

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102817\
 Data File : BF099945.D
 Acq On : 29 Oct 2017 6:04
 Operator : SJ/JU
 Sample : I6059-04
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 101817-TIDO-F-U-S

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-BF102317.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.998	493	495	508	rBV	110216	159965	5.58%	0.870%
2	5.410	562	565	587	rBV	731059	893304	31.14%	4.860%
3	6.428	735	738	753	rBV	632181	762265	26.58%	4.147%
4	6.557	757	760	767	rVB	1675326	1517820	52.92%	8.258%
5	6.787	795	799	808	rBV	681342	614833	21.44%	3.345%
6	6.939	821	825	842	rBV	2105775	2045029	71.30%	11.126%
7	7.357	892	896	916	rBV	1469085	1536951	53.59%	8.362%
8	8.069	1013	1017	1028	rBV	963131	822784	28.69%	4.477%
9	8.628	1109	1112	1124	rBV	135616	119678	4.17%	0.651%
10	9.145	1195	1200	1203	rBV	2973705	2868225	100.00%	15.605%
11	9.539	1264	1267	1274	rBV	244366	202912	7.07%	1.104%
12	9.822	1311	1315	1325	rVB	985604	822477	28.68%	4.475%
13	10.492	1425	1429	1433	rBV	50759	39054	1.36%	0.212%
14	10.533	1433	1436	1441	rBV	494590	402041	14.02%	2.187%
15	10.616	1447	1450	1465	rBV	848760	773405	26.96%	4.208%
16	11.316	1565	1569	1576	rVB	673805	642726	22.41%	3.497%
17	12.698	1801	1804	1808	rVB	146348	127862	4.46%	0.696%
18	12.898	1833	1838	1841	rBV	2008116	1871846	65.26%	10.184%
19	13.404	1921	1924	1927	rBV2	106770	85042	2.96%	0.463%
20	13.774	1983	1987	1998	rVB	71931	77627	2.71%	0.422%
21	13.963	2015	2019	2027	rVB	592008	524032	18.27%	2.851%
22	15.021	2195	2199	2203	rBV	47930	50328	1.75%	0.274%
23	15.386	2255	2261	2264	rBV	86404	104695	3.65%	0.570%
24	15.421	2264	2267	2276	rVB	407751	510150	17.79%	2.776%
25	15.804	2327	2332	2337	rVB	97138	131746	4.59%	0.717%
26	16.286	2408	2414	2422	rBV	107595	181918	6.34%	0.990%
27	16.851	2504	2510	2522	rVB	77982	152708	5.32%	0.831%
28	17.521	2614	2624	2632	rBV2	65389	156038	5.44%	0.849%
29	18.321	2752	2760	2773	rVB2	36144	105496	3.68%	0.574%
30	19.262	2911	2920	2929	rBV3	23604	77091	2.69%	0.419%

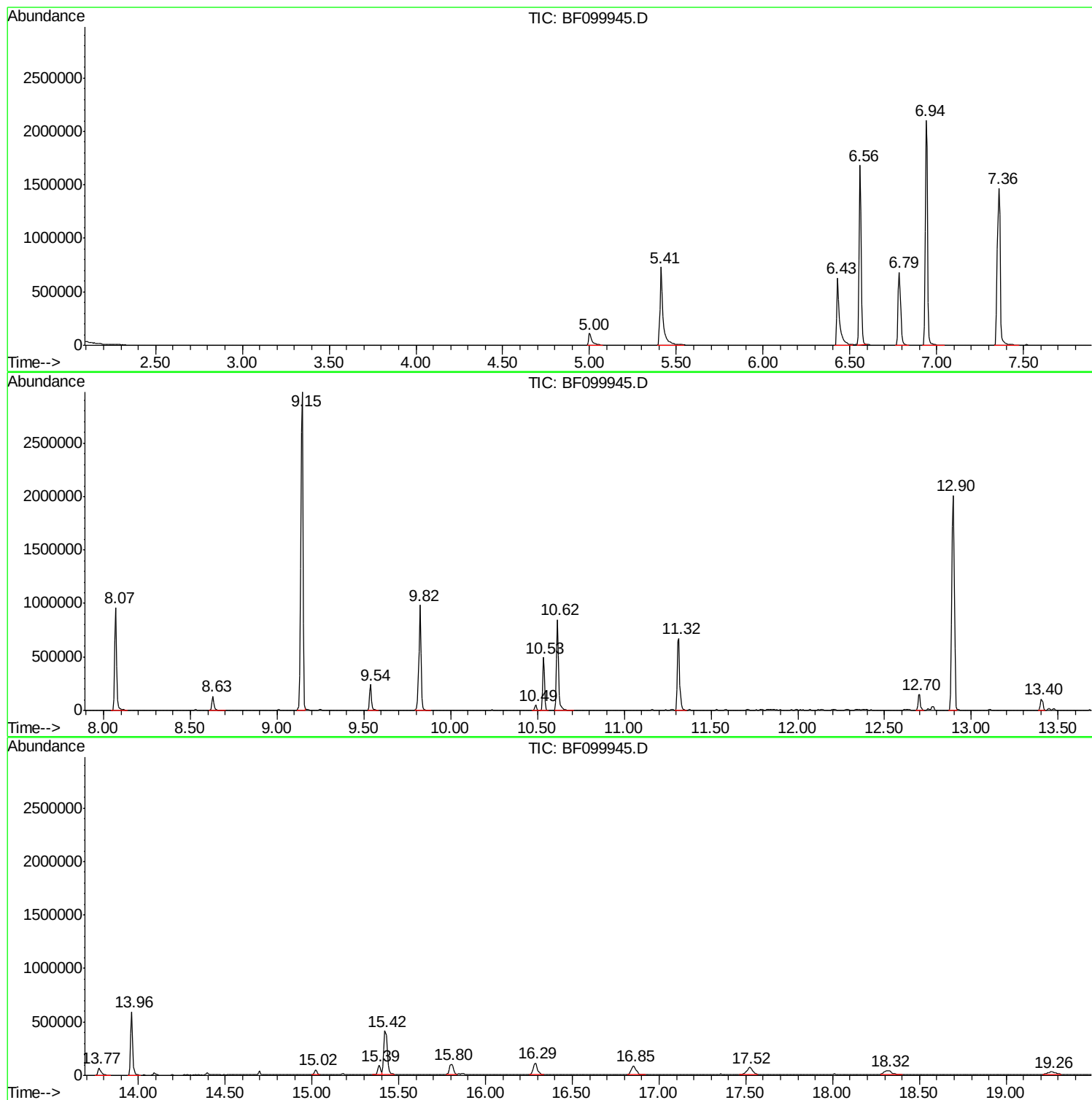
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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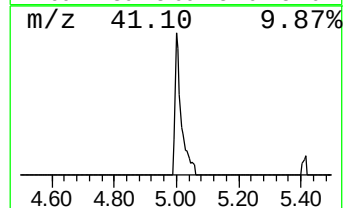
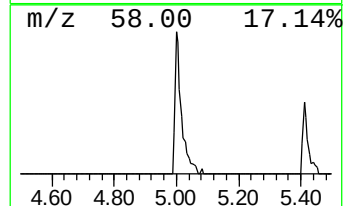
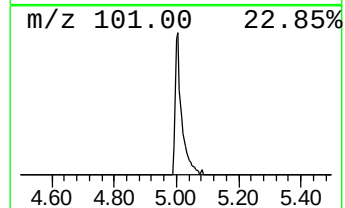
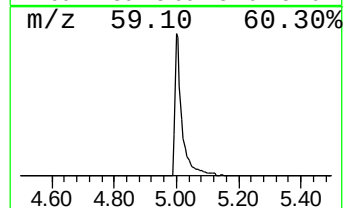
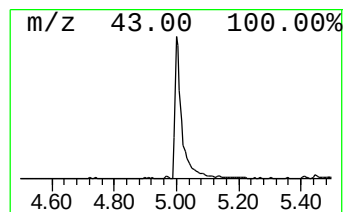
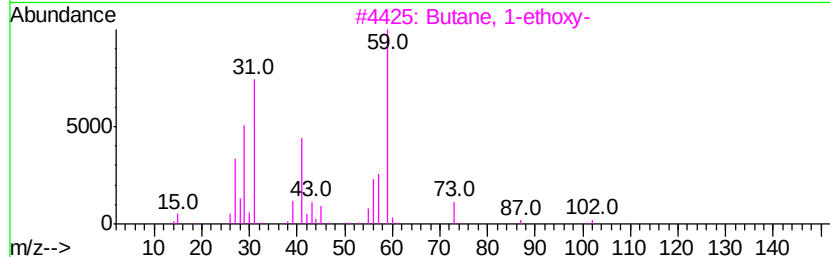
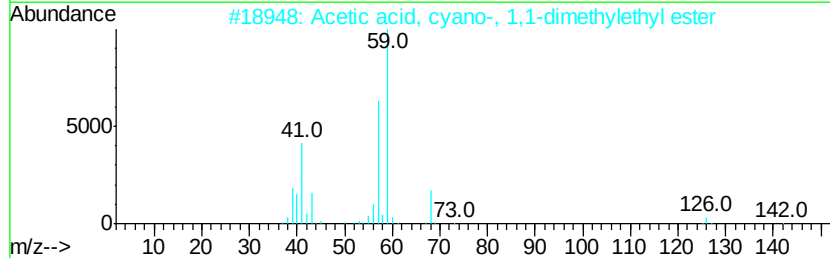
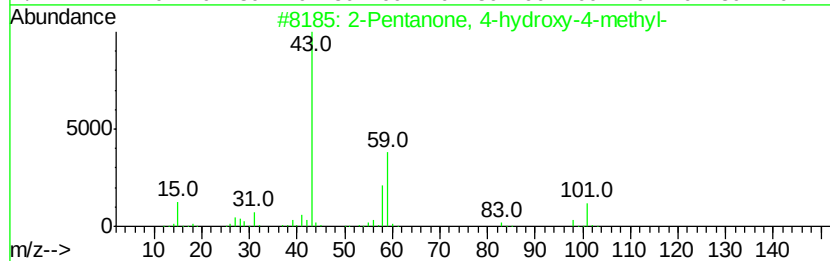
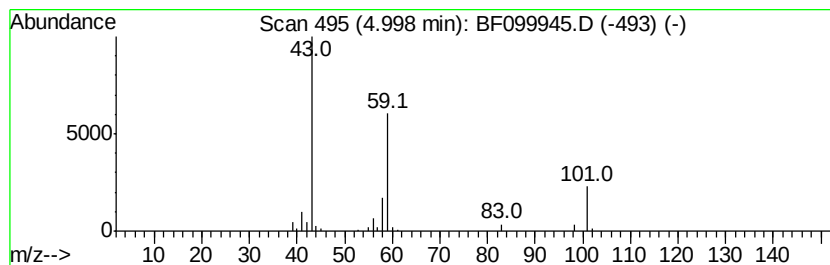
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	5.20 ng	159965	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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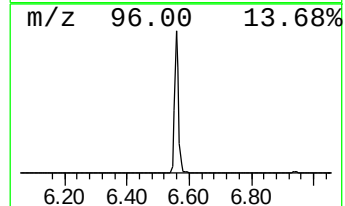
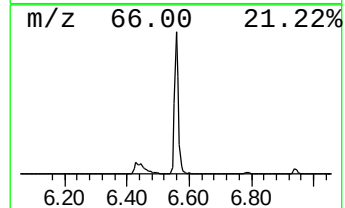
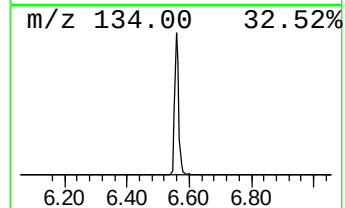
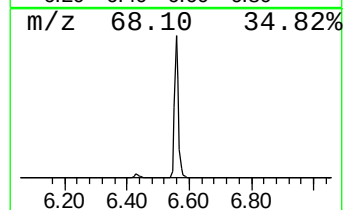
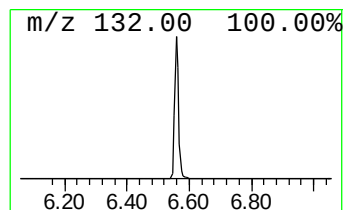
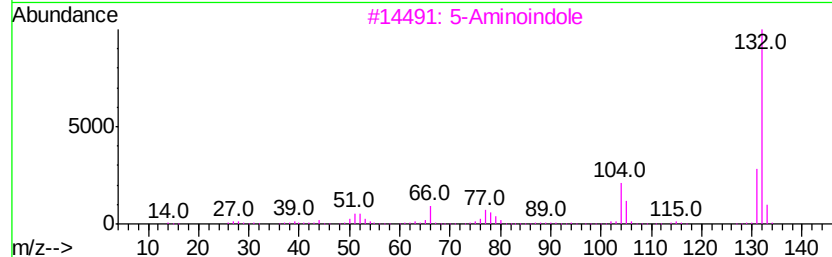
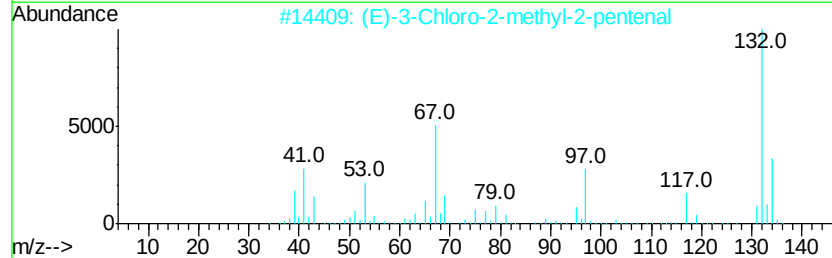
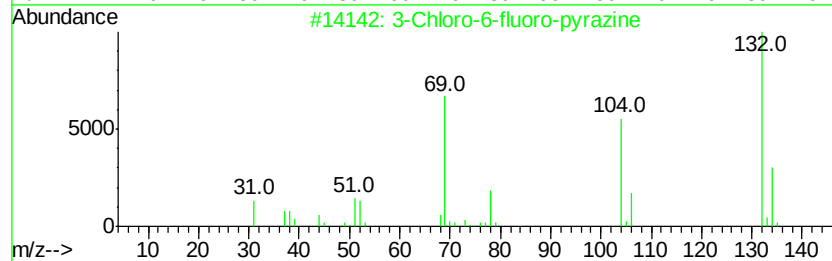
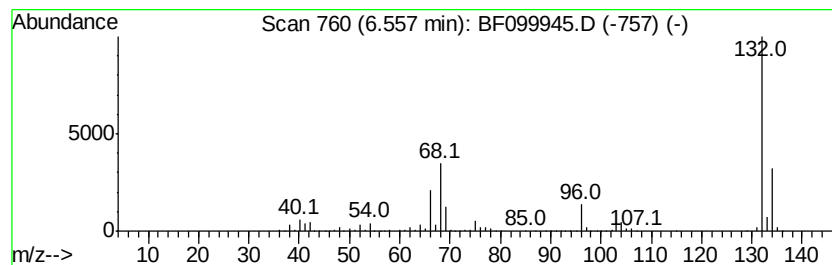
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Peak Number 2 unknown6.56 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	49.37 ng	1517820	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	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	10



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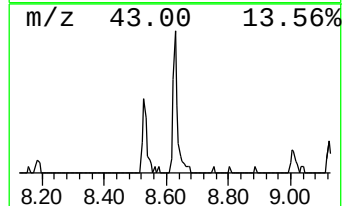
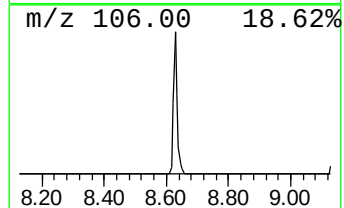
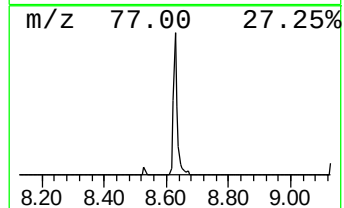
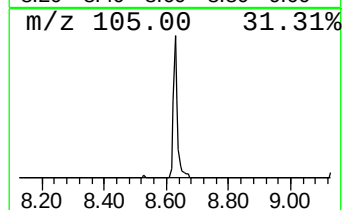
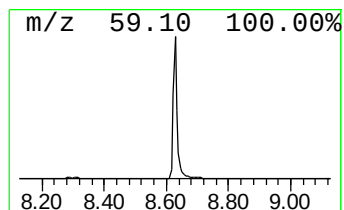
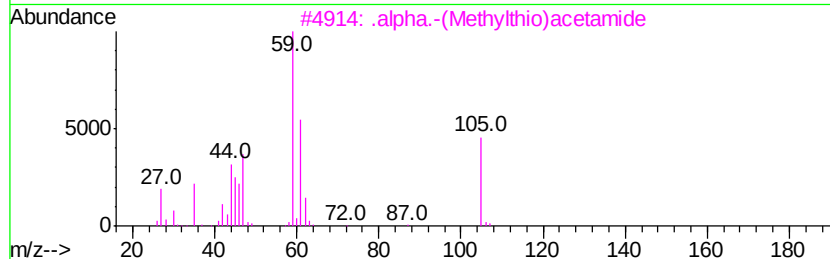
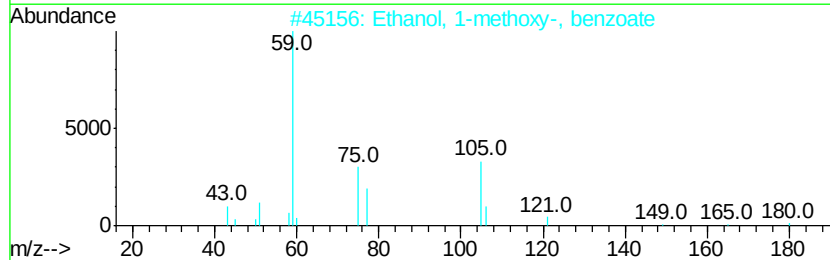
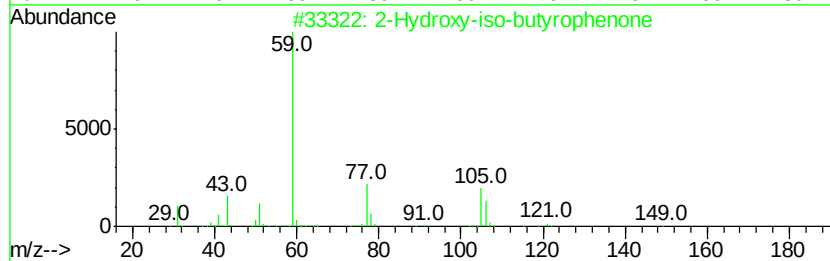
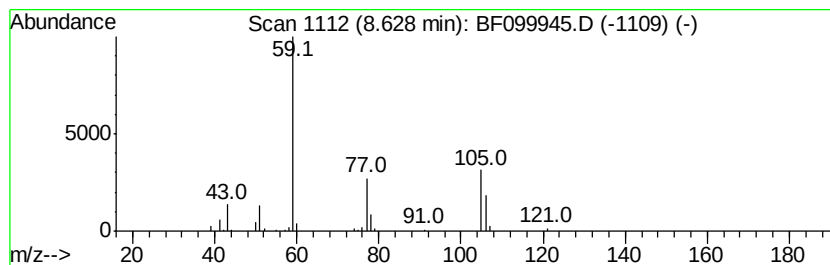
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Peak Number 4 2-Hydroxy-iso-butyrophenone Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.63	2.91 ng	119678	Napthalene-d8	8.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Hydroxy-iso-butyrophenone	164	C10H12O2	007473-98-5	86
2		Ethanol, 1-methoxy-, benzoate	180	C10H12O3	051835-44-0	50
3		.alpha.-(Methylthio)acetamide	105	C3H7NOS	022551-24-2	7
4		Formamide, N-methyl-	59	C2H5NO	000123-39-7	7
5		Guanidine	59	CH5N3	000113-00-8	7



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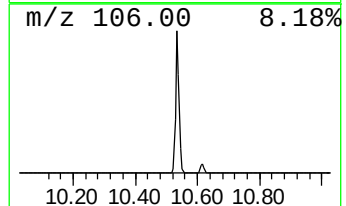
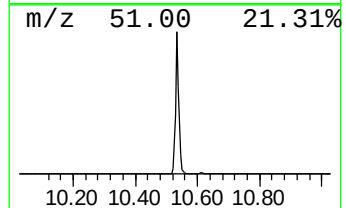
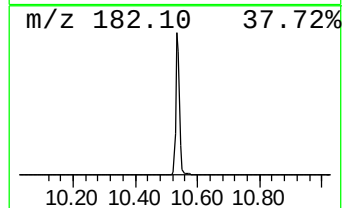
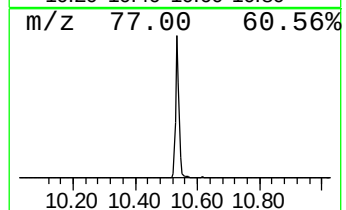
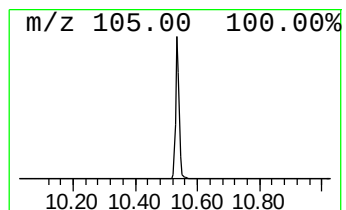
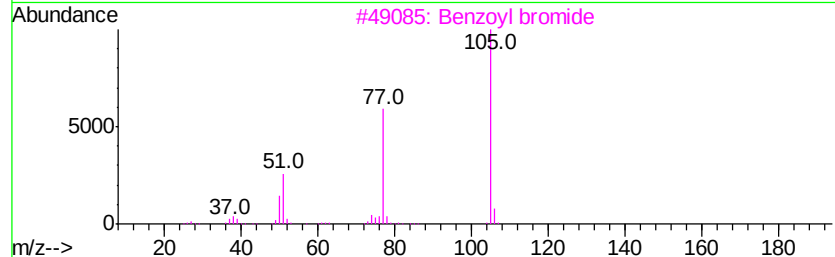
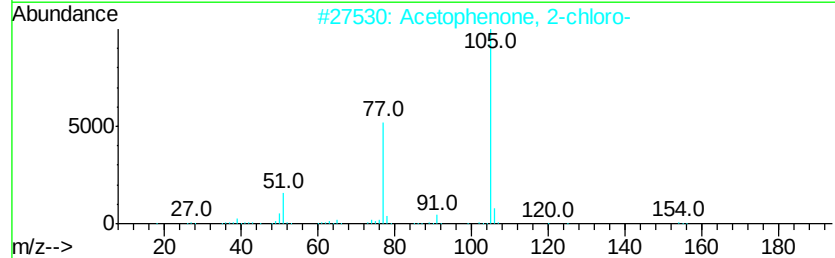
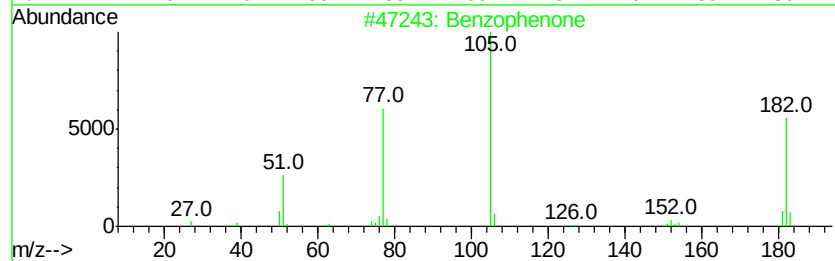
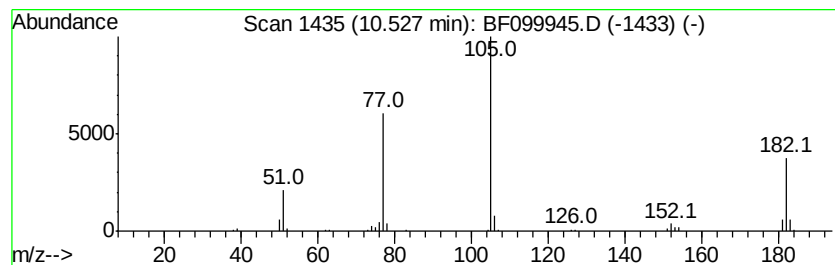
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Peak Number 5 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.53	9.78 ng	402041	Acenaphthene-d10	9.82
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzophenone	182 C13H10O	000119-61-9 96
2		Acetophenone, 2-chloro-	154 C8H7ClO	000532-27-4 49
3		Benzoyl bromide	184 C7H5BrO	000618-32-6 47
4		Vinyl benzoate	148 C9H8O2	000769-78-8 47
5		S-Benzoyl(thiohydroxylamine)	153 C7H7NOS	025740-80-1 47



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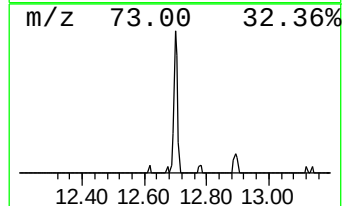
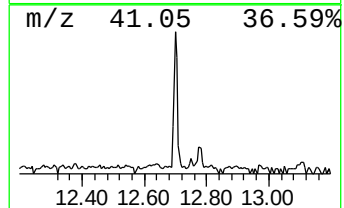
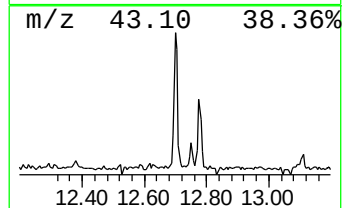
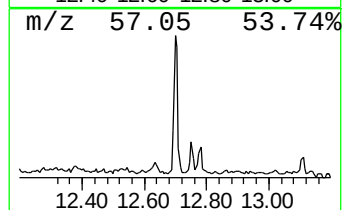
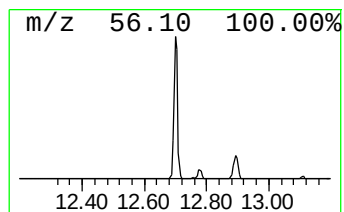
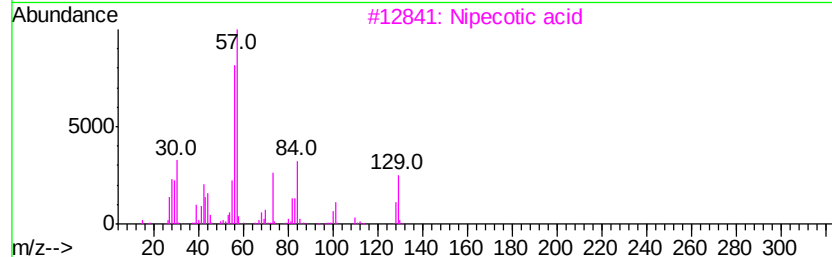
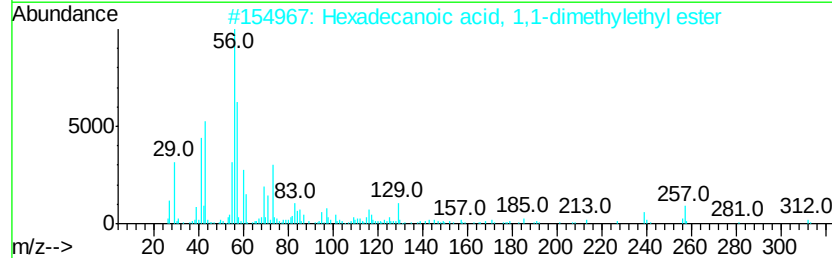
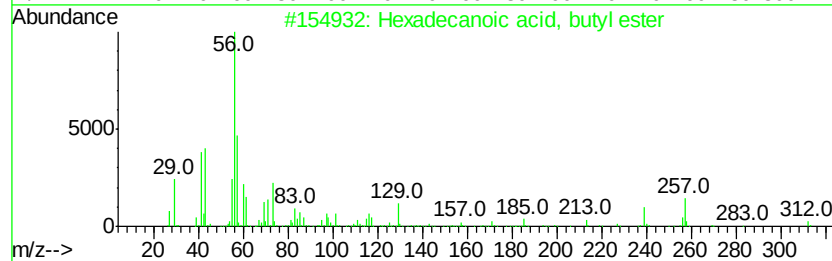
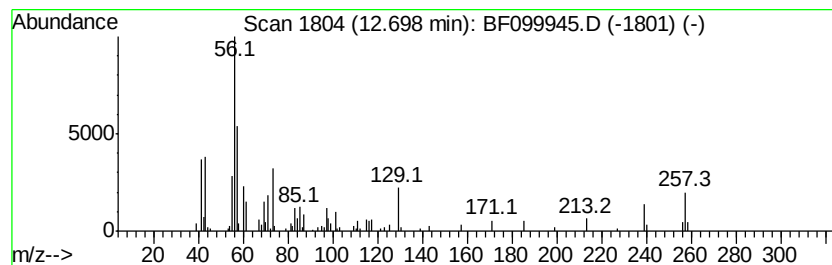
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Peak Number 6 Hexadecanoic acid, butyl ester Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	4.88 ng	127862	Chrysene-d12	13.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	91
2		Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	91
3		Nipecotic acid	129	C6H11NO2	000498-95-3	47
4		Cyclopentanecarboxylic acid, 2-a...	129	C6H11NO2	037910-65-9	38
5		1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	30



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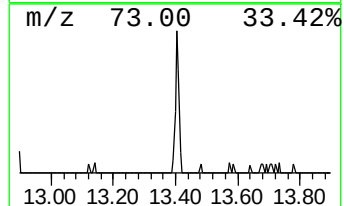
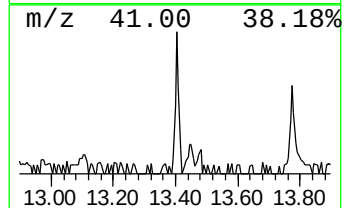
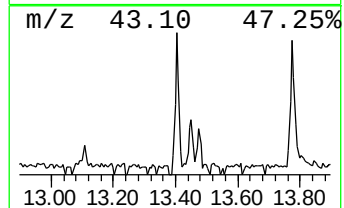
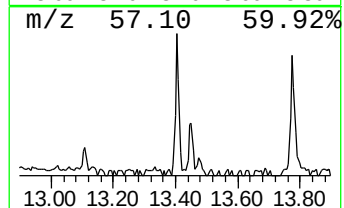
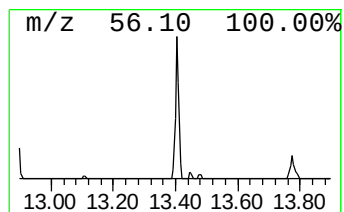
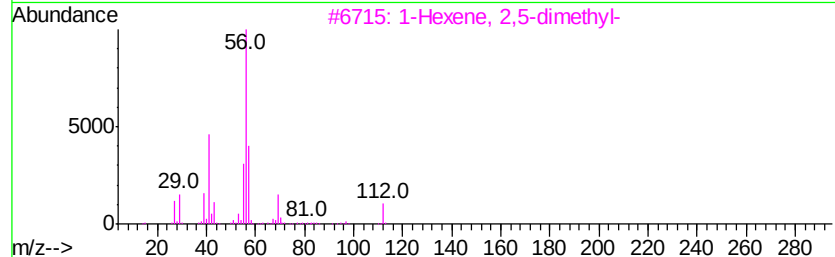
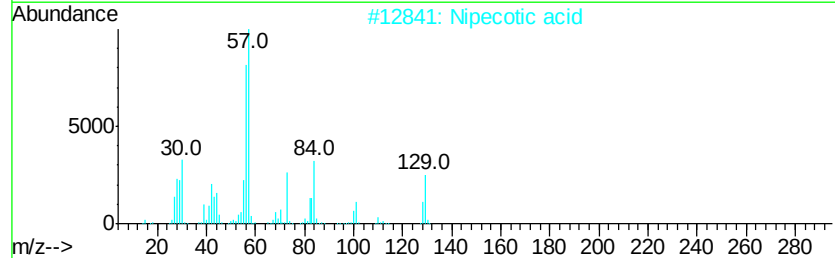
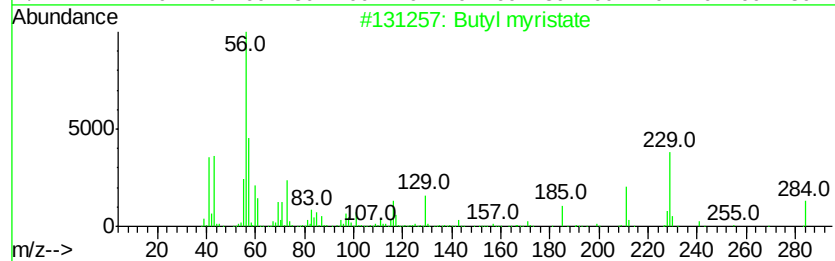
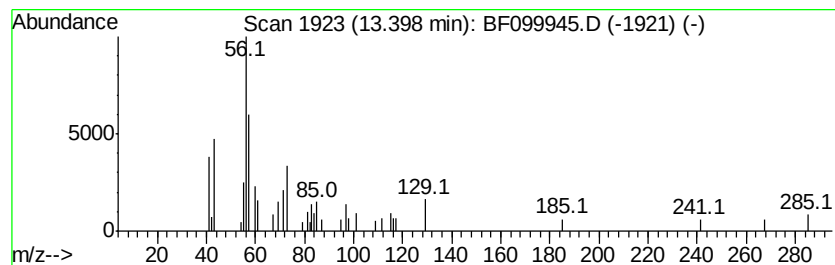
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ClientSampleId :
101817-TIDO-F-U-S

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Butyl myristate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.40	3.25 ng	85042	Chrysene-d12		13.96
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl myristate	284	C18H36O2	000110-36-1	59
2	Nipecotic acid	129	C6H11NO2	000498-95-3	50
3	1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	30
4	Cyclohexanol, 4-amino-, trans-	115	C6H13NO	027489-62-9	27
5	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	27



Data Path : Z:\HPCHEM1\BNA F\DATA\BF102817\
Data File : BF099945.D
Acq On : 29 Oct 2017 6:04
Operator : SJ/JU
Sample : I6059-04
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
101817-TIDO-F-U-S

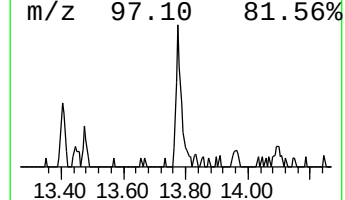
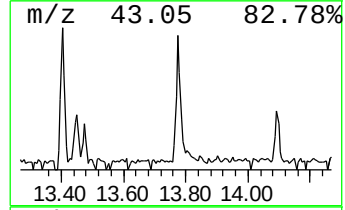
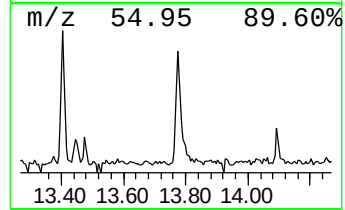
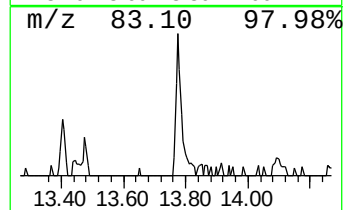
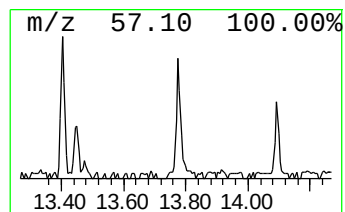
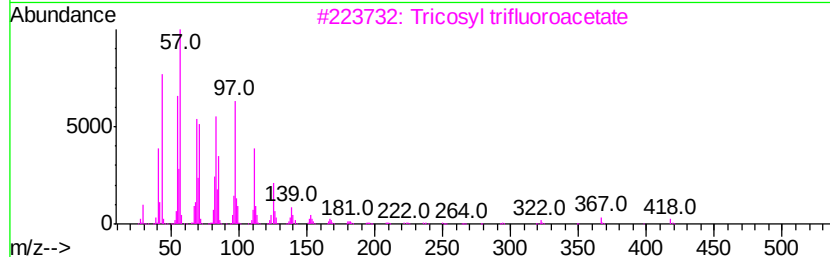
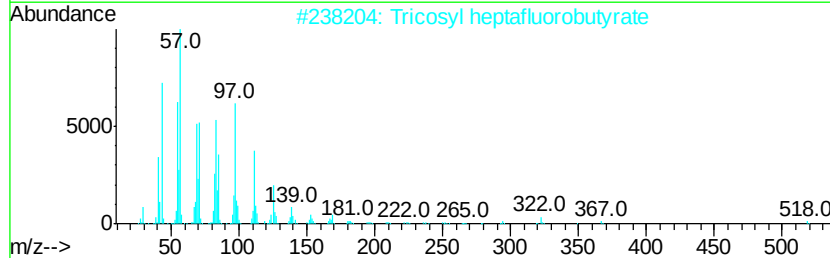
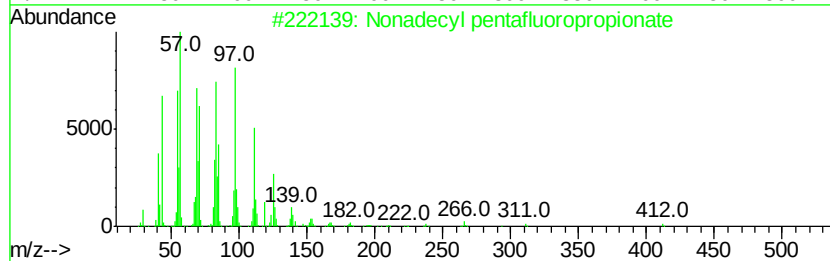
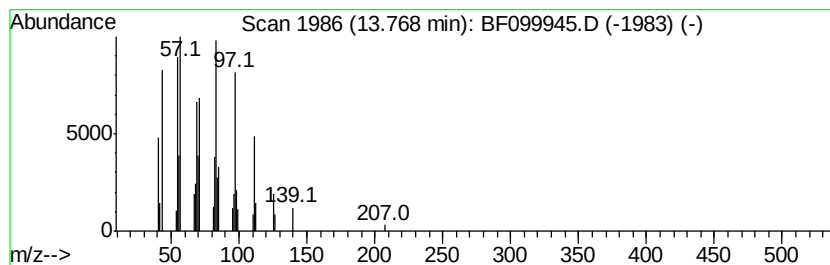
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Nonadecyl pentafluoropropio... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	2.96 ng	77627	Chrysene-d12	13.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
2		Tricosyl heptafluorobutyrate	536	C27H47F7O2	1000351-83-4	91
3		Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	91
4		Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	91
5		Heneicosyl heptafluorobutyrate	508	C25H43F7O2	1000351-83-8	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102817\
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Acq On : 29 Oct 2017 6:04
Operator : SJ/JU
Sample : I6059-04
Misc :
ALS Vial : 35 Sample Multiplier: 1

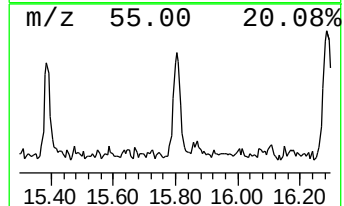
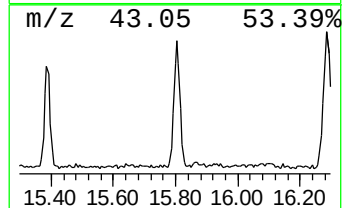
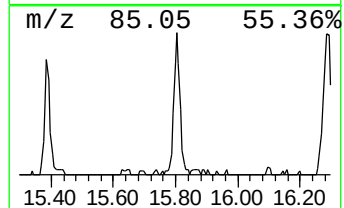
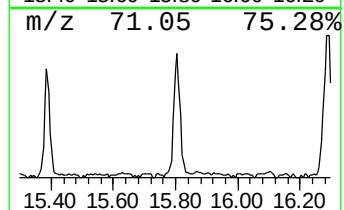
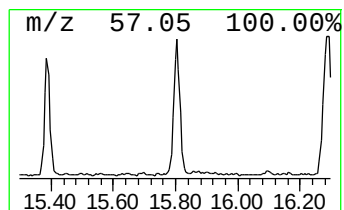
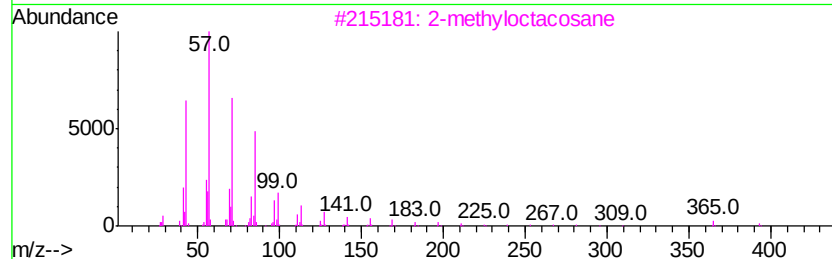
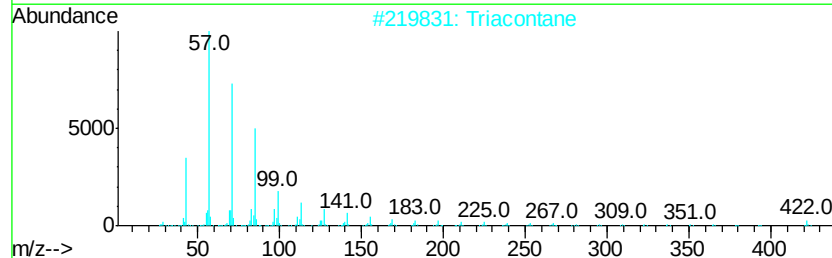
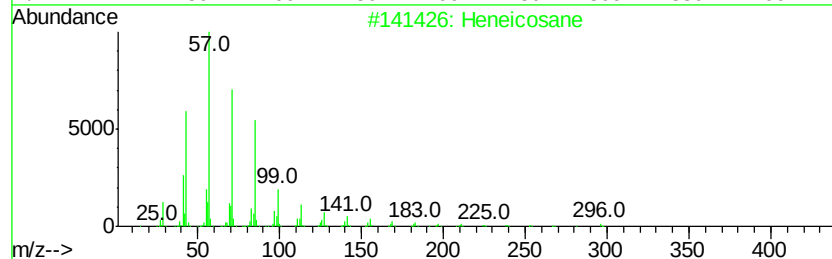
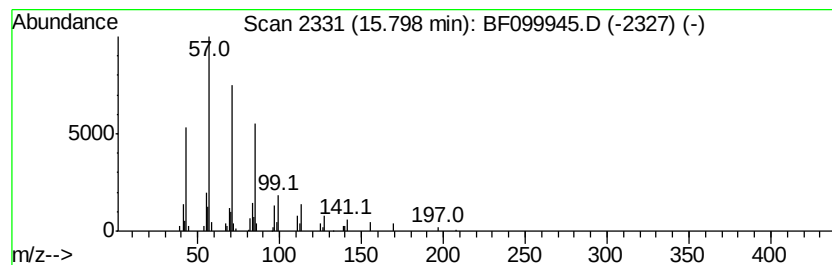
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.80	5.16 ng	131746	Perylene-d12	15.42
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heneicosane	296 C21H44	000629-94-7	91
2	Triacontane	422 C30H62	000638-68-6	91
3	2-methyloctacosane	408 C29H60	1000376-72-8	91
4	Hexadecane, 1-iodo-	352 C16H33I	000544-77-4	90
5	2-methyltetracosane	352 C25H52	1000376-72-6	90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF102817\
Data File : BF099945.D
Acq On : 29 Oct 2017 6:04
Operator : SJ/JU
Sample : I6059-04
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
101817-TIDO-F-U-S

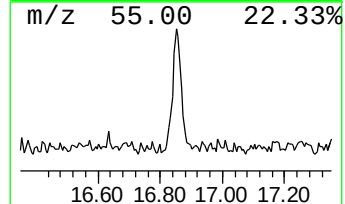
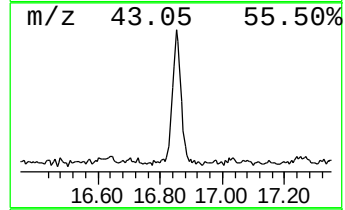
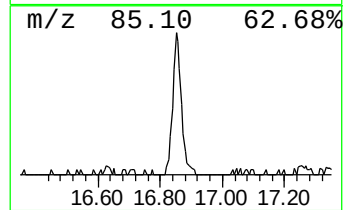
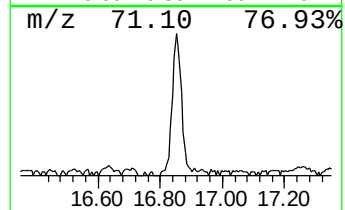
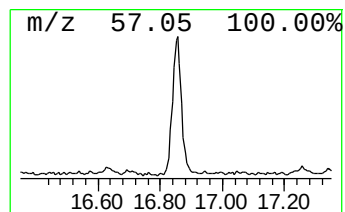
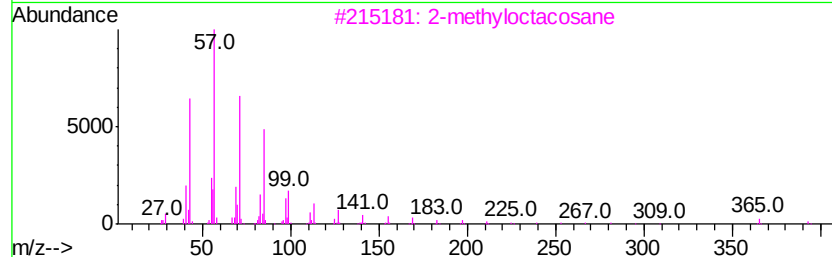
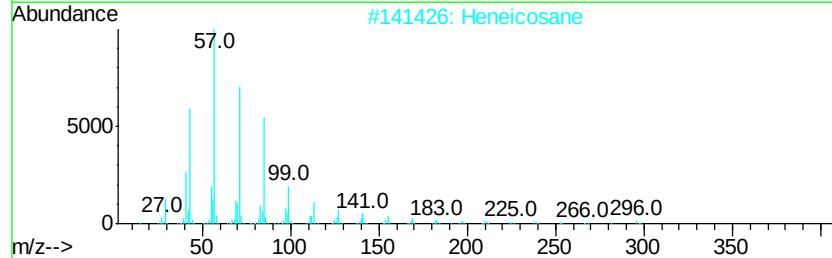
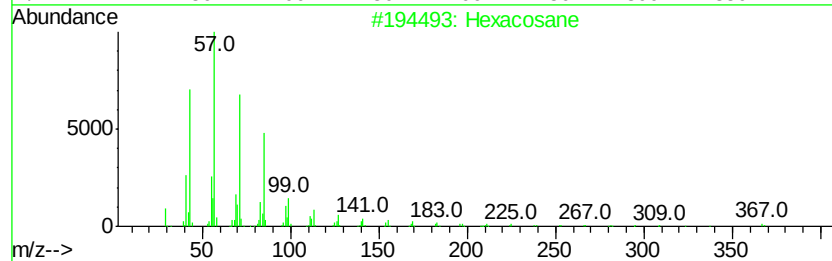
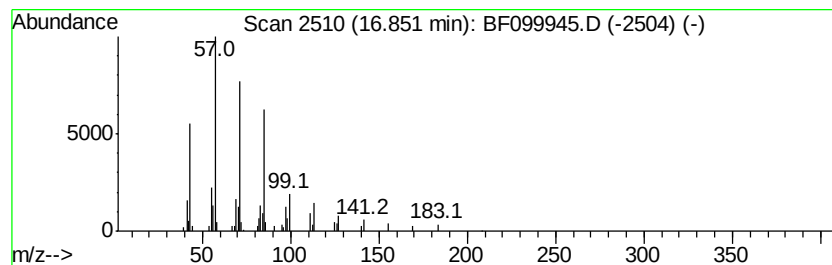
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.85	5.99 ng	152708	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	90
4		Heptadecane	240	C17H36	000629-78-7	87
5		10-Methylnonadecane	282	C20H42	056862-62-5	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF102817\
 Data File : BF099945.D
 Acq On : 29 Oct 2017 6:04
 Operator : SJ/JU
 Sample : I6059-04
 Misc :
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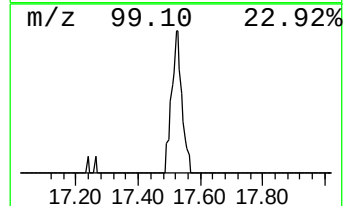
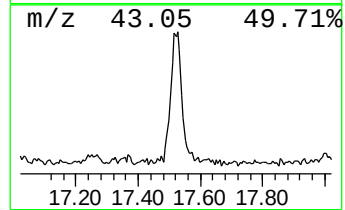
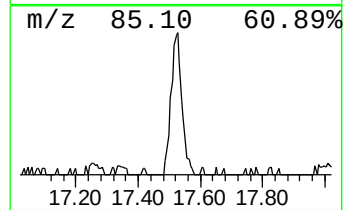
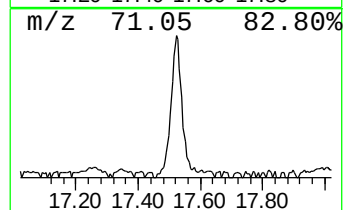
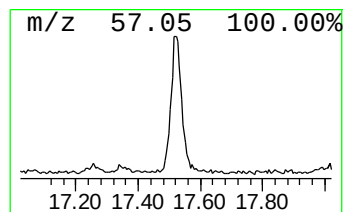
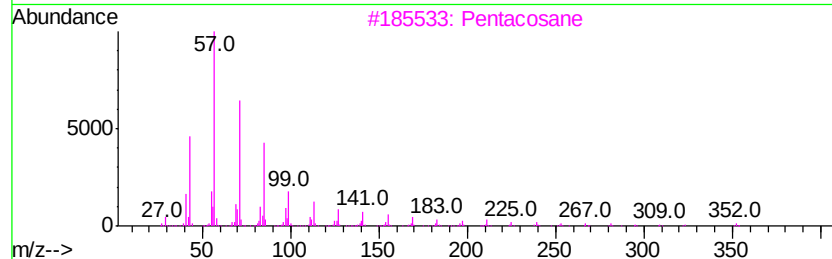
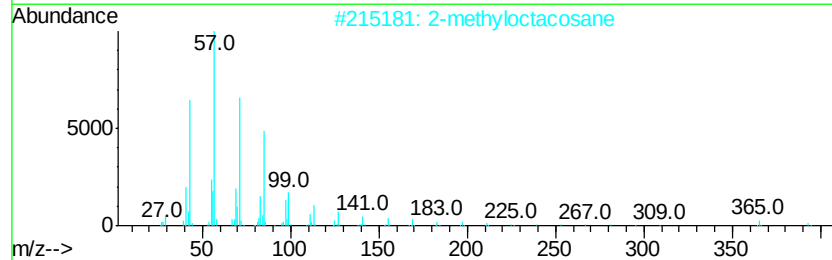
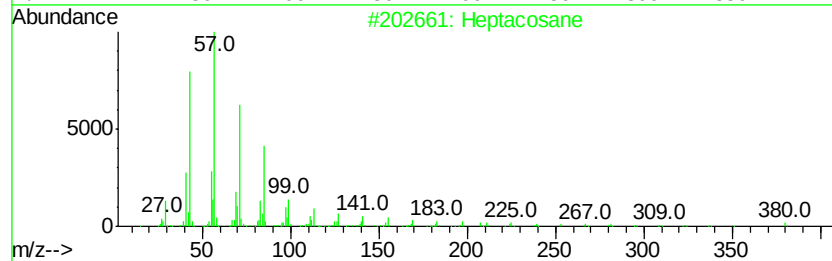
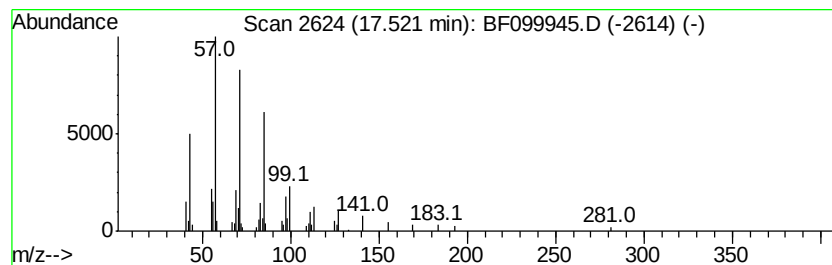
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 12 Heptacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.52	6.12 ng	156038	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	90
3		Pentacosane	352	C25H52	000629-99-2	90
4		Hexacosane	366	C26H54	000630-01-3	87
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102817\
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Acq On : 29 Oct 2017 6:04
Operator : SJ/JU
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Misc :
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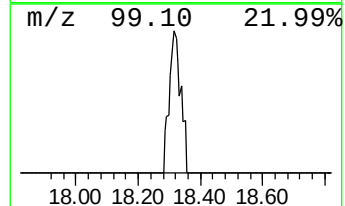
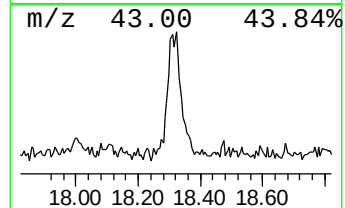
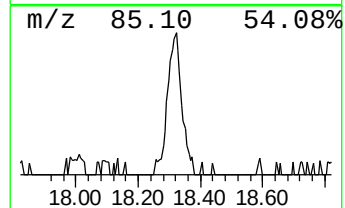
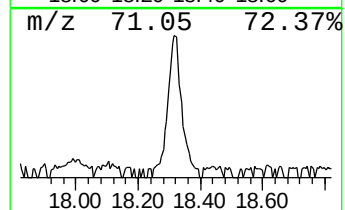
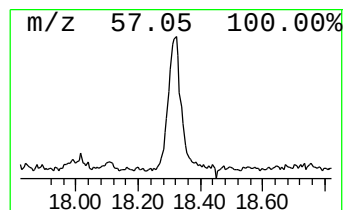
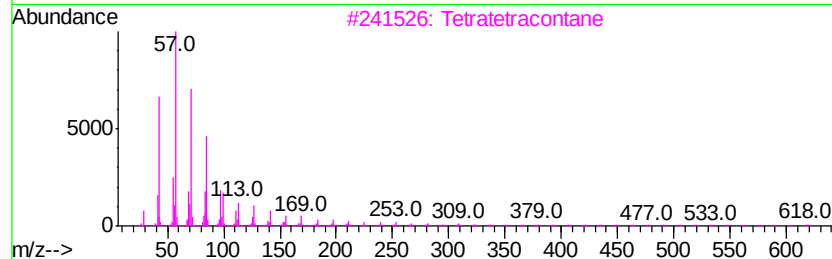
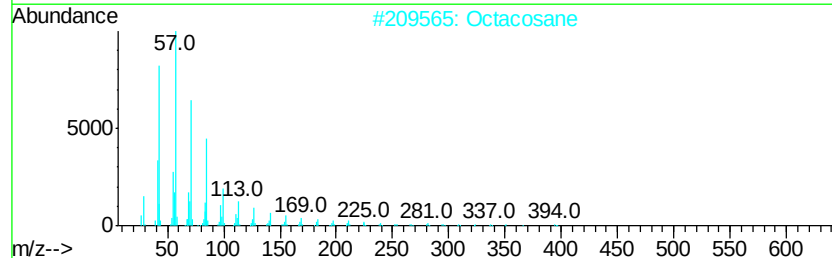
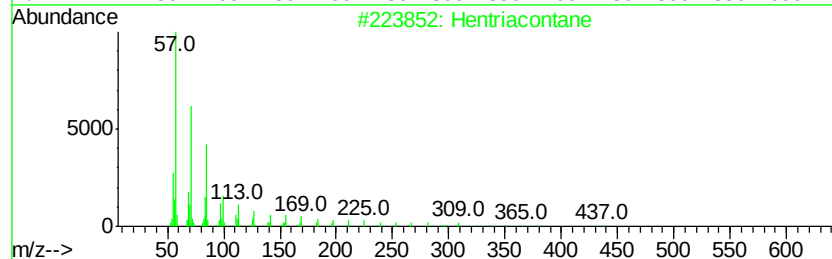
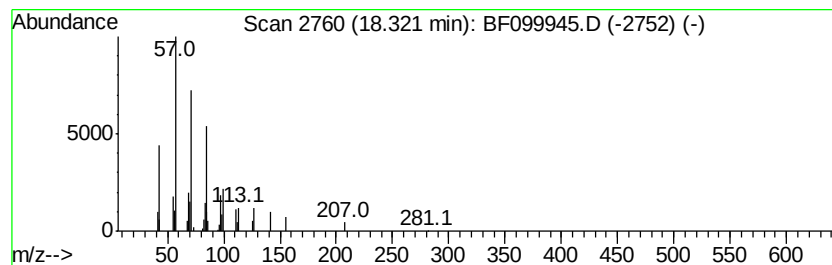
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hentriacontane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.32	4.14 ng	105496	Perylene-d12	15.42

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hentriacontane	437	C31H64	000630-04-6	91
2			Octacosane	394	C28H58	000630-02-4	91
3			Tetratetracontane	619	C44H90	007098-22-8	90
4			2-methyltetracosane	352	C25H52	1000376-72-6	90
5			2-methyloctacosane	408	C29H60	1000376-72-8	90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF102817\
Data File : BF099945.D
Acq On : 29 Oct 2017 6:04
Operator : SJ/JU
Sample : I6059-04
Misc :
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Instrument :
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ClientSampleId :
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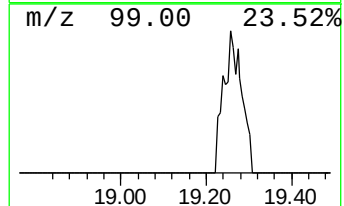
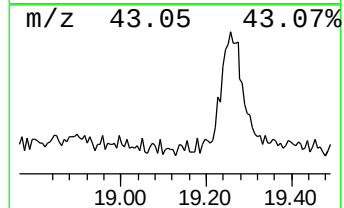
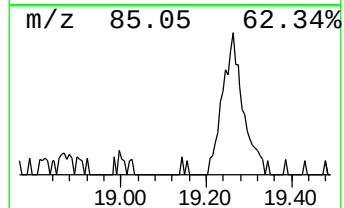
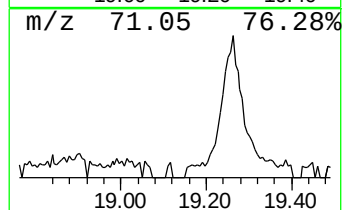
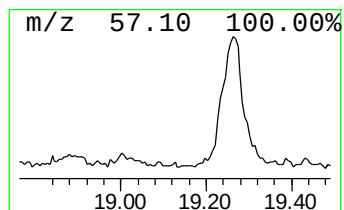
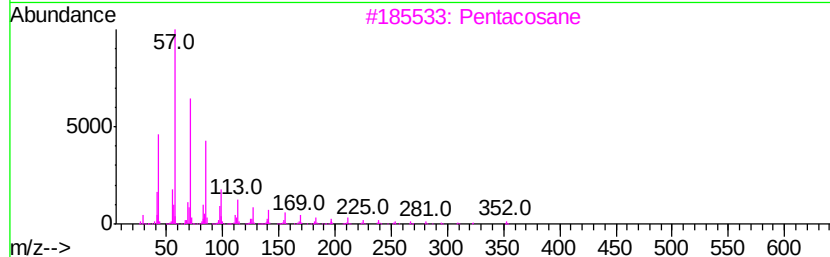
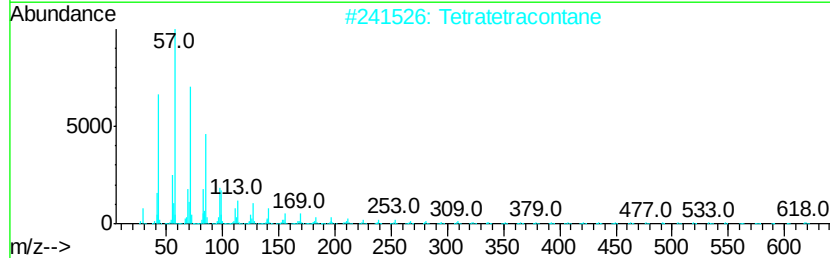
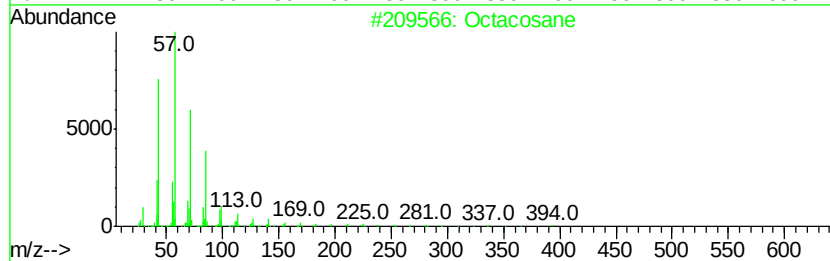
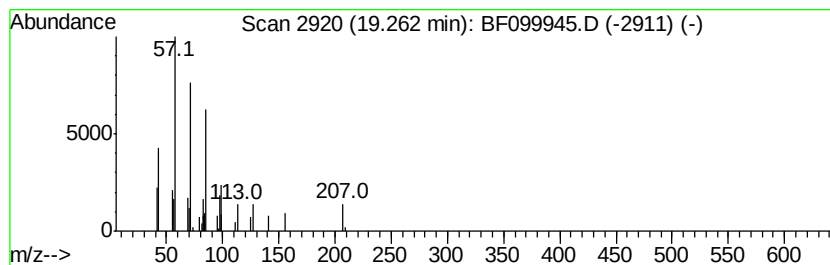
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Octacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.26	3.02 ng	77091	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	90
2		Tetratetracontane	619	C44H90	007098-22-8	90
3		Pentacosane	352	C25H52	000629-99-2	90
4		Hentriacontane	437	C31H64	000630-04-6	90
5		triacontane	422	C30H62	000638-68-6	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102817\
 Data File : BF099945.D
 Acq On : 29 Oct 2017 6:04
 Operator : SJ/JU
 Sample : I6059-04
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.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	5.2 ng		159965	1	6.79	614833	20.0
unknown6.56	6.56	49.4 ng		1517820	1	6.79	614833	20.0
2-Hydroxy-iso-but...	8.63	2.9 ng		119678	2	8.07	822784	20.0
Benzophenone	10.53	9.8 ng		402041	3	9.82	822477	20.0
Hexadecanoic acid...	12.70	4.9 ng		127862	5	13.96	524032	20.0
Butyl myristate	13.40	3.3 ng		85042	5	13.96	524032	20.0
Nonadecyl pentafl...	13.77	3.0 ng		77627	5	13.96	524032	20.0
Heneicosane	15.80	5.2 ng		131746	6	15.42	510150	20.0
Hexacosane	16.85	6.0 ng		152708	6	15.42	510150	20.0
Heptacosane	17.52	6.1 ng		156038	6	15.42	510150	20.0
Hentriacontane	18.32	4.1 ng		105496	6	15.42	510150	20.0
Octacosane	19.26	3.0 ng		77091	6	15.42	510150	20.0