Package ‘cellGrowth’

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Title Fitting cell population growth models
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Description This package provides functionalities for the fitting of cell population growth models on experimental OD curves.
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**bandwidthCV**

**Bandwidth cross-validation**

**Description**
Perform cross-validation to detect optimal bandwidth.

**Usage**

```r
bandwidthCV(well, fileParser = readYeastGrower, 
getWellIds = getWellIdsTecan, 
bandwidths = seq(0.5 * 3600, 10 * 3600, length.out = 30), 
nFold = 10, nWell = 100, cutoff = 0.95, 
calibration = identity, scaleY = log2)
```

**Arguments**

- **well**: well dataframe. See `wellDataFrame`.
- **fileParser**: Converts the file generated by the machine to proper R format. See `readYeastGrower` for details.
- **getWellIds**: function or vector. If function its parameter is the return value of fileParser. It should return a vector containing the well ids (e.g. A01, A02, ...). You can set the well ids vector directly. See `getWellIdsTecan`.
- **bandwidths**: vector of bandwidths to test on
- **nFold**: integer. In how many parts is the sample divided for cross-validation?
- **nWell**: integer. How many wells out of the well dataframe will be used for cross validation?
- **cutoff**: scalar between 0 and 1. See details.
- **calibration**: function or list of functions. If function, calibration is applied to all raw data. If list, the well dataframe must contain a column `machine`. Depending on that column the according function in the list is applied to the raw data. See details
- **scaleY**: function applied to the calibrated data.
Details

This function needs a few minutes time. The "optimal" bandwidth is the largest bandwidth which is in 95% (cutoff parameter) of all cases within one standard deviation of the best bandwidth. This should make the derivative of the fitted curve more robust. The raw values from the machine might not be directly optical densities (OD), which is needed to infer doubling time. Calibration functions for each machine can be provided to map raw values into OD using the calibration parameter.

Value

list with entries

- bandwidth: "optimal" bandwidth
- well: well dataframe
- bandwidths: tested bandwidths
- err2: squared error
- err2std: Standard deviation of squared error
- muStd: Standard deviation of max growth rate
- oneStdOfMini: bandwidths within one std of best

Author(s)

Julien Gagneur and Andreas Neudecker

Examples

```r
folder <- system.file("extdata", package="cellGrowth")
well <- wellDataFrame(file.path(folder, "plateLayout.txt"), file.path(folder,"machineRun.txt"))

## for a fast example, we use nWell = 1 here. Use a large number (default 100) for real life applications
bw <- bandwidthCV(well, nWell=1)
```

---

**Baranyi growth model**

Description

Baranyi growth model as defined in Kelly et al.

Usage

```r
baranyi(x, mu, l, z0, zmax)
```

Arguments

- `x`: numeric vector: time points for which log(OD) must be computed
- `mu`: numeric scalar: maximal growth rate parameter
- `l`: numeric scalar: time lag parameter
- `z0`: numeric scalar: minimal log(OD) parameter
- `zmax`: numeric scalar: maximal log(OD) parameter
Value

numeric vector: log(OD) for the time points given in x

Author(s)

Julien Gagneur

References


Examples

```r
x = 1:1000
y = baranyi(x, mu=0.01, l=200, z0=1, zmax=5)
plot(x,y)
```

Description

Fit a cell growth curve

Usage

```r
fitCellGrowth(x, z, model = "locfit",
locfit.h = 3 * 60 * 60, locfit.deg = 2,
relative.height.at.lag = 0.1)
```

Arguments

- `x`: numeric vector: time points
- `z`: numeric vector: log(OD)
- `model`: which model to fit.
- `locfit.h`: numeric: h parameter (window size) in call to `locfit`. The default value is set to three hours assuming x given in seconds. You can detect a better bandwidth by calling `bandwidthCV`
- `locfit.deg`: numeric: deg parameter (polynomials degree) in call to `locfit`
- `relative.height.at.lag`: Parameter used by `guessCellGrowthParams`

Details

For the non-parametric "locfit" model, local regression is done by a call to `locfit`. The returned maximum growth rate values the maximum value of the fitted derivative over the data points. For the parametric models "logistic", "gompertz", "rosso" and "baranyi", the function does a non-least square fit by calling `nls`. Initial parameters values are generated by `guessCellGrowthParams`. The returned maximum growth rate values the `mu` parameter of these models.
Value

Fit as returned by \texttt{locfit} for the "locfit" model and as returned by \texttt{nls} for the "logistic", "gompertz", "rosso" and "baranyi" models. The returned value also has an attribute \texttt{maxGrowthRate} valuing the inferred maximum growth rate as well as an attribute \texttt{pointOfMaxGrowthRate} valuing the datapoint at which the growth rate is maximal. Also, it has an attribute \texttt{max} valuing the inferred maximum among the time points as well as \texttt{pointOfMax} valuing the datapoint of max fitted value. It gets the additional class \texttt{cellCurveFit} assigned.

Author(s)

Julien Gagneur and Moritz Matthey

See Also

\texttt{nls, locfit, baranyi, gompertz, logistic, rosso, guessCellGrowthParams, fitCellGrowths}

Examples

\begin{verbatim}
x = 1:1000
z = gompertz(x, mu=0.01, l=200, z0=1, zmax=5) + rnorm(length(x),sd=0.1)
f = fitCellGrowth(x, z, model = "gompertz")
floc = fitCellGrowth(x, z, model = "locfit", locfit.h=500)
plot(x,z, main="simulated data\nGompertz model")
lines(x, predict(f,x), lwd=2, col="red")
lines(x, predict(floc,x), lwd=2, col="blue")
legend( "right", legend=c("gompertz fit", "locfit"), lwd=1, col=c("red","blue") )
\end{verbatim}

Description

Fit growth curves for multiple wells

Usage

\begin{verbatim}
fitCellGrowth(well, plot.folder = NULL, 
model = "locfit", xlab = "time", 
ylab = expression(log2(Absorption)), scaleX = 1, 
scaleY = log2, calibration = identity, 
fileParser = readYeastGrower, 
getWellIds = getWellIdsTecan, locfit.h = 3 * 60 * 60, 
bandwidths = seq(0.5 * 3600, 10 * 3600, length.out = 30), 
nFold = 10, nWell = 100, cutoff = 0.95, ...)
\end{verbatim}

Arguments

\begin{itemize}
  \item well \hspace{1cm} data.frame with mandatory columns directory, filename, well. See \texttt{wellDataFrame}
  \item plot.folder \hspace{1cm} see details
  \item model \hspace{1cm} model to choose for fitting growth curve
\end{itemize}
fileParser Converts the file generated by the machine to proper R format. See `readYeastGrower` for details.

xlab plot parameter

ylab plot parameter

scaleX useful if you want to get the doubling in another unit, e.g. days instead of seconds.

calibration function applied to the calibrated data.

calibration function or list of functions. If function, calibration is applied to all raw data. If list, the well dataframe must contain a column machine. Depending on that column the according function in the list is applied to the raw data. See details

getWellIds function or vector. If function its parameter is the return value of fileParser. It should return a vector containing the well ids (e.g. A01, A02, ...). You can set the well ids vector directly. See `getWellIdsTecan`.

locfit.h bandwidth parameter for local polynomial fitting. If set to "bandwidthCV" bandwidth is automatically selected through `bandwidthCV` bandwidths passed to `bandwidthCV` if `locfit.h="bandwidthCV"

nFold passed to `bandwidthCV` if `locfit.h="bandwidthCV"

nWell passed to `bandwidthCV` if `locfit.h="bandwidthCV"

cutoff passed to `bandwidthCV` if `locfit.h="bandwidthCV"

... Parameter is passed to `fitCellGrowth`

Details

Essentially a wrapper for `fitCellGrowth`. The function gets a well object and fits a growth curve on all wells. It computes the doubling frequency observed in a well and extracts the maximal growth rate (1/Minimal doubling time). The raw values from the machine might not be directly optical densities (OD), which is needed to infer doubling time. Calibration functions for each machine can be provided to map raw values into OD using the calibration parameter. If the parameter plot.folder is set, the function creates a folder within plot.folder for each file in the well object. For each well a plot is written into that folder, named well_id.png.

Value
dataframe with entries

maxGrowthRate maximal growth rate

pointOfMaxGrowthRate datapoint where growth rate is maximal

max inferred maximum among the time points

pointOfMax datapoint of the max fitted value

Author(s)

Julien Gagneur and Andreas Neudecker

See Also

`fitCellGrowth`
**getRowColumn**

*Description*

Converts well ids to row and column

*Usage*

getRowColumn(wellId)

*Arguments*

- **wellId** vector of well ids

*Value*

vector of lists containing row and column

*Author(s)*

Andreas Neudecker

*Examples*

getRowColumn(c("A01","B05"))

---

**getWellIdsTecan**

*Description*

The aliases are generated by extracting the information from parsed data of the file generated by the tecan machine. See `readYeastGrower` and `readGenios`.

*Usage*

getWellIdsTecan(data)

*Arguments*

- **data** parsed data of the file. See `readYeastGrower` and `readGenios`
Value

vector containing the aliases

Author(s)

Julien Gagneur, Andreas Neudecker

Examples

data <- readYeastGrower( system.file("extdata", "Plate1_YPFruc.txt", package="cellGrowth"))
ids <- getWellIdsTecan(data)

gompertz(x, mu, l, z0, zmax)

Arguments

x numeric vector: time points for which log(OD) must be computed
mu numeric scalar: maximal growth rate parameter
l numeric scalar: time lag parameter
z0 numeric scalar: minimal log(OD) parameter
zmax numeric scalar: maximal log(OD) parameter

Value

numeric vector: log(OD) for the time points given in x

Author(s)

Julien Gagneur

References


Examples

x = 1:1000
y = gompertz(x, mu=0.01, l=200, z0=1, zmax=5)
plot(x,y)
**guessCellGrowthParams**  
*Guess growth models parameters*

**Description**

Guess initial parameters values for growth models

**Usage**

```r
guessCellGrowthParams(x, z, relative.height.at.lag = 0.1)
```

**Arguments**

- `x` numeric vector: time points
- `z` codenumeric vector: log(OD)
- `relative.height.at.lag` numeric scalar (see Details)

**Details**

The `relative.height.at.lag` parameter should be close to the relative height of the point, where the curve reaches its maximal slope. If the fitting fails, try to set this parameter to a different value.

**Value**

A list with entries:

- `mu` numeric scalar: maximal growth rate parameter
- `l` numeric scalar: time lag parameter
- `z0` numeric scalar: minimal log(OD) parameter
- `zmax` numeric scalar: maximal log(OD) parameter

**Author(s)**

Julien Gagneur

**Examples**

```r
x <- 1:1000
z <- gompertz(x, mu=0.01, l=200, z0=1, zmax=5)+rnorm(length(x),mean=0,sd=0.25)
guess <- guessCellGrowthParams(x,z,relative.height.at.lag=0.5)
fit <- nls(z~gompertz(x,mu,l,z0,zmax),start=guess)
plot(x,z)
lines(x,predict(fit,x),lwd=2,col="red")
```
logistic 

Logistic growth model

Description

Logistic growth model as defined in Zwietering et al.

Usage

logistic(x, mu, l, z0, zmax)

Arguments

x numeric vector: time points for which log(OD) must be computed
mu numeric scalar: maximal growth rate parameter
l numeric scalar: time lag parameter
z0 numeric scalar: minimal log(OD) parameter
zmax numeric scalar: maximal log(OD) parameter

Value

numeric vector: log(OD) for the time points given in x

Author(s)

Julien Gagneur

References


Examples

x = 1:1000
y = logistic(x, mu=0.01, l=200, z0=1, zmax=5)
plot(x,y)
plot.cellGrowthFit  
*Generic plot function for datatype cellGrowthFit*

**Description**
Plot of a growth curve showing raw data and fitted curve

**Usage**

```r
plot.cellGrowthFit(x, scaleX = 1, xlab = "time", ylab = "log2(OD)", lwd = 0.5, ...)
```

**Arguments**
- `x`: growth curve object. See `fitCellGrowth`
- `scaleX`: scalar affecting the scaling of the x-axis.
- `xlab`: plot parameter
- `ylab`: plot parameter
- `lwd`: plot parameter
- `...`: optional plot parameters passed to the plot function

**Author(s)**
Andreas Neudecker

**Examples**

```r
# Parse file
dat = readYeastGrower( system.file("extdata", "Plate1_YPFruc.txt", package="cellGrowth") )

# fit
n <- names( dat$OD)[36]
fit <- fitCellGrowth(x=dat$time,z=log2(dat$OD[n]), model = "locfit")
plot(fit)
```

---

plot.well  
*Generic plot function for datatype well*

**Description**
Plots well plate as lattice `xyplot`.

**Usage**

```r
plot.well(x, file = NULL, labelColumn = NULL, calibration = identity, ...)
```

---
plotPlate

Arguments

- **x**: the well object
- **file**: which plate file to plot? If NULL (default) the first file is taken.
- **labelColumn**: column in the well object to take label for the wells from
- **calibration**: function or list of functions. If calibration is a function it is applied to all raw data. If it is a list, the well dataframe must contain a column named `machine`. Depending on that column the according function in the list is applied to the raw data.
- **...**: optional plot parameters, see details

Details

This function calls `plotPlate` for the plate `plate`. The `...` parameter is passed to the `plotPlate` function.

Author(s)

Andreas Neudecker

plotPlate  
 Plot of a well plate

Description

Plot of a well plate directly from a file using a lattice xyplot

Usage

```r
plotPlate(file, labels = NULL, 
          fileParser = readYeastGrower, 
          getWellIds = getWellIdsTecan, calibration = identity, 
          extractRowColumn = getRowColumn, cex = 0.05, 
          scaleX = 1, scaleY = log2, strip.lines = 1.05, 
          strip.cex = 0.8, xlab = "time", ylab = "log2(OD)", 
          main = basename(file), 
          scales = list(x = list(rot = 45)), ...) 
```

Arguments

- **file**: file name
- **labels**: vector of characters indicating the label of the wells
- **fileParser**: the file parser which reads the file generated by the machine
- **getWellIds**: function or vector. If getWellIds is a function its parameter is the parsed data of the file parsed by `fileParser`. It should return a vector containing the well identifiers, e.g. A01, A02, .. You can as well set the well identifiers as a vector directly
- **calibration**: calibration function applied to the raw data (before `scaleY` is applied)
- **extractRowColumn**: function which converts well identifiers into row and column names
readGenios

<table>
<thead>
<tr>
<th>Parameter</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>cex</td>
<td>plot parameter</td>
</tr>
<tr>
<td>scaleX</td>
<td>factor which scales the x-axis</td>
</tr>
<tr>
<td>scaleY</td>
<td>function how to convert the y-axis (e.g. log2)</td>
</tr>
<tr>
<td>strip.lines</td>
<td>height in lines of the labels</td>
</tr>
<tr>
<td>strip.cex</td>
<td>text-size of the labels</td>
</tr>
<tr>
<td>xlab</td>
<td>plot parameter</td>
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<tr>
<td>ylab</td>
<td>plot parameter</td>
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<tr>
<td>main</td>
<td>plot parameter</td>
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<tr>
<td>scales</td>
<td>plot parameter</td>
</tr>
<tr>
<td>...</td>
<td>optional plot parameter. See details</td>
</tr>
</tbody>
</table>

Details

All plot parameters are passed to the `xyplot` function

Author(s)

Andreas Neudecker

Examples

```r
plotPlate( system.file("extdata", "tecan_genios.txt", package="cellGrowth"), fileParser=readGenios)
```

---

readGenios  
**Read Tecan Genios data files**

Description

Read raw data file form Tecan Genios instrument

Usage

```r
readGenios(file)
```

Arguments

- `file` filename

Value

a list with entries:

- `time` a numeric vector of time points
- `OD` a data.frame vector of measured OD. The colnames are the well names.
- `read` a numeric vector of read numbers
- `temperature` a numeric vector of temperatures
- `header` a character vector: the header of the file
# Description

Read raw data file from Yeast Grower software

## Usage

```r
readYeastGrower(file)
```

## Arguments

- `file`: filename

## Value

A list with entries:

- `time`: a numeric vector of time points
- `OD`: a data.frame vector of measured OD. The colnames are the well names.
- `read`: a numeric vector of read numbers
- `temperature`: a numeric vector of temperatures
- `header`: a character vector: the header of the file

## Author(s)

Julien Gagneur

## See Also

- `readGenios`
Examples

# Get file names
# Parse file
dat = readYeastGrower( system.file("extdata", "Plate1_YPFruc.txt", package="cellGrowth") )

# fit
n <- names( dat$OD)[36]
fit <- fitCellGrowth(x=dat$time, z=log2(dat$OD[[n]]), model = "locfit")
plot(fit)

---

**Rosso**

*Rosso growth model*

**Usage**

```r
rosso(x, mu, l, z0, zmax)
```

**Arguments**

- `x`: vector: time points for which log(OD) must be computed
- `mu`: scalar: maximal growth rate parameter
- `l`: scalar: time lag parameter
- `z0`: scalar: minimal log(OD) parameter
- `zmax`: scalar: maximal log(OD) parameter

**Details**

Rosso model is $z0$ if $x \leq l$ $zmax - \log( 1 + (\exp(zmax-z0) - 1)\exp(-mu*(x-l)) )$ otherwise

**Value**

vector: log(OD) for the time points given in `x`

**Author(s)**

Julien Gagneur

**Examples**

```r
x = 1:1000
y = rosso(x, mu=0.01, l=200, z0=1, zmax=5)
plot(x,y)
```
standardWellId  Make standard names for well ids

Description
Make standard names for well in 96 well plates

Usage
standardWellId(wellId)

Arguments
wellId  vector of well ids

Details
A1 -> A01 A01 -> A01

Value
standard well name

Author(s)
Julien Gagneur

Examples
standardWellId(c("A1", "B01", "H2"))

wellDataFrame  Create a well data frame

Description
Load a plate layout file and a file specifying the machine runs

Usage
wellDataFrame(plateLayoutFile, machineRunFile)

Arguments
plateLayoutFile  a file containing the plate layout. The file must contain a column named plate and a column named well
machineRunFile  a file containing the machine runs The file must contain columns named directory, filename and plate specifying the directory and filename of the data for the corresponding run. The column use is optional. If present, only rows with use == TRUE are put into the dataframe.
wellDataFrame

Details

See the provided example files for the layout and machine run file formats.

Value

an object of class well and data.frame

Author(s)

Andreas Neudecker

Examples

plateLayout <- system.file("extdata", "plateLayout.txt", package="cellGrowth")
machineRun <- system.file("extdata", "machineRun.txt", package="cellGrowth")
well <- wellDataFrame(plateLayout,machineRun)
plot(well,plate=1)
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