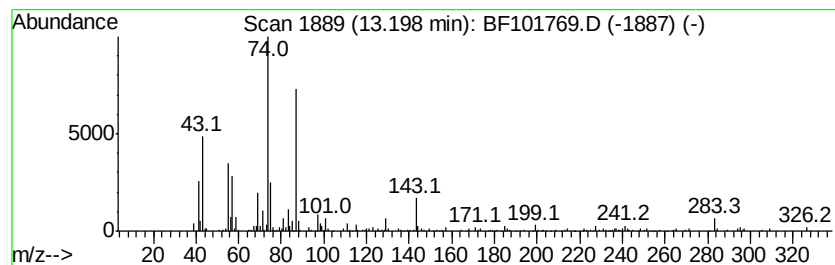
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Eicosanoic acid, methyl ester Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.20	9.30 ng	410502	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosanoic acid, methyl ester	326	C21H42O2	001120-28-1	98
2		Methyl stearate	298	C19H38O2	000112-61-8	96
3		Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	95
4		Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	93
5		Methyl 18-methylnonadecanoate	326	C21H42O2	1000352-20-6	92



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122717\
Data File : BF101769.D
Acq On : 27 Dec 2017 18:22
Operator : SJ/JU
Sample : I6685-09 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

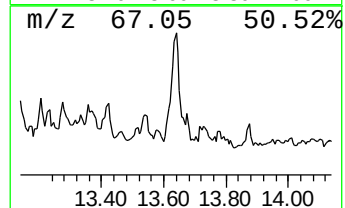
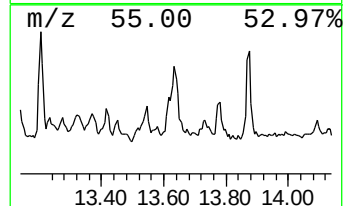
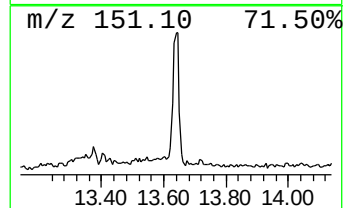
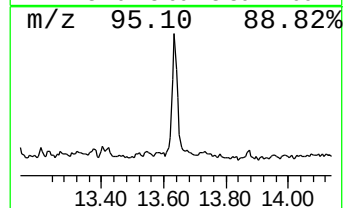
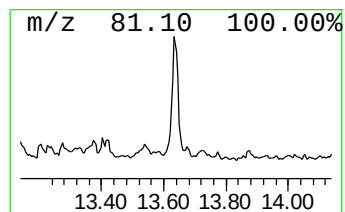
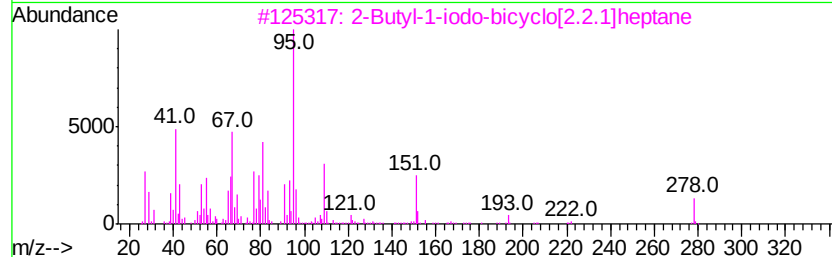
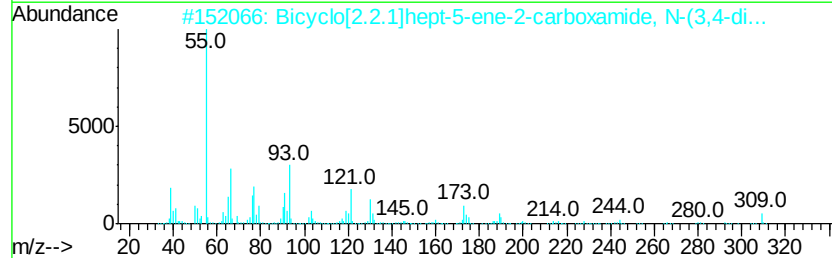
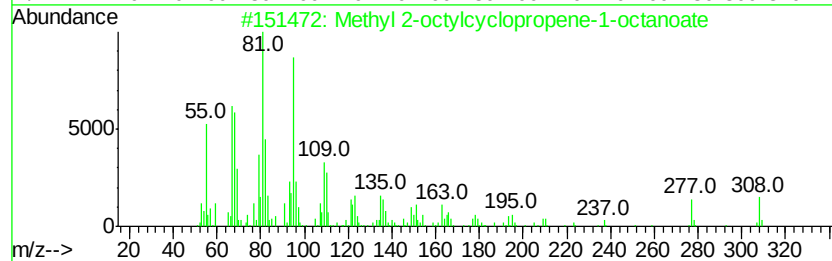
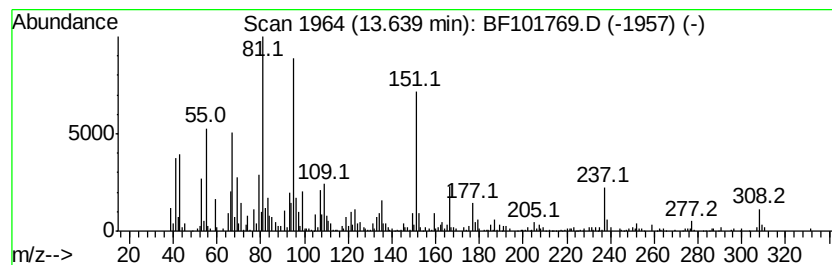
Instrument :
BNA_F
ClientSampleId :
AU-05-122617-C

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Methyl 2-octylcyclopropene-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.64	13.65 ng	602207	Chrysene-d12	13.97		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methyl 2-octylcyclopropene-1-oct...	308	C20H36O2	003220-60-8	50
2		Bicyclo[2.2.1]hept-5-ene-2-carbo...	309	C18H19N3O2	1000265-40-3	46
3		2-Butyl-1-iodo-bicyclo[2.2.1]hep...	278	C11H19I	1000190-59-7	43
4		2(3H)-Benzofuranone, 3a,4,5,7a-t...	166	C10H14O2	033722-72-4	25
5		2-n-Heptylfuran	166	C11H18O	003777-71-7	18



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122717\
Data File : BF101769.D
Acq On : 27 Dec 2017 18:22
Operator : SJ/JU
Sample : I6685-09 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AU-05-122617-C

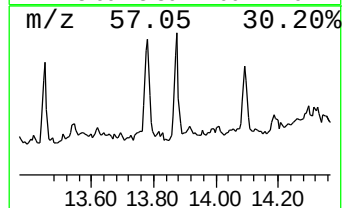
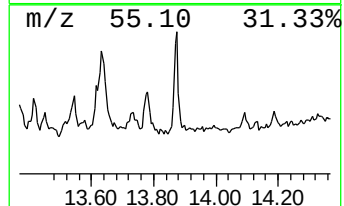
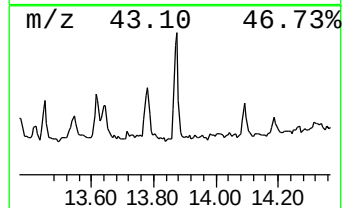
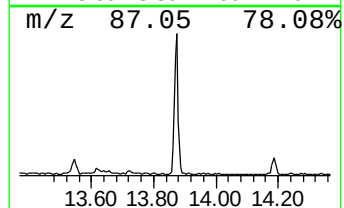
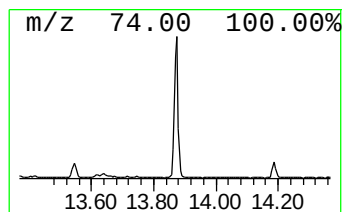
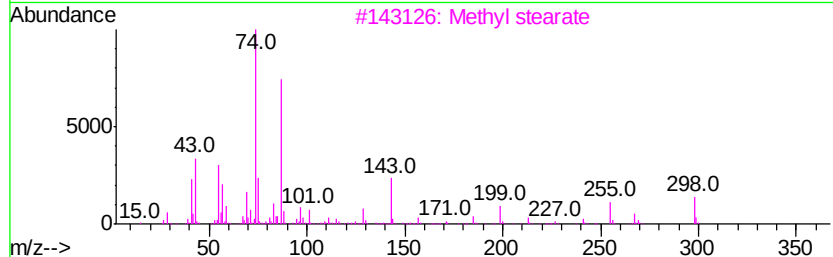
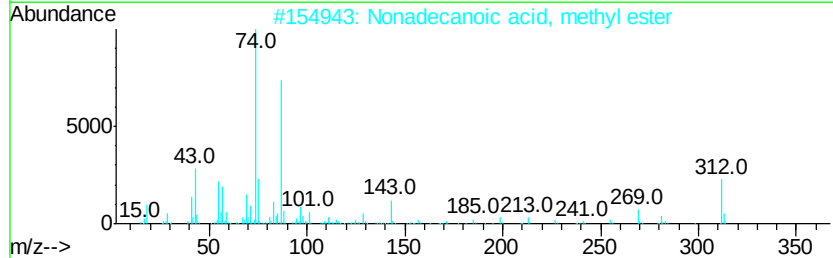
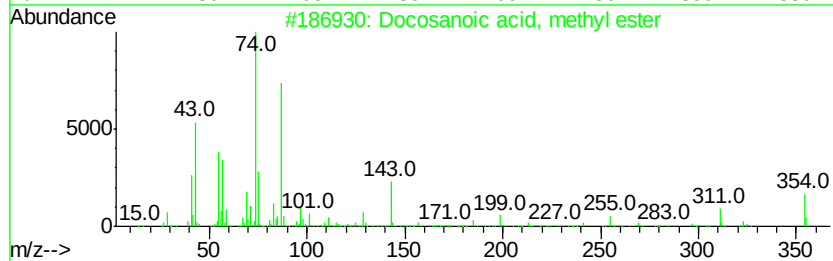
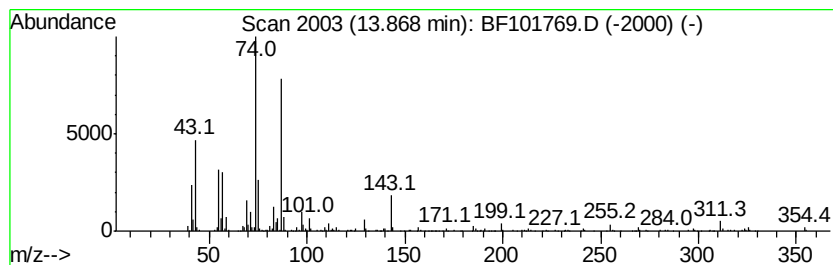
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Docosanoic acid, methyl ester Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	8.46 ng	373282	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosanoic acid, methyl ester	354	C23H46O2	000929-77-1	98
2		Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	96
3		Methyl stearate	298	C19H38O2	000112-61-8	87
4		Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	86
5		Heneicosanoic acid, methyl ester	340	C22H44O2	006064-90-0	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122717\
 Data File : BF101769.D
 Acq On : 27 Dec 2017 18:22
 Operator : SJ/JU
 Sample : I6685-09 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-05-122617-C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.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...	5.00	4.5	ng	221830	1	6.79	992990	20.0
unknown6.56	6.56	40.1	ng	1990630	1	6.79	992990	20.0
Tetradecane	8.65	3.3	ng	202658	2	8.07	1217600	20.0
Hexadecane	9.74	4.5	ng	266422	3	9.83	1192690	20.0
Tetratetracontane	10.71	7.6	ng	405442	4	11.32	1070790	20.0
Hexadecanoic acid...	11.70	127.6	ng	6829450	4	11.32	1070790	20.0
n-Hexadecanoic acid	11.83	3.4	ng	181244	4	11.32	1070790	20.0
9,12-Octadecadien...	12.40	287.0	ng	15364100	4	11.32	1070790	20.0
Methyl stearate	12.49	66.1	ng	3538290	4	11.32	1070790	20.0
Methyl 10-trans,1...	12.72	5.9	ng	261639	5	13.97	882351	20.0
cis-9-Hexadeceno...	13.13	7.2	ng	318395	5	13.97	882351	20.0
Eicosanoic acid, ...	13.20	9.3	ng	410502	5	13.97	882351	20.0
Methyl 2-octylcyc...	13.64	13.7	ng	602207	5	13.97	882351	20.0
Docosanoic acid, ...	13.87	8.5	ng	373282	5	13.97	882351	20.0