Package: afmToolkit (via r-universe)

September 2, 2024

Title Functions for Atomic Force Microscope Force-Distance Curves Analysis

Version 0.0.1

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Description Set of functions for analyzing Atomic Force Microscope (AFM) force-distance curves. It allows to obtain the contact and unbinding points, perform the baseline correction, estimate the Young's modulus, fit up to two exponential decay function to a stress-relaxation / creep experiment, obtain adhesion energies. These operations can be done either over a single F-d curve or over a set of F-d curves in batch mode.

Depends R (>= 3.2.2), ggplot2

Imports stats, utils, minpack.lm, grid, gridExtra, scales, dplyr, DBI, assertthat, tibble

License GPL

LazyData true

RoxygenNote 5.0.1

NeedsCompilation no

Repository CRAN

Date/Publication 2017-04-03 15:30:48 UTC

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`afmAdhesionEnergy` *Adhesion Energy*

Description

Finds the adhesion and the full detach energies from the retract segment of the AFM F-d curve.

Usage

```
afmAdhesionEnergy(afmdata, width = 1, lagdiff = width, mul, mdj = NULL)
```

Arguments

`afmdata`	An afmdata or afmexperiment class variables. Baseline correction should have been done already.
`width`	Width of the window for the local regression (in vector position units)
`lagdiff`	Lag for estimating the differences in Delta (or slopes) signal. By default it takes the same value as the window with.
`mul`	Multiplier for the calculating the threshold inthe estimation of jumps and peaks in the Delta signal
`mdj`	Minimum distance between jumps. If none is given then it will be set equal to `width`

Value

An afmdata class variable which will consist on the original input afmdata variable plus a new list named AdhEner with the following fields:

Points Array containing the indices of the retract segment where the adhesion begins, the unbinding event takes place and the adhesion ends.

Energies Data frame with three columns: E1adh, E2adh and Etotal, being the first one the energy from the begining of the adhesion until the unbinding event, then second one the energy from the unbinding event until the full detachment of the tip, and the third one, the sum of them.

Examples

```
path <- path.package("afmToolkit")
data <- afmReadJPK("force-save-JPK-3h.txt.gz", path = path)
data <- afmContactPoint(data, width = 20, mul1 = 1, mul2 = 10)
data <- afmDetachPoint(data, width = 20, mul1 = 2, mul2 = 30)
data <- afmBaselineCorrection(data)
data <- afmAdhesionEnergy(data, width = 20, mul = 10)
str(data$AdhEner)
```

---

afmBaselineCorrection *Performs a baseline correction to an AFM F-z curve*

Description

This function performs the baseline correction to an AFM F-z curve within an afmdata structure.

It substracts a best fit line to the cuve: for the approach and contact segments, it fits a line to the approach curve points where for which |z|>ZPointApp and for the retract segment, it fits a line to the retract curve where |z|> ZpointRet.

If no ZPointApp is given and the contact point has been already estimated (via afmContactPoint() function), then it is found as

$$ZPointApp = 0.7ContactPoint + 0.3max(Z)$$

Usage

```
afmBaselineCorrection(afmdata, ZPointApp = NULL, ZPointRet = NULL,
fitpause = c("approach","retract","none"), vsTime = FALSE)
```

Arguments

afmdata	An afmdata structure.
ZPointApp	Point in the approach segment of the curve that defines the approach baseline
ZPointRet	Point in the retract segment of the curves that defines the retract baseline

fitpause	Behaviour for the baseline correction at the pause segment: if "approach" (default), the pause segment is correted using the best line fit done on the approach segment, if "retract" the best line fit of the retract segment is used, if "none", no baseline correction is done on the pause segment.
vsTime	Logical. If TRUE then the baseline correction is performed following the Force vs time approach described by S. Moreno-Flores (*Moreno Flores (2016)*).

Value

afmdata An afmdata structure identical to the one in the input, but with an additional ForceCorrected column in the data dataframe of the afmdata structure.

References

Moreno Flores (2016). Baseline correction of AFM force curves in the force-time representation. Microscopy Research and Technique, 79, (11), pp. 1045-1049.

Examples

```
AFMcurve <- afmReadJPK("force-save-JPK-2h.txt.gz", path = path.package("afmToolkit"))
ZPointApp <- 6.43e-6
ZPointRet <- 6.45e-6
AFMcurve <- afmBaselineCorrection(AFMcurve,ZPointApp = ZPointApp,ZPointRet = ZPointRet)
plot(AFMcurve)

# Without providing ZPointApp
AFMcurve <- afmReadJPK("force-save-JPK-3h.txt.gz", path = path.package("afmToolkit"))
AFMcurve <- afmContactPoint(AFMcurve,width = 10,mul1 = 1,mul2 = 20,
                            loessSmooth = FALSE)
AFMcurve <- afmBaselineCorrection(AFMcurve)
plot(AFMcurve)
```

---

afmContactPoint *Contact point*

Description

Find the contact point in for the Force-Distance curve following the local regression and two thresholds methods described in Microscopy Research and Technique 2013 (see reference).

Usage

```
afmContactPoint(afmdata, width = 1, mul1, mul2, lagdiff = width, Delta = TRUE,
  loessSmooth = FALSE)
```

Arguments

afmdata	A Force-Distance curve with the afmdata structure. It should be a list with at least the 'data' field with a data frame of at least 4 columns.
width	Width of the window for the local regression (in vector position units)
mul1	First multiplier for the first alarm threshold
mul2	Second multiplier for the second alarm threshold
lagdiff	Lag for estimating the differences in Delta (or slopes) signal. By default it takes the same value as the window with.
Delta	Logical. If TRUE, then the statistic for determining the contact point is the differences between two consecutive values of the slope of the local regression line. If FALSE then the slope itself is used.
loessSmooth	Logical If TRUE, a loess smoothing (via loess.smooth()) is done prior to the determination of the contact point. The span of the smoothing is 0.05 (5 approach segment.

Value

An afmdata class variable which will consist on the original input afmdata variable plus a new list named CP with the following fields:

CP The contact point value.

iCP The position in the array for the contact point value.

delta The delta signal.

noise The noise of the delta signal

References

Benitez R., Moreno-Flores S., Bolos V. J. and Toca-Herrera J.L. (2013). "A new automatic contact point detection algorithm for AFM force curves". Microscopy research and technique, **76** (8), pp. 870-876.

See Also

afmDetachPoint

Examples

```
path <- path.package("afmToolkit")
data <- afmReadJPK("force-save-JPK-3h.txt.gz", path = path)
width <- 20
mul1 <- 1
mul2 <- 10
data <- afmContactPoint(data, width = width, mul1 = mul1, mul2 = mul2)
## Not run:
plot(data, segment = "approach") + geom_vline(xintercept = data$CP$CP, lty = 2)

## End(Not run)
```

afmdata *AFM data*

Description

This function creates an afmdata structure, which is as list with at least one field called data which is a data frame with a valid AFM data, that is, at least 3 variables called "Z", "Force", and "Segment".

Usage

```
afmdata(data, dstr = "Z", Fstr = "Force", Segstr = "Segment", tstr = "Time",
 params = list(SpringConstant = numeric(), curvename = NULL ))
```

Arguments

data	A data frame consisting in 3 or 4 columns. A minimum of "Z" (or "distance"), "Force" and "Segment". Optionally a fourth column with "Time" could be added.
dstr	Character string with the posible names for the distance variable.
Fstr	Character string with the posible names for the force variable.
Segstr	Character string with the posible names for the Segment variable.
tstr	Character string with the posible names for the time variable.
params	A list that may contain parameters describing the F-d curve. At least will contain the SpringConstant and the curvename, being the former the cantilever spring constant and the latter a F-d curve ID. Function afmReadJPK will try to obtain the spring constant from the file header and the curvename from the data file name.

Value

An object of class afmdata

See Also

afmexperiment

Examples

```
#Making some artifical data following a L-J 12-6 potential
n <- 1000
z <- seq(from = 9e-3, to = 1e-1, length.out = n )
u0 <- 1e-5
z0 <- 1e-2
Force <- -u0*(12*z0^6/z^7-12*z0^12/z^13)
Segment <- rep("approach",n)
AFMcurve <- afmdata(data.frame(Z = z, Force = Force, Segment  = Segment))
plot(AFMcurve)
```

afmDetachPoint *Detach point*

Description

Find the detach point (or unbinding point) for the Force-Distance curve following the local regression and two thresholds methods described in Microscopy Research and Technique 2013 (see reference).

The procedure is similar to the one used by the afmContactPoint() function for obtaining the contact point.

Usage

```
afmDetachPoint(afmdata,width=1,mul1,mul2, lagdiff = width, Delta=TRUE,
  loessSmooth = FALSE)
```

Arguments

afmdata	A Force-Distance curve with the afmdata structure. It should be a list with at least the 'data' field with a data frame of at least 4 columns.
width	Width of the window for the local regression (in vector position units)
mul1	First multiplier for the first alarm threshold
mul2	Second multiplier for the second alarm threshold
lagdiff	Lag for estimating the differences in Delta (or slopes) signal. By default it takes the same value as the window with.
Delta	Logical. If TRUE, then the statistic for determining the contact point is the differences between two consecutive values of the slope of the local regression line. If FALSE then the slope itself is used.
loessSmooth	Logical If TRUE, a loess smoothing (via loess.smooth()) is done prior to the determination of the contact point. The span of the smoothing is 0.05 (5 approach segment.

Value

An afmdata class variable which will consist on the original input afmdata variable plus a new list named DP with the following fields:

DP The detach point value.

iDP The position in the array for the detach point value.

delta The delta signal.

noise The noise of the delta signal

References

Benitez R., Moreno-Flores S., Bolos V. J. and Toca-Herrera J.L. (2013). "A new automatic contact point detection algorithm for AFM force curves". Microscopy research and technique, **76** (8), pp. 870-876.

See Also

afmContactPoint

Examples

```
data <- afmReadJPK("force-save-JPK-3h.txt.gz", path = path.package("afmToolkit"))
width <- 10
mul1 <- 2
mul2 <- 40
data <- afmDetachPoint(data, width = width, mul1 = mul1, mul2 = mul2)
## Not run:
plot(data, segment = "retract") + geom_vline(xintercept = data$DP$DP, lty = 2)

## End(Not run)
```

---

afmExpDecay *Exponential decay fit*

Description

Fits a viscoelastic exponential decay in a Force-Relaxation or Creep experiments as described in Nanotechnology 2010 (see references).

Usage

```
afmExpDecay(afmdata, nexp = 2, tmax = NULL, type = c("CH","CF"), plt = TRUE,
  ...)
```

Arguments

afmdata	An object of afmdata class with a **pause** segment and a **Time** column in the data dataframe.
nexp	Number of expontials in the Prony series to be fitted. Currently only one or two exponentials are supported. Default is 2.
tmax	Maximum time considered in the relaxation curve. It defaults to Inf, meaning that the whole pause segment is considered.
type	Type of the experiment. Can be either "CH" (Constant Height) for a force-relaxation experiment or "CF" (Constant Force) for a creep experiment. Default is type = "CH".
plt	Logical. If TRUE (default) then a plot of the pause segment with the overlay of the fit is shown.

`...` Options passed to the `nlsM()` function from the `minpack.lm` package. At least should contain the starting values (`start = list(...)`) for the Levenberg-Mardquart nonlinear least square method.

Value

An `afmdata` class variable which will consist on the original input `afmdata` variable plus a new list named `ExpFit` with the following fields:

`expdecayModel`: A nls object returned from `nlsM()` function.

`expdecayFit`: The values predicted by the fit, returned from the `predict()` function.

References

Susana Moreno-Flores, Rafael Benitez, Maria dM Vivanco and Jose Luis Toca-Herrera (2010). "Stress relaxation and creep on living cells with the atomic force microscope: a means to calculate elastic moduli and viscosities of cell components". Nanotechnology, **21** (44), pp. 445101.

Examples

```
data <- afmReadJPK("force-save-JPK-3h.txt.gz", path = path.package("afmToolkit"))
width <- 20
mul1 <- 1
mul2 <- 10
data <- afmContactPoint(data, width = width, mul1 = mul1, mul2 = mul2)
data <- afmDetachPoint(data, width = width, mul1 = mul1, mul2 = mul2)
data <- afmBaselineCorrection(data)
data <- afmExpDecay(data, nexp = 2, type = "CH")
```

---

afmexperiment *AFM experiment*

Description

This function creates an `afmexperiment` structure, which is as list (or an array) of elements of `afmdata` class.

Usage

```
afmexperiment(data, ID=NULL)
```

Arguments

`data`	A variable of `afmdata` class, or a list of elements of `afmdata` class.
`ID`	Character string with the identifier of the `data` variable or a string array in case `data` is a list of `afmdata` variables.

Value

An object of class afmexp.

See Also

afmdata

Examples

```
dataFolder <- paste(path.package("afmToolkit"), "afmexperiment",sep = "/")
dataFiles <- list.files(dataFolder, pattern = "force", full.names = FALSE)
data <- lapply(dataFiles, afmReadJPK, path = dataFolder)
names(data) <- dataFiles
data <- afmexperiment(data)
plot(data[[1]])
```

---

afmExtract *Extract computed parameters from an afmexperiment*

Description

Extracts some parameters from an afmexperiment for an easy further analysis.

Usage

```
afmExtract(afmexperiment, params = list("YM", "AE", "ED"), opt.param = NULL)
```

Arguments

afmexperiment	Data of afmexperiment class.
params	List of parameters to extract from the data.
opt.param	Optional parameter or factor in the params field of the afmdata list to add to the data extraction.

Value

A data frame with the name of the curve and the corresponding values of the parameters extacted.

Examples

```
## Not run:
require(dplyr) # Not really necessary

# Load the data
data(batchExperiment)

# Process the afmexperiment
data <- afmContactPoint(batchExperiment, width = 50, mul1 = 1, mul2 = 10)
```

```
data <- afmDetachPoint(data, width = 50, mul1 = 1, mul2 = 10)
data <- afmBaselineCorrection(data)
data <- afmZeroPointSlope(data)
data <- afmIndentation(data)
data <- afmYoungModulus(data, thickness = 2e-7, params = list(alpha = 22))
data <- afmExpDecay(data, plt = FALSE)
data <- afmAdhesionEnergy(data, mul = 7)

# Extract the values of the parameters obtained in the analysis
afmExpParams <- afmExtract(data, opt.param = "type")

# Plotting the Young's Modulus
afmExpParams[[1]] %>% ggplot(aes(x = type, y = YM)) + geom_boxplot()
ylab("Young's Modulus  (Pa)")

## End(Not run)
```

`afmIndentation` *afmIndentation*

Description

This function computes the deformation of the sample from the calibrated Force-Distance curve, by substracting Z to the Zero Force Point calculated with afmZeroPointSlope function.

Usage

```
afmIndentation(afmdata)
```

Arguments

`afmdata`	An `afmdata` object. It should be a valid afmdata object upon which the Contact Point, the baseline correction and the Zero Force Point must have been calculated first (using functions `afmContactPoint()`, `afmBaselineCorrection()`) and `afmZeroPointSlope()`

Value

Returns a list with one field:

`afmdata`: An afmdata class in which a `Indentation` column is added in the `data` field.

Examples

```
data <- afmReadJPK("force-save-JPK-3h.txt.gz", path = path.package("afmToolkit"))
data <- afmContactPoint(data, width = 20, mul1 = 1, mul2 = 20)
data <- afmDetachPoint(data, width = 40, mul1 = 3, mul2 = 40)
data <- afmBaselineCorrection(data)
data <- afmZeroPointSlope(data, segment = "approach")
data <- afmIndentation(data)
head(data$data)
```

afmReadJPK *Read Nanowizard JPK ascii file*

Description

Read an ascii JPK file.

Reads an ascii JPK file with one to three headers.

Usage

```
afmReadJPK(filename, path = "", FColStr = "Vertical",
ZColStr = "Height (measured & smoothed)", tColStr = "Segment Time")
```

Arguments

filename	String with the name of the jpk file.
path	Path to the folder where the file is.
FColStr	String with a pattern identifying the Force column.
ZColStr	String with a pattern identifying the Z column.
tColStr	String with a pattern identifying the Time column.

Value

A list containing a field 'data' which is a data frame

Examples

```
data <- afmReadJPK("force-save-JPK-3h.txt.gz",path = path.package("afmToolkit"))
str(data)
```

afmReadJPKFolder *Read all Nanowizard JPK ascii files in a folder*

Description

Read all JPK ascii files in a given folder. It searches for all files containing a given patter (".txt" by default) and uses the afmReadJPJ function.

Usage

```
afmReadJPKFolder(folder, pattern = ".txt", ...)
```

Arguments

folder	Name of the folder containing the jpk files.
pattern	Pattern that will identify the jok files (".txt" by default).
...	Other parameters passed to afmReadJPK function.

Value

An afmexperiment class data structure with all F-d curves.

Examples

```
folder <- paste(path.package("afmToolkit"), "afmexperiment",sep = "/")
data <- afmReadJPKFolder(folder = folder)
str(data)
```

---

afmReadVeeco *Read Bruke Nanoscope Veeco ascii file*

Description

Read an ascii Veeco file.

Reads an ascii Veeco file with one or two segments.

Usage

```
afmReadVeeco(filename, path = "")
```

Arguments

filename	String with the name of the jpk file.
path	Path to the folder where the file is.

Value

A list containing a field 'data' which is a data frame

Examples

```
data <- afmReadVeeco("veeco_file.txt.gz",path = path.package("afmToolkit"))
str(data)
```

afmReadVeecoFolder *Read all Bruke Nanoscope Veeco ascii files in a folder*

Description

Read all Veeco ascii files in a given folder. It searches for all files containing a given patter (".txt" by default) and uses the afmReadVeeco function.

Usage

```
afmReadVeecoFolder(folder, pattern = ".txt")
```

Arguments

folder	Name of the folder containing the Veeco files.
pattern	Pattern that will identify the Veeco files (".txt" by default).

Value

An afmexperiment class data structure with all F-d curves.

Examples

```
folder <- paste(path.package("afmToolkit"), "veecoFolder",sep = "/")
data <- afmReadVeecoFolder(folder = folder)
str(data)
```

afmYoungModulus *afmYoungModulus*

Description

This function computes the Young's Modulus of the sample from the approach curve using Hertz's contact model for a pyramidal tip.

Usage

```
afmYoungModulus(afmdata, thickness = NULL, model = "Hertz", geometry =
  c("pyramid","paraboloid"), silent = TRUE, params)
```

Arguments

afmdata	An afmdata object. It should be a valid afmdata object upon which the Contact Point, the baseline correction and the Zero Force Point and the Indentation must have been calculated first (using functions afmContactPoint(), afmBaselineCorrection(), afmZeroPointSlope(), and afmIndentation())
thickness	Thickness (in m) of the surface. The Force - Indentation fit will be done for values of the Indentation variable smaller than the thickness. If no value is given, it will be done for all values in the curve for which the Indentation is negative.
model	Contact mechanics model to be used. Currently only Hertz's pure elastic model is available.
geometry	Geometry of the tip. Currently only pyramidal (default) and paraboloid geometries are implemented.
silent	Logical value. If FALSE it prints the fit model summary (via summary.lm()). Default value is TRUE
params	A list containing different parameters of the model: e.g. nu (Poisson's ratio) or alpha (internal angle, in degrees, of the pyramidal tip) or R (tip radius, in the paraboloid geometry)

Value

An afmdata class variable which will consist on the original input afmdata variable plus a new list named YoungModulus with the following fields:

YoungModulus The Young's modulus value (in Pa).

fitYM The Force vs Indentation^2 fit as an lm object.

fitdata The subset of the data used in the fit.

Examples

```
data <- afmReadJPK("force-save-JPK-2h.txt.gz", path = path.package("afmToolkit"))
data <- afmContactPoint(data, width = 20, mul1 = 1, mul2 = 20)
data <- afmDetachPoint(data, width = 40, mul1 = 3, mul2 = 40)
data <- afmBaselineCorrection(data)
data <- afmZeroPointSlope(data, segment = "approach")
data <- afmIndentation(data)
data <- afmYoungModulus(data, thickness = 1e-8, params = list(alpha = 22),
                        silent = TRUE)
print(data$YoungModulus$YoungModulus)
```

---

afmZeroPointSlope *Zero Force Point and Slope*

Description

This function finds the point of zero force (real contact point) and the slope of the contact part of the Force-Distance curve.

Usage

```
afmZeroPointSlope(afmdata, fstar = 0, segment = c("approach", "retract"))
```

Arguments

afmdata	An afmdata object. It should be a valid afmdata object upon which the Contact Point and the baseline correction must have been calculated first (using functions afmContactPoint() and afmBaselineCorrection())
fstar	Value such that fstar * sd is to be considered as zero Force, where sd is the standard deviation of Force at the basline. It takes fstar = 0 as default value, meaning that zero force is actually zero.
segment	The segment on which everything is calculated.

Value

An afmdata class variable which will consist on the original input afmdata variable plus a new list named Slopes with the following fields: Z0Point: Point of zero force. Slope: Slope of the best fit line in the contact part of the Force-Distance curve.

Examples

```
data <- afmReadJPK("force-save-JPK-2h.txt.gz", path = path.package("afmToolkit"))
data <- afmContactPoint(data, width = 20, mul1 = 1, mul2 = 20)
data <- afmDetachPoint(data, width = 40, mul1 = 3, mul2 = 40)
data <- afmBaselineCorrection(data)
data <- afmZeroPointSlope(data, segment = "approach")
## Not run:
plot(data, segment = "approach") + geom_vline(xintercept = data$Slopes$Z0Point, lty = 2)

## End(Not run)
```

---

append.afmdata *Append to an* afmdata *list.*

Description

This function appends a list to an existing afmdata structure. It is used internally by several afm* functions when attaching the results to the input afmdata variable. This function should not be used directly unless by experienced users.

Usage

```
append.afmdata(afmdata, x, name = NULL)
```

Arguments

afmdata	The afmdata to which the new list is going to be joined.
x	A list to be appended.
name	The name of new field of the resulting afmdata object. If none is given, it is the same as x.

Value

The new list of class afmdata

---

batchExperiment *Example of an afmexperiment data class.*

Description

An afmexperiment list containing 14 afmdata Force-distance experiments. Each experiment has three segments ("approach", "pause" and "retract") and they are divided in two groups depending on the covering of the sample ("CHI" for Chitosan, and "PAH" for Polyallylamine hydrochloride).

Usage

```
batchExperiment
```

Format

An afmexperiment class consisting on a list of 14 afmdata class elements each one having the following fields:

data Data frame with the data itself with a variable number of rows (between 4692 and 6142) and 4 variables:

- **Z** Distance (in meters)
- **Force** Force (in Newtons)
- **Time** Time starting at the begining of each segment (in seconds)
- **Segment** Segment of the Force-distance curve (factor: "approach", "pause", "retract")

params List with the following fields describing the experiment:

- **SpringConstant** Cantilever spring constant (in N/m)
- **curvename** Name of the original AFM data file from which the data was obtained
- **type** Type of sample covering: "CHI" for Chitosan, and "PAH" for Polyallylamine hydrochloride

is.afmdata *Afmdata check.*

Description

Checks wether an R object is an afmdata or not.

Usage

```
is.afmdata(x)
```

Arguments

x	Any **R** object.

Value

Returns TRUE if its argument is an afmdata (that is, has "afmdata" amongst its classes) and FALSE otherwise.

is.afmexperiment *Afmexperiment check.*

Description

Checks wether an R object is an afmexperiment or not.

Usage

```
is.afmexperiment(x)
```

Arguments

x	Any **R** object.

Value

Returns TRUE if its argument is an afmdata (that is, has "afmexperiment" amongst its classes) and FALSE otherwise.

plot.afmdata *Plot an afmdata object*

Description

Plots an afmdata object.

Usage

```
## S3 method for class 'afmdata'
plot(x, y = NULL, vs = "Z", segment = "all", ...)
```

Arguments

x	An object of afmdata class.
y	Variable added for compatibility with plot.
vs	The variable for the x-axis. May take the values "Time" or "Z". It defaults to "Z", plotting thus a Force-Distance curve. If vs is set to "Time", then it plots a Force-Time curve.
segment	The segment of the curve to be plotted. If segment = "all" then all segments of the curve are plotted. Possible values are: "approach", "pause", "retract" and "all".
...	Additional parameters to be pased to the ggplot functions.

Examples

```
# Loading the data
path <- path.package("afmToolkit")
data <- afmReadJPK("force-save-JPK-3h.txt.gz", path = path)
# Standard plot (out of the box)
plot(data)
# Computing the contact and detach points
data <- afmContactPoint(data, width = 20, mul1 = 1, mul2 = 10)
data <- afmDetachPoint(data, width = 40, mul1 = 3, mul2 = 20)
# Making the baseline correction
data <- afmBaselineCorrection(data)
# Plot once the baseline correction is done
plot(data)
# Plotting only retract segment
plot(data, segment = "retract")
# Plotting the pause segment: Force vs Time
plot(data, segment = "pause", vs = "Time")
```

summary.afmdata *Summary of an* afmdata *class object.*

Description

This function sumarizes the main features of an afmdata object and, optionnaly plots all segments available with all parameters estimated.

Usage

```
## S3 method for class 'afmdata'
summary(object, plt = TRUE, ...)
```

Arguments

object	An object of afmdata class.
plt	Logical variable. If TRUE plots all available segments with all available data.
...	Additional arguments (for compatibility with summary)

Examples

```
## Not run: path <- path.package("afmToolkit")
data <- afmReadJPK("force-save-JPK-3h.txt.gz", path = path)
data <- afmContactPoint(data, width = 20, mul1 = 1, mul2 = 10)
data <- afmDetachPoint(data, width = 20, mul1 = 2, mul2 = 30)
data <- afmBaselineCorrection(data)
data <- afmAdhesionEnergy(data, width = 20, mul = 10)
data <- afmZeroPointSlope(data, segment = "approach")
data <- afmIndentation(data)
data <- afmYoungModulus(data, thickness = 1e-7, params = list(alpha = 22),
                        silent = TRUE)
data <- afmExpDecay(data, nexp = 2, type = "CH")
summary(data)
## End(Not run)
```

windowedFit *Linear fit in a running window*

Description

This is an internal function used by the afmContactPoint and afmDetachPoint functions. It computes the slopes of a linear fit to the data in a window of a given radius. This function should not be used directly unless by experienced users.

Usage

```
windowedFit(X, width)
```

Arguments

X	Least squares matrix on the form [1 z Force], according to input parameters in function lm.fit
width	Width of the window for the local regression (in vector position units)

Value

OUT A vector of length nrow(X)-2*width, containing with the slopes of the fits.

Examples

```
n <- 100
x <- seq(0,2*pi,length.out = n)
y = sin(x)+0.1*rnorm(n)
X <- matrix(c(rep(1,n),x,y),nrow = n,ncol = 3)
width <- 5
b <- windowedFit(X,width)
plot(x[(width+1):(n-width)],b,xlab = "x",ylab = "y",type = "l")
lines(x,y,col = "red")
legend("bottomleft",c("Slopes","Signal"),col = c(1,2),lty = 1)
```

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