

Package ‘BacArena’

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Title Modeling Framework for Cellular Communities in their Environments

Version 1.8

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Description Can be used for simulation of organisms living in communities. Each organism is represented individually and genome scale metabolic models determine the uptake and release of compounds. Biological processes such as movement, diffusion, chemotaxis and kinetics are available along with data analysis techniques.

URL <https://BacArena.github.io/>

BugReports <https://github.com/euba/BacArena/issues>

Depends R (>= 3.0.0), sybil (>= 1.3.0), ReacTran (>= 1.4.2), deSolve (>= 1.12), Matrix (>= 1.2)

Imports igraph, methods, utils, stats, graphics, ggplot2, reshape2, glpkAPI, plyr, Rcpp

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addDefaultMed	<i>Add default medium of an organism to arena.</i>
---------------	--

Description

The generic function `addDefaultMed` uses the lower bounds defined in an organism's model file to compose minimal medium.

Usage

```
addDefaultMed(object, org, unit = "mM")
```

```
## S4 method for signature 'Arena'
addDefaultMed(object, org, unit = "mM")
```

Arguments

<code>object</code>	An object of class <code>Arena</code> .
<code>org</code>	An object of class <code>Organism</code>
<code>unit</code>	A character used as chemical unit to set the amount of the substances to be added (valid values are: <code>mmol/cell</code> , <code>mmol/cm2</code> , <code>mmol/arena</code> , <code>mM</code>)

addEssentialMed	<i>Add minimal medium of an organism to arena.</i>
-----------------	--

Description

The generic function `addEssentialMed` uses flux variability analysis to determine a essential growth medium components.

Usage

```
addEssentialMed(object, org, only_return = FALSE, limit = 10)
```

```
## S4 method for signature 'Arena'
addEssentialMed(object, org, only_return = FALSE,
  limit = 10)
```

Arguments

object	An object of class Arena.
org	An object of class Organism
only_return	Set true if essential metabolites should only be returned but not added to arena
limit	A metabolite is considered as essential if its remove whould decrease biomass growth below limit (between 0,100; default 10%)

addEval	<i>Function for adding a simulation step</i>
---------	--

Description

The generic function addEval adds results of a simulation step to an Eval object.

Usage

```
addEval(object, arena, replace = F)
```

```
## S4 method for signature 'Eval'
addEval(object, arena, replace = F)
```

Arguments

object	An object of class Eval.
arena	An object of class Arena.
replace	A boolean variable indicating if the last simulation step should be replaced by the new simulation step arena.

Details

The function addEval can be used in iterations to manipulate an Arena object and store the results in an Eval object.

See Also

[Eval-class](#) and [Arena-class](#)

Examples

```

data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
eval <- simEnv(arena,5)
addEval(eval,arena)

```

addOrg

Add individuals to the environment

Description

The generic function addOrg adds individuals to the environment.

Usage

```

addOrg(object, specI, amount = 1, x = NULL, y = NULL,
       posmat = NULL, biomass = NA, n0 = NULL, n = NULL, m0 = NULL,
       m = NULL)

```

```

## S4 method for signature 'Arena'
addOrg(object, specI, amount = 1, x = NULL,
       y = NULL, posmat = NULL, biomass = NA, n0 = NULL, n = NULL,
       m0 = NULL, m = NULL)

```

Arguments

object	An object of class Arena.
specI	An object of class Organism.
amount	A numeric number giving the number of individuals to add.
x	A numeric vector giving the x positions of individuals on the grid.
y	A numeric vector giving the y positions of individuals on the grid.
posmat	A As an alternative to parameter x, y, a matrix with coordinates can be specified
biomass	A numeric vector giving the starting biomass of the individuals. (unit: fg)
n0	Start column of matrix to take free positions from (default 1)
n	End column of matrix to take free positions from (default arena@m)
m0	Start row of matrix to take free positions from (default 1)
m	End row of matrix to take free positions from (default arena@n)

Details

The arguments x and y should be in the same length as the number of organisms added (given by the argument amount).

See Also

[Arena-class](#) and [Bac-class](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms

# Alternative way: adding organisms by giving matrix with positions
arena <- Arena(n=20,m=20)
mat <- matrix(sample(c(0,1), 400, replace = TRUE), nrow = 20, ncol = 20)
bac <- Bac(Ec_core)
arena <- addOrg(arena, specI=bac, posmat = mat)
```

addSubs

Add substances to the environment

Description

The generic function addSubs adds specific substances to the environment.

Usage

```
addSubs(object, smax = 0, mediac = object@mediac, difunc = "pde",
        pde = "Diff2d", difspeed = 6.7e-06, unit = "mmol/cell",
        add = TRUE, diffmat = NULL, template = FALSE, Dgrid = NULL,
        Vgrid = NULL, addAnyway = FALSE)

## S4 method for signature 'Arena'
addSubs(object, smax = 0, mediac = object@mediac,
        difunc = "pde", pde = "Diff2d", difspeed = 6.7e-06,
        unit = "mmol/cell", add = TRUE, diffmat = NULL, template = FALSE,
        Dgrid = NULL, Vgrid = NULL, addAnyway = FALSE)
```

Arguments

object	An object of class Arena.
smax	A numeric vector indicating the maximum substance concentration per grid cell.
mediac	A character vector giving the names of substances, which should be added to the environment (the default takes all possible substances).
difunc	A character vector ("pde", "cpp" or "r") describing the function for diffusion.
pde	Choose diffusion transport reaction to be used (default is diffusion only)

difspeed	A number indicating the diffusion speed (given by number of cells per iteration).
unit	A character used as chemical unit to set the amount of the substances to be added (valid values are: mmol/cell, mmol/cm2, mmol/arena, mM)
add	A boolean variable defining whether the amount of substance should be summed or replaced
diffmat	A matrix with spatial distributed initial concentrations (if not set, a homogenous matrix using smax is created)
template	True if diffmat matrix should be used as tempalte only (will be multiplied with smax to obtain cocentrations)
Dgrid	A matrix indicating the diffusion speed in x and y direction (given by cm ² /h).
Vgrid	A number indicating the advection speed in x direction (given by cm/h).
addAnyway	If true substance will be added even if there is no connection (i.e. exchanges) with organisms

Details

If nothing but object is given, then all possible substrates are initilized with a concentration of 0. Afterwards, [changeSub](#) can be used to modify the concentrations of specific substances.

See Also

[Arena-class](#) and [changeSub](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,20,c("EX_glc(e)", "EX_o2(e)", "EX_pi(e)")) #add glucose, o2, pi
```

Arena-class

Structure of the S4 class "Arena"

Description

Structure of the S4 class Arena to represent the environment in which Organisms and Substances interact.

Slots

- orgdat A data frame collecting information about the accumulated biomass, type, phenotype, x and y position for each individual in the environment.
- specs A list of organism types and their associated parameters.
- media A list of objects of class [Substance-class](#) for each compound in the environment.
- phenotypes A list of unique phenotypes (metabolites consumed and produced), which occurred in the environment.
- mediac A character vector containing the names of all substances in the environment.
- tstep A number giving the time (in h) per iteration.
- stir A boolean variable indicating if environment should be stirred. If true, bacteria move to random positions within the environment and substances have a uniform concentration value.
- mflux A vector containing highly used metabolic reactions within the arena
- exchanges A data.frame containing last exchanges of each organism.
- shadow A vector containing shadow prices of metabolites present in the arena
- n A number giving the horizontal size of the environment.
- m A number giving the vertical size of the environment.
- Lx A number giving the horizontal grid size in cm.
- Ly A number giving the vertical grid size in cm.
- gridgeometry A list containing grid geometry parameter
- seed An integer referring to the random number seed used to be reproducible
- scale A numeric defining the scale factor used for intern unit conversion.
- models A list containing Objects of class `sybil::modelorg` which represent the genome scale metabolic models
- occupyM A matrix indicating grid cells that are obstacles
- sublb A data matrix containing positions with amounts of substance for all organism

Arena-constructor *Constructor of the S4 class [Arena-class](#)*

Description

Constructor of the S4 class [Arena-class](#)

Usage

```
Arena(Lx = NULL, Ly = NULL, n = 100, m = 100,
      seed = sample(1:10000, 1), ...)
```

Arguments

Lx	A number giving the horizontal grid size in cm.
Ly	A number giving the vertical grid size in cm.
n	A number giving the horizontal size of the environment.
m	A number giving the vertical size of the environment.
seed	An integer referring to the random number seed used to be reproducible
...	Arguments of Arena-class

Bac-class	<i>Structure of the S4 class "Bac"</i>
-----------	--

Description

Structure of the S4 class Bac inheriting from class [Organism-class](#) representing bacterial cells.

Slots

deathrate	A numeric value giving the factor by which the biomass should be reduced in every iteration if no growth is possible (default (E.coli): 0.21 pg)
minweight	A numeric value giving the growth limit at which the organism dies. (default (E.coli): 0.083 pg)
cellarea	A numeric value indicating the surface that one organism occupies (default (E.coli): 4.42 μm^2)
maxweight	A numeric value giving the maximal dry weight of single organism (default (E.coli): 1.172 pg)
chem	A character vector indicating name of substance which is the chemotaxis attractant. Empty character vector if no chemotaxis.

Bac-Constructor	<i>Constructor of the S4 class Bac-class</i>
-----------------	--

Description

Constructor of the S4 class [Bac-class](#)

Usage

```
Bac(model, chem = "", ...)
```

Arguments

model	model
chem	A character vector indicating name of substance which is the chemotaxis attractant. Empty character vector if no chemotaxis.
...	Arguments of Organism

Value

Object of class [Bac-class](#)

BacArena	<i>BacArena: An Agent-Based Modeling Framework for Cellular Communities</i>
----------	---

Description

The BacArena package provides six classes: Arena (subclass Eval), Organism (subclasses Bac, Human) and Substance. Accordingly there are three categories of important functions: Arena, Organism and Substance.

Arena functions

The Arena functions ...

Organism functions

The Organism functions ...

Substance functions

The Substance functions ...

cellgrowth	<i>Function implementing a growth model of a human cell</i>
------------	---

Description

The generic function cellgrowth implements different growth models for an object of class Human.

Usage

```
cellgrowth(object, population, j, occupyM, fbasol, tstep)
```

```
## S4 method for signature 'Human'
```

```
cellgrowth(object, population, j, occupyM, fbasol, tstep)
```

Arguments

object	An object of class <code>Human</code> .
population	An object of class <code>Arena</code> .
j	The number of the iteration of interest.
occupyM	A matrix indicating grid cells that are obstacles
fbasol	Problem object according to the constraints and then solved with <code>optimizeProb</code> .
tstep	A number giving the time intervals for each simulation step.

Details

Linear growth of organisms is implemented by adding the calculated growthrate by `optimizeLP` to the already present growth value. Exponential growth of organisms is implemented by adding the calculated growthrate multiplied with the current growth calculated by `optimizeLP` plus to the already present growth value.

Value

Boolean variable of the *j*th individual indicating if individual died.

See Also

[Human-class](#), [growLin](#) and [growExp](#)

changeDiff

Change substance concentration patterns in the environment

Description

The generic function `changeDiff` changes specific substance concentration patterns in the environment.

Usage

```
changeDiff(object, newdiffmat, mediac)
```

```
## S4 method for signature 'Arena'
changeDiff(object, newdiffmat, mediac)
```

Arguments

object	An object of class <code>Arena</code> .
newdiffmat	A matrix giving the new gradient matrix of the specific substances in the environment.
mediac	A character vector giving the names of substances, which should be added to the environment (the default takes all possible substances).

Details

This function can be used to add gradients of specific substances in the environment. The default conditions in changeSubs assumes an equal concentration in every grid cell of the environment.

See Also

[Arena-class](#) and [changeSub](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,30) #add all substances with no concentrations.
gradient <- matrix(1:200,20,20)
arena <- changeDiff(arena,gradient,c("EX_glc(e)", "EX_o2(e)", "EX_pi(e)"))
# add substances glucose, oxygen and phosphate
```

changeFobj

Function for changing the objective function of the model

Description

The generic function changeFobj changes the objective function, which is used for the linear programming in optimizeLP.

Usage

```
changeFobj(object, new_fobj, model, alg = "fba")

## S4 method for signature 'Human'
changeFobj(object, new_fobj, model, alg = "fba")
```

Arguments

object	An object of class Human.
new_fobj	A character vector giving the reaction name of the new objective function.
model	The original model structure which is converted into a problem object used for the next optimization.
alg	A character vector giving the algorithm which should be used for the optimization (default is flux balance analysis).

Details

To avoid the bias to just one particular objective function, the objective can be changed dynamically in this function.

See Also

[Human-class](#) and [optimizelP](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
human <- Human(Ec_core,deathrate=0.05,
               minweight=0.05,growtype="exponential") #initialize a bacterium
changeFobj(human, 'EX_glc(e)',Ec_core)
```

changeOrg

Change organisms in the environment

Description

The generic function `changeOrg` changes organisms in the environment.

Usage

```
changeOrg(object, neworgdat)

## S4 method for signature 'Arena'
changeOrg(object, neworgdat)
```

Arguments

<code>object</code>	An object of class <code>Arena</code> .
<code>neworgdat</code>	A data frame with new information about the accumulated biomass, type, phenotype, x and y position for each individual in the environment.

Details

The argument `neworgdat` contains the same information as the `orgdat` slot of [Arena-class](#). The `orgdat` slot of an `Arena` object can be used to create `neworgdat`.

See Also

[Arena-class](#) and [addOrg](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
neworgdat <- arena@orgdat #get the current orgdat
neworgdat <- neworgdat[-1,] #remove the first individual
arena <- changeOrg(arena,neworgdat)
```

changeSub	<i>Change substances in the environment</i>
-----------	---

Description

The generic function `changeSub` changes specific substances in the environment.

Usage

```
changeSub(object, smax, mediac, unit = "mmol/cell")

## S4 method for signature 'Arena'
changeSub(object, smax, mediac, unit = "mmol/cell")
```

Arguments

<code>object</code>	An object of class <code>Arena</code> .
<code>smax</code>	A number or vector of numbers indicating the maximum substance concentration per grid cell.
<code>mediac</code>	A character vector giving the names of substances, which should be added to the environment (the default takes all possible substances).
<code>unit</code>	A character used as chemical unit to set the amount of the substances to be added (valid values are: <code>mmol/cell</code> , <code>mmol/cm2</code> , <code>mmol/arena</code> , <code>mM</code>)

Details

If nothing but `object` is given, then all possible substrates are initialized with a concentration of 0. Afterwards, [changeSub](#) can be used to modify the concentrations of specific substances.

See Also

[Arena-class](#) and [addSubs](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena) #add all substances with no concentrations.
arena <- changeSub(arena,20,c("EX_glc(e)","EX_o2(e)","EX_pi(e)"))
#add substances glucose, oxygen and phosphate
```

checkCorr	<i>Function to show correlations of a simulated organism or substrate</i>
-----------	---

Description

The generic function checkCorr returns the correlation matrix of several objects.

Usage

```
checkCorr(object, corr = NULL, tocheck = list())
```

```
## S4 method for signature 'Eval'  
checkCorr(object, corr = NULL, tocheck = list())
```

Arguments

object	An object of class Eval.
corr	A correlation matrix (getCorrM)
tocheck	A list with substrate, reactions or organism names whose correlations should be shown

Details

Returns correlation matrix which can be used for statistical analysis

See Also

[Eval-class](#) and [getCorrM](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model  
bac <- Bac(Ec_core,deathrate=0.05,  
           minweight=0.05,growthtype="exponential") #initialize a bacterium  
arena <- Arena(n=20,m=20) #initialize the environment  
arena <- addOrg(arena,bac,amount=10) #add 10 organisms  
arena <- addSubs(arena,40) #add all possible substances  
eval <- simEnv(arena,5)  
checkCorr(eval, tocheck="o2")
```

checkPhen	<i>Function for checking phenotypes in the environment</i>
-----------	--

Description

The generic function checkPhen checks and adds the phenotypes of organisms in the environment.

Usage

```
checkPhen(object, org, cutoff = 1e-06, fbasol)
```

```
## S4 method for signature 'Arena'  
checkPhen(object, org, cutoff = 1e-06, fbasol)
```

Arguments

object	An object of class Arena.
org	An object of class Organism.
cutoff	A number giving the cutoff for values of the objective function and fluxes of exchange reactions.
fbasol	Problem object according to the constraints and then solved with optimizeProb.

Details

The phenotypes are defined by flux through exchange reactions, which indicate potential differential substrate usages. Uptake of substances are indicated by a negative and production of substances by a positive number.

Value

Returns a number indicating the number of the phenotype in the phenotype list.

See Also

[Arena-class](#) and [getPhenotype](#)

checkPhen_par *Function for checking phenotypes in the environment*

Description

The generic function checkPhen_par checks and adds the phenotypes of organisms in the environment.

Usage

```
checkPhen_par(object, org, cutoff = 1e-06, fbasol)
```

```
## S4 method for signature 'Arena'
checkPhen_par(object, org, cutoff = 1e-06, fbasol)
```

Arguments

object	An object of class Arena.
org	An object of class Organism.
cutoff	A number giving the cutoff for values of the objective function and fluxes of exchange reactions.
fbasol	Problem object according to the constraints and then solved with optimizeProb.

chemotaxis *Function for chemotaxis of bacteria to their preferred substrate*

Description

The generic function chemotaxis implements a bacterial movement in the Moore neighbourhood to the highest substrate concentration.

Usage

```
chemotaxis(object, population, j, chemo, occupyM)
```

```
## S4 method for signature 'Bac'
chemotaxis(object, population, j, chemo, occupyM)
```

Arguments

object	An object of class Bac.
population	An object of class Arena.
j	The number of the iteration of interest.
chemo	The vector that contains the preferred substrate.
occupyM	A matrix indicating grid cells that are obstacles

Details

Bacteria move to a position in the Moore neighbourhood which has the highest concentration of the preferred substrate, which is not occupied by other individuals. The preferred substance is given by slot chem in the Bac object. If there is no free space the individuals stays in the same position. If the concentration in the Moore neighbourhood has the same concentration in every position, then random movement is implemented.

See Also

[Bac-class](#) and [emptyHood](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05, chem = "EX_o2(e)",
          minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
chemotaxis(bac,arena,1, "EX_glc(e)", arena@occupyM)
```

colpal1

Color palette

Description

color dictionary of 269 maximally distinct colors from all previous colors

Usage

```
colpal1
```

Format

A vector with 269 color hex-codes

Source

<http://godsnottwherogodsnott.blogspot.ru/2013/11/kmeans-color-quantization-seeding.html>

colpal2

Color palette

Description

20 optimally distinct colors

Usage

colpal2

Format

A vector with 20 color hex-codes

Source

<http://graphicdesign.stackexchange.com/questions/3682/where-can-i-find-a-large-palette-set-of-cont>

colpal3

Color palette

Description

K. Kelly (1965): Twenty-two colors of maximum contrast. // Color Eng., 3(6), 1965

Usage

colpal3

Format

A vector with 22 color hex-codes

Source

www.iscc.org/pdf/PC54_1724_001.pdf

colpal4

Color palette

Description

26 distinct colors (Zeileis(2009): Escaping RGBland: Selecting Colors for Statistical Graphics)

Usage

colpal4

Format

A vector with 26 color hex-codes

Source

<http://eeecon.uibk.ac.at/~zeileis/papers/Zeileis+Hornik+Murrell-2009.pdf>

colpal5

Color palette

Description

64 distinct colors

Usage

colpal5

Format

A vector with 64 color hex-codes

Source

<http://graphicdesign.stackexchange.com/questions/3682/where-can-i-find-a-large-palette-set-of-cont>

colpal6	<i>Color palette</i>
---------	----------------------

Description

64 distinct colors

Usage

```
colpal6
```

Format

A vector with 64 color hex-codes

Source

<http://graphicdesign.stackexchange.com/questions/3682/where-can-i-find-a-large-palette-set-of-cont>

constrain	<i>Function for constraining the models based on metabolite concentrations</i>
-----------	--

Description

The generic function constrain changes the constraints of the model representation of an organism.

Usage

```
constrain(object, reacts, lb, dryweight, tstep, scale, j, cutoff = 1e-06)
```

```
## S4 method for signature 'Organism'
constrain(object, reacts, lb, dryweight, tstep, scale,
          j, cutoff = 1e-06)
```

Arguments

object	An object of class Organisms.
reacts	A character vector giving the names of reactions which should be constrained.
lb	A numeric vector giving the constraint values of lower bounds (e.g. available metabolite concentrations)
dryweight	A number giving the current dryweight of the organism.
tstep	A number giving the time intervals for each simulation step.

scale	A numeric defining the scaling (units for linear programming has to be in certain range)
j	debugging index to track cell
cutoff	value used to define numeric accuracy while interpreting optimization results

Details

The constraints are calculated according to the flux definition as mmol/(gDW*hr) with the parameters dryweight and tstep.

Value

Returns the lower bounds, which carry the constraints and names of relevant reactions.

See Also

[Organism-class](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
org <- Organism(Ec_core,deathrate=0.05,
               minweight=0.05,growtype="exponential") #initialize an organism
lobnds <- constrain(org,org@medium,org@lbnd[org@medium],1,1,1,1,1)
```

consume

Function to account for the consumption and production of substances

Description

The generic function consume implements the consumption and production of substances based on the flux of exchange reactions of organisms

Usage

```
consume(object, sublb, cutoff = 1e-06, bacnum, fbasol)

## S4 method for signature 'Organism'
consume(object, sublb, cutoff = 1e-06, bacnum,
        fbasol)
```

Arguments

object	An object of class Organisms.
sublb	A vector containing the substance concentrations in the current position of the individual of interest.
cutoff	A number giving the cutoff value by which value of objective function is considered greater than 0.
bacnum	Integer indicating the number of bacteria individuals per gridcell
fbasol	Problem object according to the constraints and then solved with optimizeProb.

Details

The consumption is implemented by adding the flux of the exchange reactions to the current substance concentrations.

Value

Returns the updated vector containing the substance concentrations in the current position of the individual of interest.

See Also

[Organism-class](#)

Examples

NULL

createGradient	<i>Change substance concentration patterns in the environment according to a gradient</i>
----------------	---

Description

The generic function createGradient changes specific substance concentration patterns in the environment.

Usage

```
createGradient(object, mediac, position, smax, steep, add = FALSE,
              unit = "mmol/cell")
```

```
## S4 method for signature 'Arena'
createGradient(object, mediac, position, smax, steep,
              add = FALSE, unit = "mmol/cell")
```


Arguments

object	An object of class Arena.
mediac	A character vector giving the names of substances, which should be added to the environment (the default takes all possible substances).
position	A character vector giving the position (top, bottom, right and left) of the gradient.
smax	A number giving the maximum concentration of the substance.
steep	A number between 0 and 1 giving the steepness of the gradient (concentration relative to the arena size).
add	A boolean variable defining whether the amount of substance should be summed or replaced
unit	A character used as chemical unit to set the amount of the substances to be added (valid values are: mmol/cell, mmol/cm2, mmol/arena, mM)

Details

This function can be used to add gradients of specific substances in the environment.

See Also

[Arena-class](#) and [changeSub](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,30) #add all substances with no concentrations.
arena <- createGradient(arena,smax=50,mediac=c("EX_glc(e)","EX_o2(e)","EX_pi(e)"),
                       position='top',steep=0.5, add=FALSE)
```

dat2mat

Function for transforming the organism data frame to a presence/absence matrix of organisms

Description

The generic function dat2mat simulates the event of mixing all substrates and organisms in the environment.

Usage

```
dat2mat(object)
```

```
## S4 method for signature 'Arena'
dat2mat(object)
```

Arguments

object An object of class Arena.

Value

Returns the presence/absence matrix of organisms on the grid based on the orgdat slot of the Arena class.

See Also

[Arena-class](#) and [getSublb](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
occmat <- dat2mat(arena)
image(occmat)
```

diffuse

Function for diffusion

Description

The generic function `diffuse` computes the media distribution via diffusion

Usage

```
diffuse(object, lrw, sublb, verbose = TRUE)

## S4 method for signature 'Arena'
diffuse(object, lrw, sublb, verbose = TRUE)
```

Arguments

object An object of class Arena.

lrw A numeric value needed by solver to estimate array size (by default lrw is estimated in the `simEnv()` by the function `estimate_lrw()`)

sublb A matrix with the substrate concentration for every individual in the environment based on their x and y position.

verbose Set to false if no status messages should be printed.

diffusePDE *Function for diffusion of the Substance matrix*

Description

The generic function `diffusePDE` implements the diffusion by the solving diffusion equation.

Usage

```
diffusePDE(object, init_mat, gridgeometry, lrw = NULL, tstep)
```

```
## S4 method for signature 'Substance'
diffusePDE(object, init_mat, gridgeometry,
  lrw = NULL, tstep)
```

Arguments

<code>object</code>	An object of class <code>Substance</code> .
<code>init_mat</code>	A matrix with values to be used by the diffusion.
<code>gridgeometry</code>	A list specifying the geometry of the Arena
<code>lrw</code>	A numeric value needed by solver to estimate array size (by default <code>lrw</code> is estimated in <code>simEnv()</code> by the function <code>estimate_lrw()</code>)
<code>tstep</code>	A numeric value giving the time step of integration

Details

Partial differential equation is solved to model 2d diffusion process in the arena.

See Also

[Substance-class](#) and [diffuseR](#)

diffuseR *Function for naive diffusion (neighbourhood) of the Substance matrix*

Description

The generic function `diffuseR` implements the diffusion in the Moore neighbourhood in R.

Usage

```
diffuseR(object)
```

```
## S4 method for signature 'Substance'
diffuseR(object)
```

Arguments

object An object of class Substance.

Details

The diffusion is implemented by iterating through each cell in the grid and taking the cell with the lowest concentration in the Moore neighbourhood to update the concentration of both by their mean.

See Also

[Substance-class](#) and [diffusePDE](#)

diffuse_par *Function for parallelized diffusion*

Description

The generic function `diffuse_par` computes the media distribution via diffusion in parallel

Usage

```
diffuse_par(object, lrw, cluster_size, sublb)
```

```
## S4 method for signature 'Arena'
diffuse_par(object, lrw, cluster_size, sublb)
```

Arguments

object An object of class Arena.

lrw A numeric value needed by solver to estimate array size (by default lrw is estimated in the `simEnv()` by the function `estimate_lrw()`)

cluster_size Amount of cores to be used

sublb A matrix with the substrate concentration for every individual in the environment based on their x and y position.

emptyHood	<i>Function to check if there is a free place in the Moore neighbourhood</i>
-----------	--

Description

The generic function emptyHood gives a free space which is present in the Moore neighbourhood of an individual of interest.

Usage

```
emptyHood(object, pos, n, m, x, y, occupyM)
```

```
## S4 method for signature 'Organism'  
emptyHood(object, pos, n, m, x, y, occupyM)
```

Arguments

object	An object of class Organisms.
pos	A dataframe with all occupied x and y positions
n	A number giving the horizontal size of the environment.
m	A number giving the vertical size of the environment.
x	A number giving the x position of the individual of interest in its environment.
y	A number giving the y position of the individual of interest in its environment.
occupyM	A matrix indicating grid cells that are obstacles

Value

Returns the free position in the Moore neighbourhood, which is not occupied by other individuals. If there is no free space NULL is returned.

See Also

[Organism-class](#)

Examples

```
NULL
```

Eval-class	<i>Structure of the S4 class "Eval"</i>
------------	---

Description

Structure of the S4 class Eval inheriting from class [Arena-class](#) for the analysis of simulations.

Slots

medlist A list of compressed medium concentrations (only changes of concentrations are stored) per time step.

simlist A list of the organism features per time step.

mfluxlist A list of containing highly used metabolic reactions per time step.

shadowlist A list of containing shadow prices per time step.

subchange A vector of all substrates with numbers indicating the degree of change in the overall simulation.

exchangeslist A list of containing exchanges per time step.

Eval-constructor	<i>Constructor of the S4 class Eval-class</i>
------------------	---

Description

Constructor of the S4 class [Eval-class](#)

Usage

Eval(arena)

Arguments

arena An object of class Arena.

evalArena	<i>Function for plotting spatial and temporal change of populations and/or concentrations</i>
-----------	---

Description

The generic function evalArena plots heatmaps from the simulation steps in an Eval object.

Usage

```
evalArena(object, plot_items = "Population", phencol = F,
          retdata = F, time = (seq_along(object@simlist) - 1),
          show_legend = TRUE, legend_pos = "left")

## S4 method for signature 'Eval'
evalArena(object, plot_items = "Population",
          phencol = F, retdata = F, time = (seq_along(object@simlist) - 1),
          show_legend = TRUE, legend_pos = "left")
```

Arguments

object	An object of class Eval.
plot_items	A character vector giving the name of the items which should be plotted such as the population structure and several metabolites.
phencol	A boolean variable indicating if the phenotypes of the organisms in the environment should be integrated as different colors in the population plot.
retdata	A boolean variable indicating if the data used to generate the plots should be returned.
time	A numeric vector giving the simulation steps which should be plotted.
show_legend	A boolean variable indicating if a legend should be shown.
legend_pos	Position of the legend.

Details

If phencol is TRUE then different phenotypes of the same organism are visualized by varying colors, otherwise different organism types are represented by varying colors. The parameter retdata can be used to access the data used for the returned plots to create own custom plots.

Value

Returns several plots of the chosen plot items. Optional the data to generate the original plots can be returned.

See Also

[Eval-class](#) and [Arena-class](#)

Examples

```

data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
eval <- simEnv(arena,5)
evalArena(eval)
## Not run:
## if animation package is installed a movie of the simulation can be stored:
library(animation)
saveVideo({evalArena(eval)},video.name="Ecoli_sim.mp4")

## End(Not run)

```

extractMed	<i>Function for re-constructing a medium concentrations from simulations</i>
------------	--

Description

The generic function `extractMed` re-constructs a list of vectors of medium concentrations from a simulation step in an `Eval` object.

Usage

```

extractMed(object, time = length(object@medlist),
           mediac = object@mediac)

## S4 method for signature 'Eval'
extractMed(object, time = length(object@medlist),
           mediac = object@mediac)

```

Arguments

<code>object</code>	An object of class <code>Eval</code> .
<code>time</code>	A number giving the simulation step of interest.
<code>mediac</code>	A character vector giving the names of substances, which should be added to the environment (the default takes all possible substances).

Details

Medium concentrations in slot `medlist` of an object of class `Eval` store only the changes of concentrations in the simulation process. The function `extractMed` reconstructs the original and un-compressed version of medium concentrations.

Value

Returns a list containing concentration vectors of all medium substances.

See Also

[Eval-class](#) and [Arena-class](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
eval <- simEnv(arena,5)
med5 <- extractMed(eval,5)
```

findFeeding

Function for investigation of feeding between phenotypes

Description

The generic function findFeeding

Usage

```
findFeeding(object, dict = NULL, tcut = 5, scut = NULL,
            org_dict = NULL, legendpos = "topleft", lwd = 1)

## S4 method for signature 'Eval'
findFeeding(object, dict = NULL, tcut = 5,
            scut = NULL, org_dict = NULL, legendpos = "topleft", lwd = 1)
```

Arguments

object	An object of class Eval.
dict	List defining new substance names. List entries are interpreted as old names and the list names as the new ones.
tcut	Integer giving the minimal mutual occurrence to be considered (dismiss very seldom feedings)
scut	substance names which should be ignored
org_dict	A named list/vector with names that should replace (eg. unreadable) IDs
legendpos	A character variable declaring the position of the legend
lwd	Line thickness scale in graph

Value

Graph (igraph)

findFeeding2

Function for investigation of feeding between phenotypes

Description

The generic function findFeeding2

Usage

```
findFeeding2(object, time, mets, rm_own = T, ind_threshold = 0,
             collapse = F)
```

```
## S4 method for signature 'Eval'
findFeeding2(object, time, mets, rm_own = T,
             ind_threshold = 0, collapse = F)
```

Arguments

object	An object of class Eval.
time	A numeric vector giving the simulation steps which should be plotted.
mets	Character vector of substance names which should be considered
rm_own	A boolean flag indicating if interactions within same species should be plotted
ind_threshold	A number indicating the threshold of individuals to be considered as producers/consumers
collapse	A boolean flag indicating if all phenotypes for every species should be collapsed to either producers or consumers

Value

Graph (igraph)

findFeeding3	<i>Function for investigation of feeding between phenotypes</i>
--------------	---

Description

The generic function findFeeding3

Usage

```
findFeeding3(object, time, mets, plot = TRUE, cutoff = 1e-06)
```

```
## S4 method for signature 'Eval'
findFeeding3(object, time, mets, plot = TRUE,
  cutoff = 1e-06)
```

Arguments

object	An object of class Eval.
time	A numeric vector giving the simulation steps which should be plotted.
mets	Character vector of substance names which should be considered
plot	Should the graph also be plotted?
cutoff	Accuracy of crossfeeding interaction (minimal flux to be considered)

Value

Graph (igraph)

findFeeding3rep	<i>Function for investigation of cross feeding patterns of replicate simulations</i>
-----------------	--

Description

The generic function findFeeding3rep investigates the cross feeding patterns of replicate simulations

Usage

```
findFeeding3rep(simlist, time, mets, plot = TRUE, mfunction = "mean")
```

Arguments

simlist	A list with objects of class Eval.
time	A numeric vector giving the simulation steps which should be plotted.
mets	Character vector of substance names which should be considered
plot	Should the graph also be plotted?
mfunction	Function by which the replicate simulations should be combined e.g. "mean" or "median"

Value

Graph (igraph)

findInArena	<i>Function for searching a keyword in arena organisms and media</i>
-------------	--

Description

The generic function findInArena tries to find information (e.g. full names) about a specific keyword

Usage

```
findInArena(object, pattern, search_rea = TRUE, search_sub = TRUE)
```

```
## S4 method for signature 'Arena'
findInArena(object, pattern, search_rea = TRUE,
  search_sub = TRUE)
```

Arguments

object	An object of class Arena.
pattern	A pattern for searching
search_rea	Only search for reactions
search_sub	Only search for substances

Examples

```
data(Ec_core)
bac <- Bac(Ec_core)
arena <- Arena(n=20,m=20)
arena <- addOrg(arena,bac,amount=10)
findInArena(arena, "acetate")
```

findrBiomass	<i>Find biomass reaction in model</i>
--------------	---------------------------------------

Description

Helper function to search for biomass reaction in available reactions of a model

Usage

```
findrBiomass(model, keys = c("biom", "cpd11416"))
```

Arguments

model	Object of class <code>sybil::modelorg</code> containing the genome scale metabolic model
keys	Vector with strings which are used to find biomass reaction in model

Value

Vector with reaction ids for biomass reaction(s)

findRxnFlux	<i>Function to get all reactions fluxes that are associated with the metabolite of a given exchange reactions</i>
-------------	---

Description

The generic function `findRxnFlux` returns a matrix with the flux for each organism and the reaction that is using the metabolite of the given exchange reaction

Usage

```
findRxnFlux(object, ex, time, print_reactions = FALSE,
             drop_unused = TRUE)
```

```
## S4 method for signature 'Eval'
```

```
findRxnFlux(object, ex, time, print_reactions = FALSE,
             drop_unused = TRUE)
```

Arguments

object	An object of class <code>Eval</code> .
ex	An exchange reaction of which the metabolite should be shared for in all reactions
time	the time point of the simulation which should be considered
print_reactions	If true the detailed definition of each reactions is printed
drop_unused	If true then inactive reactions will be excluded

Details

Returns a list with the minimum and maximum substance usage for each time point.

See Also

[Eval-class](#) and [simEnv](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
eval <- simEnv(arena,5)
fluxlist <- findRxnFlux(eval, "EX_h(e)", 5)
```

flushSubs

Remove all substances in the environment

Description

The generic function flushSubs removes specific substances in the environment.

Usage

```
flushSubs(object)

## S4 method for signature 'Arena'
flushSubs(object)
```

Arguments

object An object of class Arena.

See Also

[Arena-class](#) and [addSubs](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena, smax=40) #add all substances with no concentrations.
arena <- changeSub(arena,20,c("EX_glc(e)", "EX_o2(e)", "EX_pi(e)"))
```

```
#add substances glucose, oxygen and phosphate
arena <- flushSubs(arena) #remove all created substance concentrations
```

fluxVarSim	<i>Function to compute flux variability analysis on an simulation object to get min/max of substance usage</i>
------------	--

Description

The generic function fluxVarSim returns a list with the minimum and maximum substance usage of all individuals for each simulation step.

Usage

```
fluxVarSim(object, rnd)

## S4 method for signature 'Eval'
fluxVarSim(object, rnd)
```

Arguments

object	An object of class Eval.
rnd	An integer giving the decimal place to which min/max flux should be rounded.

Details

Returns a list with the minimum and maximum substance usage for each time point.

See Also

[Eval-class](#) and [simEnv](#)

getArena	<i>Function for re-constructing an Arena object from a simulation step</i>
----------	--

Description

The generic function getArena re-constructs an Arena object from a simulation step within an Eval object.

Usage

```
getArena(object, time = (length(object@medlist) - 1))

## S4 method for signature 'Eval'
getArena(object, time = (length(object@medlist) - 1))
```

Arguments

object An object of class Eval.
 time A number giving the simulation step of interest.

Details

The function addEval can be used to manipulate an Arena object from a simulation step to modify the subsequent simulation steps.

Value

Returns an object of class Arena containing the organisms and substance conditions in simulation step time.

See Also

[Eval-class](#) and [Arena-class](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
eval <- simEnv(arena,5)
arena5 <- getArena(eval,5)
```

 getCorrM

Function to compute and return correlation matrix

Description

The generic function getCorrM returns the correlation matrix of several objects.

Usage

```
getCorrM(object, reactions = TRUE, bacs = TRUE, substrates = TRUE)

## S4 method for signature 'Eval'
getCorrM(object, reactions = TRUE, bacs = TRUE,
          substrates = TRUE)
```


Arguments

object	An object of class Eval.
reactions	A boolean indicating whether reactions should be included in correlation matrix
bacs	A boolean indicating whether bacteria should be included in correlation matrix
substrates	A boolean indicating whether substrates should be included in correlation matrix

Details

Returns correlation matrix which can be used for statistical analysis

Value

correlation matrix

See Also

[Eval-class](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
eval <- simEnv(arena,5)
getCorrM(eval)
```

getPhenoMat

Function for getting a matrix of phenotypes from the dataset

Description

The generic function getPhenoMat reconstructs a matrix with the usage of exchange reactions of the different organisms in the environment.

Usage

```
getPhenoMat(object, time = "total", sparse = F)
```

```
## S4 method for signature 'Eval'
```

```
getPhenoMat(object, time = "total", sparse = F)
```

Arguments

object	An object of class Eval.
time	An integer indicating the time step to be used (default value is character "total")
sparse	A boolean indicating whether zero entries should be removed from return matrix

Details

The phenotypes are defined by flux through exchange reactions, which indicate potential differential substrate usages.

Value

Returns a matrix with different phenotypes of the organism as rows and all possible exchange reactions as columns. A value of 1 means secretion, 2 means uptake and 0 means no usage of the substance of interest.

See Also

[Eval-class](#) and [getPhenotype](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
eval <- simEnv(arena,5)
phenmat <- getPhenoMat(eval)
```

getPhenotype

Function to extract the phenotype of an organism object

Description

The generic function getPhenotype implements an identification of organism phenotypes.

Usage

```
getPhenotype(object, cutoff = 1e-06, fbasol, par = FALSE)

## S4 method for signature 'Organism'
getPhenotype(object, cutoff = 1e-06, fbasol,
             par = FALSE)
```

Arguments

object	An object of class Organisms.
cutoff	A number giving the cutoff value by which value of objective function is considered greater than 0.
fbasol	Problem object according to the constraints and then solved with optimizeProb.
par	A boolean indicating if running in parallel mode.

Details

The phenotypes are defined by flux through exchange reactions, which indicate potential differential substrate usages. Uptake of substances is indicated by a negative and production of substances by a positive number.

Value

Returns the phenotype of the organisms where the uptake of substances is indicated by a negative and production of substances by a positive number

See Also

[Organism-class](#), [checkPhen](#) and [minePheno](#)

getSubHist	<i>Function to get timeline of a substance</i>
------------	--

Description

The generic function getSubHist returns list with amount of substance for each timestep

Usage

```
getSubHist(object, sub, unit = "fmol/cell")

## S4 method for signature 'Eval'
getSubHist(object, sub, unit = "fmol/cell")
```

Arguments

object	An object of class Eval.
sub	Vector with substances.
unit	Unit to be used

getSublb	<i>Function for calculated the substrate concentration for every organism</i>
----------	---

Description

The generic function `getSublb` calculates the substrate concentration for every individual in the environment based on their x and y position.

Usage

```
getSublb(object)

## S4 method for signature 'Arena'
getSublb(object)
```

Arguments

`object` An object of class `Arena`.

Value

Returns the substrate concentration for every individual in the environment with substrates as well as x and y positions as columns and rows for each organism.

See Also

[Arena-class](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
sublb <- getSublb(arena)
```

getVarSubs *Function to get varying substances*

Description

The generic function getVarSubs returns ordered list of substances that showed variance during simulation

Usage

```
getVarSubs(object, show_products = TRUE, show_substrates = TRUE,
           cutoff = 1e-06, size = NULL)
```

```
## S4 method for signature 'Eval'
getVarSubs(object, show_products = FALSE,
           show_substrates = FALSE, cutoff = 1e-06, size = NULL)
```

Arguments

object	An object of class Eval.
show_products	A boolean indicating if only products should be shown
show_substrates	A boolean indicating if only substrates should be shown
cutoff	Value used to define numeric accuracy while interpreting optimization results
size	Maximal number of returned substances (default: show all)

growExp *Function for letting organisms grow exponentially*

Description

The generic function growExp implements a growth model of organisms in their environment.

Usage

```
growExp(object, biomass, fbasol, tstep)
```

```
## S4 method for signature 'Organism'
growExp(object, biomass, fbasol, tstep)
```

Arguments

object	An object of class Organisms.
biomass	A number indicating the current biomass, which has to be updated.
fbasol	Problem object according to the constraints and then solved with optimizeProb.
tstep	A number giving the time intervals for each simulation step.

Details

Exponential growth of organisms is implemented by adding the calculated growthrate multiplied with the current growth calculated by `optimizeLP` plus to the already present growth value

Value

Returns the updated biomass of the organisms of interest.

See Also

[Organism-class](#) and [optimizeLP](#)

growLin	<i>Function for letting organisms grow linearly</i>
---------	---

Description

The generic function `growLin` implements a growth model of organisms in their environment.

Usage

```
growLin(object, biomass, fbasol, tstep)
```

```
## S4 method for signature 'Organism'
growLin(object, biomass, fbasol, tstep)
```

Arguments

<code>object</code>	An object of class <code>Organisms</code> .
<code>biomass</code>	A number indicating the current biomass, which has to be updated.
<code>fbasol</code>	Problem object according to the constraints and then solved with <code>optimizeProb</code> .
<code>tstep</code>	A number giving the time intervals for each simulation step.

Details

Linear growth of organisms is implemented by adding the calculated growthrate by `optimizeLP` to the already present growth value.

Value

Returns the updated biomass of the organisms of interest.

See Also

[Organism-class](#) and [optimizeLP](#)

growth	<i>Function implementing a growth model of a bacterium</i>
--------	--

Description

The generic function growth implements different growth models for an object of class Bac.

Usage

```
growth(object, population, j, occupyM, fbasol, timestep)
```

```
## S4 method for signature 'Bac'  
growth(object, population, j, occupyM, fbasol, timestep)
```

Arguments

object	An object of class Bac.
population	An object of class Arena.
j	The index of the organism of interest in orgdat.
occupyM	A matrix indicating grid cells that are obstacles
fbasol	Problem object according to the constraints and then solved with optimizeProb.
tstep	A number giving the time intervals for each simulation step.

Details

Linear growth of organisms is implemented by adding the calculated growthrate by optimizeLP to the already present growth value. Exponential growth of organisms is implemented by adding the calculated growthrate multiplied with the current growth calculated by optimizeLP plus to the already present growth value

Value

Boolean variable of the jth individual indicating if individual died.

See Also

[Bac-class](#), [growLin](#) and [growExp](#)

growth_par	<i>Function implementing a growth model of a bacterium</i>
------------	--

Description

The generic function growth_par implements different growth models for an object of class Bac.

Usage

```
growth_par(object, population, j, fbasol, tstep)
```

```
## S4 method for signature 'Bac'
growth_par(object, population, j, fbasol, tstep)
```

Arguments

object	An object of class Bac.
population	An object of class Arena.
j	The index of the organism of interest in orgdat.
fbasol	Problem object according to the constraints and then solved with optimizeProb.
tstep	A number giving the time intervals for each simulation step.

Value

A list

HeatMapFeeding	<i>Function for plotting heatmap of feeding between two selected cell models</i>
----------------	--

Description

The generic function HeatMapFeeding returns a heatmap between two selected cell models. If there is feeding, the heatmap show only the producer. The compounds are originated from the most varying substances of the first cell model called "speciesA".

Usage

```
HeatMapFeeding(object, speciesA, speciesB, var_nr)
```

```
## S4 method for signature 'Eval'
HeatMapFeeding(object, speciesA, speciesB, var_nr)
```


Arguments

object	An object of class Eval.
speciesA	The sequence number of the first cell model in names(object@specs)
speciesB	The sequence number of the second cell model in names(object@specs)
var_nr	The number of the most varying speciesA's substances following plotSpecActivity structure

Value

Heatmap (ggplot2)

Examples

```
sim <- sihumi_test
HeatMapFeeding(object = sim, speciesA = 1, speciesB = 2, var_nr = 78)
```

Human-class

Structure of the S4 class "Human"

Description

Structure of the S4 class Human inheriting from class [Organism-class](#) representing human cells.

Slots

objective A character vector representing the current reaction which should be used as an objective function for the flux balance analysis.

Human-constructor

Constructor of the S4 class [Human-class](#)

Description

Constructor of the S4 class [Human-class](#)

Usage

```
Human(model, objective = model@react_id[which(model@obj_coef != 0)],
      speed = 0, ...)
```

Arguments

model	model
objective	A character vector representing the current reaction which should be used as an objective function for the flux balance analysis.
speed	A integer vector representing the speed by which bacterium is moving (given by cell per iteration).
...	Arguments of Organism

Value

Object of class [Human-class](#)

lsd	<i>Computer standard deviation lower bound</i>
-----	--

Description

Helper function to get lower error bounds in plotting

Usage

```
lsd(y)
```

Arguments

y	Vector with numbers
---	---------------------

lysis	<i>Lysis function of organismal cells by adding biomass_compounds to the medium</i>
-------	---

Description

The generic function `lysis` implements cell lysis by the stoichiometric concentration of the biomass compounds of organisms to the concentration of substances in the environment

Usage

```
lysis(object, sublb, factor = object@minweight)
```

```
## S4 method for signature 'Organism'
```

```
lysis(object, sublb, factor = object@minweight)
```

Arguments

object	An object of class Organisms.
sublb	A vector containing the substance concentrations in the current position of the individual of interest.
factor	A number given the factor with which the biomass compound concentrations are multiplied to achieve the final concentration which is added to the environment

Details

Lysis is implemented by taking the intersect between biomass compounds and the substances in the environment and adding the normalized stoichiometric concentrations of the biomass compounds to the medium.

Value

Returns the updated vector containing the substance concentrations in the current position of the dead individual of interest.

See Also

[Organism-class](#) and [optimizeLP](#)

Examples

NULL

minePheno	<i>Function for mining/analyzing phenotypes which occurred on the arena</i>
-----------	---

Description

The generic function minePheno mines the similarity and differences of phenotypes reconstructed by getPhenoMat for each simulation step in an Eval object.

Usage

```
minePheno(object, plot_type = "pca", legend = F, time = "total")

## S4 method for signature 'Eval'
minePheno(object, plot_type = "pca", legend = F,
           time = "total")
```

Arguments

object	An object of class Eval.
plot_type	A character vector giving the plot which should be returned (either "pca" for a principle coordinate analysis or "hclust" for hierarchical clustering).
legend	Boolean variable indicating if legend should be plotted
time	An integer indicating the time step to be used (default value is character "total")

Details

The phenotypes are defined by flux through exchange reactions, which indicate potential differential substrate usages.

Value

Returns a plot for each simulation step representing the similarity of phenotypes of organisms within the environment.

See Also

[Eval-class](#) and [getPhenoMat](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
eval <- simEnv(arena,5)
minePheno(eval)
```

move

Function for random movement of organisms

Description

The generic function move implements a random movement in the Moore neighbourhood of an individual.

Usage

```
move(object, pos, n, m, j, occupyM)

## S4 method for signature 'Organism'
move(object, pos, n, m, j, occupyM)
```

Arguments

object	An object of class Organism.
pos	A dataframe with all occupied x and y positions
n	A number giving the horizontal size of the environment.
m	A number giving the vertical size of the environment.
j	The number of the iteration of interest.
occupyM	A matrix indicating grid cells that are obstacles

Details

Organisms move in a random position the Moore neighbourhood, which is not occupied by other individuals. If there is no free space the individuals stays in the same position.

See Also

[Organism-class](#), [emptyHood](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
move(bac,n=20,m=20,j=1,pos=arena@orgdat[,c('x','y')], occupyM=arena@occupyM)
```

NemptyHood	<i>Function to check if there is a free place in the Moore neighbourhood</i>
------------	--

Description

The generic function NemptyHood gives a free space which is present in the Moore neighbourhood of an individual of interest.

Usage

```
NemptyHood(object, pos, n, m, x, y, occupyM)
```

```
## S4 method for signature 'Organism'
```

```
NemptyHood(object, pos, n, m, x, y, occupyM)
```

Arguments

object	An object of class Organisms.
pos	A dataframe with all occupied x and y positions
n	A number giving the horizontal size of the environment.
m	A number giving the vertical size of the environment.
x	A number giving the x position of the individual of interest in its environment.
y	A number giving the y position of the individual of interest in its environment.
occupyM	A matrix indicating grid cells that are obstacles

Value

Returns the free position in the Moore neighbourhood, which is not occupied by other individuals. If there is no free space NULL is returned.

See Also

[Organism-class](#)

Examples

NULL

openArena	<i>Start simulation</i>
-----------	-------------------------

Description

The function openArena can be used to start a default simulation.

Usage

```
openArena()
```

Value

Returns an object of class Eval which can be used for subsequent analysis steps.

Examples

```
sim <- openArena()
evalArena(sim, time=5, phencol = TRUE,
          plot_items=c("Population", "EX_o2(e)", "EX_for(e)", "EX_glc(e)"))
```

optimizeLP	<i>Function for computing the linear programming according to the model structure</i>
------------	---

Description

The generic function optimizeLP implements a linear programming based on the problem structure and refined constraints.

Usage

```
optimizeLP(object, lpobj = object@lpobj, lb = object@lbnd,
           ub = object@ubnd, cutoff = 1e-06, j, sec_obj = "none",
           with_shadow = FALSE)
```

```
## S4 method for signature 'Organism'
optimizeLP(object, lpobj = object@lpobj,
           lb = object@lbnd, ub = object@ubnd, cutoff = 1e-06, j,
           sec_obj = "none", with_shadow = FALSE)
```

Arguments

object	An object of class Organisms.
lpobj	A linear programming object encoding the problem to solve.
lb	A numeric vector giving the constraint values of lower bounds.
ub	A numeric vector giving the constraint values of upper bounds.
cutoff	value used to define numeric accuracy while interpreting optimization results
j	debugging index to track cell
sec_obj	character giving the secondary objective for a bi-level LP if wanted. Use "mtf" for minimizing total flux, "opt_rxn" for optimizing a random reaction, "opt_ex" for optimizing a random exchange reaction, and "sumex" for optimizing the sum of all exchange fluxes.
with_shadow	True if shadow costs should be stored (default false).

Details

The parameter for sec_obj can be used to optimize a bi-level LP with a secondary objective if wanted. This can be helpful to subselect the solution space and create less alternative optimal solution. The secondary objective can be set to "mtf" to minimize the total flux, to simulate minimal enzyme usage of an organisms. If set to "opt_rxn" or "opt_ex", the secondary objective is picked as a random reaction or exchange reaction respectively everytime a fba is performed. This means that every individual of a population will select a different secondary reaction to optimize. The "sumex" option maximizes the secretion of products.

Value

Modified problem object according to the constraints and then solved with `optimizeProb`.

See Also

[Organism-class](#), [optimizeProb](#) and [sysBiolAlg](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
org <- Organism(Ec_core,deathrate=0.05,
               minweight=0.05,growtype="exponential") #initialize a organism
org@fbasol <- optimizeLP(org)
```

Organism-class

Structure of the S4 class "Organism"

Description

Structure of the S4 class `Organism` representing the organisms present in the environment.

Slots

`lbnd` A numeric vector containing the lower bounds of the model structure.

`ubnd` A numeric vector containing the upper bounds of the model structure.

`type` A character vector containing the description of the organism.

`medium` A character vector containing all exchange reactions of the organism.

`lproj` A sybil optimization object containing the linear programming problem.

`fbasol` A list with the solutions of the flux balance analysis.

`lyse` A boolean variable indicating if the organism should lyse after death.

`feat` A list containing conditional features for the object (contains at the moment only biomass components for lysis).

`deathrate` A numeric value giving the factor by which the biomass should be reduced in every iteration if no growth is possible (default (E.coli): 0.21 pg)

`minweight` A numeric value giving the growth limit at which the organism dies. (default (E.coli): 0.083 pg)

`growtype` A character vector giving the functional type for growth (linear or exponential).

`kinetics` A List containing `Km` and `v_max` values for each reactions.

`speed` A integer vector representing the speed by which bacterium is moving (given by cell per iteration).

`cellarea` A numeric value indicating the surface that one organism occupies (default (E.coli): 4.42 μm^2)

maxweight A numeric value giving the maximal dry weight of single organism (default (E.coli): 1.172 pg)
cellweight_mean A numeric giving the mean of starting biomass (default (E.coli): 0.489 pg)
cellweight_sd A numeric giving the standard derivation of starting biomass (default (E.coli): 0.132 pg)
model Object of class `sybil::modelorg` containing the genome scale metabolic model
algo Algorithm to be used during optimization (default `fba`)
rbiomass Name of biomass reactions which is used for growth model (set automatically but needs input if objective is not biomass optimization)
limit_growth If true then an upper bound on growth will be set, see `maxweight` (default: `True`).

Organism-constructor *Constructor of the S4 class* [Organism-class](#)

Description

Constructor of the S4 class [Organism-class](#)

Usage

```
Organism(model, algo = "fba", ex = "EX_", ex_comp = NA,
          csuffix = "\\[c\\]", esuffix = "\\[e\\]", lyse = FALSE,
          feat = list(), typename = NA, setExInf = TRUE, ...)
```

Arguments

<code>model</code>	Object of class <code>sybil::modelorg</code> containing the genome scale metabolic model
<code>algo</code>	A single character string giving the name of the algorithm to use. See SYBIL_SETTINGS
<code>ex</code>	Identifier for exchange reactions
<code>ex_comp</code>	Defining exchange reactions whose compounds should be added to the medium of the arena (default: <code>all</code>)
<code>csuffix</code>	suffix for intern metabolites used by lysis function.
<code>esuffix</code>	suffix for external metabolites used by lysis function.
<code>lyse</code>	A boolean variable indicating if the organism should lyse after death.
<code>feat</code>	A list containing conditional features for the object (contains at the moment only biomass components for lysis).
<code>typename</code>	A string defining the name (set to model name in default case)
<code>setExInf</code>	Enable if all lower bounds of exchange reaction which are set to zero (i.e. no uptake possible!) should be set to -infinity
<code>...</code>	Arguments of Organism-class

Value

Object of class `Organism`

plotAbundance	<i>Plot abundances of species</i>
---------------	-----------------------------------

Description

The function plotAbundance takes a list of simulations and return a boxplot with species abundances

Usage

```
plotAbundance(simlist, time = c(NULL, NULL), col = colpal3,
  return_dat = F, use_biomass = F)
```

Arguments

simlist	A list of simulations (eval objects).
time	A vector with start and end time to be considered (default: total time)
col	Vector with color that should be used
return_dat	Should plain text mean abundances be returned? (default false)
use_biomass	If enabled then biomass is used instead of cell number

plotCurves	<i>Function for plotting the overall change as curves</i>
------------	---

Description

The generic function plotCurves plots the growth curves and concentration changes of substances from simulation steps in an Eval object.

Usage

```
plotCurves(object, medplot = object@mediac, retdata = F, remove = F,
  legend = F, graph = T)
```

```
## S4 method for signature 'Eval'
plotCurves(object, medplot = object@mediac,
  retdata = F, remove = F, legend = F, graph = T)
```

Arguments

object	An object of class Eval.
medplot	A character vector giving the name of substances which should be plotted.
retdata	A boolean variable indicating if the data used to generate the plots should be returned.
remove	A boolean variable indicating if substances, which don't change in their concentration should be removed from the plot.
legend	Boolean variable indicating if legend should be plotted
graph	True if graphic should be plotted.

Details

The parameter `retdata` can be used to access the data used for the returned plots to create own custom plots.

Value

Returns two graphs in one plot: the growth curves and the curves of concentration changes. Optional the data to generate the original plots can be returned.

See Also

[Eval-class](#) and [Arena-class](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
          minweight=0.05,growthtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
eval <- simEnv(arena,5)
plotCurves(eval)
```

plotCurves2

Function for plotting the overall change as curves with maximally distinct colors

Description

The generic function `plotCurves2` plots the growth curves and concentration changes of the most changing substances from simulation steps in an Eval object using maximally distinct colors.

Usage

```
plotCurves2(object, legendpos = "topleft", ignore = c("EX_h(e)",
  "EX_pi(e)", "EX_h2o(e)"), num = 10, phencol = FALSE,
  biomcol = FALSE, dict = NULL, subs = list(), growthCurve = TRUE,
  subCurve = TRUE)

## S4 method for signature 'Eval'
plotCurves2(object, legendpos = "topright",
  ignore = c("EX_h(e)", "EX_pi(e)", "EX_h2o(e)"), num = 10,
  phencol = FALSE, biomcol = FALSE, dict = NULL, subs = list(),
  growthCurve = TRUE, subCurve = TRUE)
```

Arguments

object	An object of class Eval.
legendpos	A character variable declaring the position of the legend
ignore	A list of character variables with substance names that could be omitted in the plot
num	An integer defining the number of substrates to be plot
phencol	Boolean variable indicating whether phenotypes should be highlighted
biomcol	A boolean indicating if biomass should be included in growth curve
dict	List defining new substance names. List entries are interpreted as old names and the list names as the new ones.
subs	List of substance names. If empty, substances with highest variance will be used.
growthCurve	True if growth curve should be shown (default TRUE)
subCurve	True if substance curve should be shown (default TRUE)

Details

The parameter `retdata` can be used to access the data used for the returned plots to create own custom plots.

Value

Returns two graphs in one plot: the growth curves and the curves of concentration changes

See Also

[Eval-class](#) and [Arena-class](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
  minweight=0.05,growthtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
```

```
arena <- addSubs(arena,40) #add all possible substances
eval <- simEnv(arena,5)
plotCurves2(eval)
```

plotFluxVar	<i>Plot population flux variations</i>
-------------	--

Description

The function plotFluxVar takes a list of simulations and metabolites, returning a plot with metabolite fluxes for each species

Usage

```
plotFluxVar(simlist, metsel)
```

Arguments

simlist	A list of simulations (eval objects).
metsel	A vector with the name of exchange reactions of interest

plotFVA	<i>Function to plot population level minimum and maximum flux from alternative optimal solutions obtained by FVA</i>
---------	--

Description

The generic function plotFVA plots population level flux results obtained from function fluxVarSim

Usage

```
plotFVA(fvares, mediac)
```

Arguments

fvares	results of FVA results to plot, obtained from function fluxVarSim
mediac	List with substances.

Details

Returns ggplot objects

plotGrowthCurve	<i>Plot growth curve for several simulations</i>
-----------------	--

Description

The function `plotGrowthCurve` takes a list of simulations and plots the time course of species with standard deviation.

Usage

```
plotGrowthCurve(simlist, time = c(NULL, NULL), ret_data = FALSE,
  use_biomass = F, specs = NULL)
```

Arguments

<code>simlist</code>	A list of simulations (eval objects).
<code>time</code>	Vector with two entries defining start and end time
<code>ret_data</code>	Set true if data should be returned
<code>use_biomass</code>	If enabled then biomass is used instead of cell number
<code>specs</code>	List of species for which a growth curve should be shown (default: all)

plotInterNum	<i>Plot number of variation in number of interactions for several simulations</i>
--------------	---

Description

The function `plotInterNum` takes a list of simulations and plots the time course of the number of metabolic interactions with standard deviation.

Usage

```
plotInterNum(simlist, title = "Variation in number of interactions",
  size = 1)
```

Arguments

<code>simlist</code>	A list of simulations (eval objects).
<code>title</code>	Title of the plot
<code>size</code>	A scaling factor for plot text and line size

plotPhenCurve	<i>Plot growth curve for several simulations</i>
---------------	--

Description

The function plotPhenCurve takes a list of simulations and plots the time course of species with standard deviation.

Usage

```
plotPhenCurve(simlist, subs, phens = NULL, time = c(NULL, NULL),
  cluster = TRUE, inAll = TRUE, col = colpal3, with_gc = FALSE,
  return_dat = FALSE)
```

Arguments

simlist	A list of simulations (eval objects).
subs	A vector of substance names that are used for phenotype clustering.
phens	If phencurve is given then phens specifies the phenotypes which could be plotted again.
time	Vector with two entries defining start and end time
cluster	True if phenotypes should be clustered/condensed.
inAll	True if only phenotypes which occur in all replicates should be considered
col	Vector with color that should be used
with_gc	True if growth curve of organisms should be included
return_dat	Should data be returned? (default false)

plotPhenNum	<i>Plot number of phenotypes curve for several simulations</i>
-------------	--

Description

The function plotPhenNum takes a list of simulations and plots the time course of the number of phenotypes with standard deviation.

Usage

```
plotPhenNum(simlist, title = "Phenotype number variation", size = 1)
```

Arguments

simlist	A list of simulations (eval objects).
title	Title of the plot
size	A scaling factor for plot text and line size

plotReaActivity *Function to plot reaction activity for every species*

Description

The generic function plotReaActivity displays the usage of reactions for all species

Usage

```
plotReaActivity(simlist, reactions = list(), spec_list = NULL,
  ret_data = FALSE)
```

Arguments

simlist	An object of class Eval or a list with objects of class Eval.
reactions	List of reaction names
spec_list	List of species names to be considered (default all)
ret_data	Set true if data should be returned

Details

Returns ggplot objects

plotShadowCost *Function to plot substance shadow costs for a specie*

Description

The generic function plotShadowCost plots substances have the highest impact on further growth (shadow cost < 0)

Usage

```
plotShadowCost(object, spec_nr = 1, sub_nr = 10, cutoff = -1,
  noplot = FALSE)
```

```
## S4 method for signature 'Eval'
plotShadowCost(object, spec_nr = 1, sub_nr = 10,
  cutoff = -1, noplot = FALSE)
```


Arguments

object	An object of class Eval.
spec_nr	Number of the specie
sub_nr	Maximal number of substances to be show
cutoff	Shadow costs should be smaller than cutoff
noplot	Do not plot

Details

Returns ggplot objects

plotSpecActivity *Function to plot substance usage for every species*

Description

The generic function plotSpecActivity displays the input/output substances with the highest variance (could also be defiened manually) for all species

Usage

```
plotSpecActivity(simlist, subs = list(), var_nr = 10,  
spec_list = NULL, ret_data = FALSE, useNames = FALSE)
```

Arguments

simlist	An object of class Eval or a list with objects of class Eval.
subs	List of substance names
var_nr	Number of most varying substances to be used (if subs is not specified)
spec_list	List of species names to be considered (default all)
ret_data	Set true if data should be returned
useNames	Use substance names instead of ids

Details

Returns ggplot objects

plotSubCurve	<i>Plot substance curve for several simulations</i>
--------------	---

Description

The function `plotSubCurve` takes a list of simulations and plots the time course of substances with standard deviation.

Usage

```
plotSubCurve(simlist, mediac = NULL, time = c(NULL, NULL),
             scol = NULL, unit = "mmol", ret_data = FALSE, num_var = 10,
             useNames = FALSE)
```

Arguments

<code>simlist</code>	A list of simulations (eval objects).
<code>mediac</code>	A vector of substances (if not specified most varying substances will be taken.)
<code>time</code>	Vector with two entries defining start and end time.
<code>scol</code>	Vector with colors that should be used.
<code>unit</code>	Unit for the substances which should be used for plotting (default: mmol)
<code>ret_data</code>	Set true if data should be returned
<code>num_var</code>	Number of varying substances to be shown (if mediac is not specified)
<code>useNames</code>	Use substance names instead of ids

Value

list of three ggplot object for further formating

plotSubDist	<i>Function to overview the spatial distribution of a substance over time.</i>
-------------	--

Description

The generic function `plotSubDist` returns a plot for every time step which shows the substance concentration in the environment.

Usage

```
plotSubDist(object, sub, times = NULL)
```

```
## S4 method for signature 'Eval'
plotSubDist(object, sub, times = NULL)
```

Arguments

object	An object of class Eval.
sub	Name of a substance.
times	Time points to be considered.

Details

Returns a plot with

plotSubDist2 *Function to overview the spatial distribution of a substance over time.*

Description

The generic function plotSubDist2 returns a plot for every time step which shows the substance concentration in the environment.

Usage

```
plotSubDist2(object, sub, times = NULL)

## S4 method for signature 'Eval'
plotSubDist2(object, sub, times = NULL)
```

Arguments

object	An object of class Eval.
sub	Name of a substance.
times	Time points to be considered.

Details

Returns a plot with

plotSubUsage	<i>Function to plot usage of substances species wise</i>
--------------	--

Description

The generic function plotSubUsage displays for given substances the quantities of absorption and production for each species

Usage

```
plotSubUsage(simlist, subs = vector(), cutoff = 0.01,  
             ret_data = FALSE)
```

Arguments

simlist	An object of class Eval or a list with objects of class Eval.
subs	List of substance names
cutoff	Total values below cutoff will be dismissed
ret_data	Set true if data should be returned

Details

Returns ggplot objects

plotSubVar	<i>Plot substance variations</i>
------------	----------------------------------

Description

The function plotSubVar takes a list of simulations and returns a barplot with most varying substances

Usage

```
plotSubVar(simlist, metssel)
```

Arguments

simlist	A list of simulations (eval objects).
metssel	A vector with the name of exchange reactions of interest

plotTotFlux	<i>Function for plotting the overall change in reaction activity</i>
-------------	--

Description

The generic function plotTotFlux plots the time course of reactions with high variation in activity for an Eval object.

Usage

```
plotTotFlux(object, legendpos = "topright", num = 20)

## S4 method for signature 'Eval'
plotTotFlux(object, legendpos = "topright", num = 20)
```

Arguments

object	An object of class Eval.
legendpos	A character variable declaring the position of the legend
num	An integer defining the number of substrates to be plot

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
eval <- simEnv(arena,5)
plotTotFlux(eval)
```

redEval	<i>Function for reducing the size of an Eval object by collapsing the medium concentrations</i>
---------	---

Description

The generic function redEval reduces the object size of an Eval object.

Usage

```
redEval(object, time = "all")

## S4 method for signature 'Eval'
redEval(object, time = 1:length(object@medlist))
```

Arguments

object	An object of class Eval.
time	A number giving the simulation step of interest.

Details

The function redEval can be used to reduce the size of an Eval object from a simulation step.

Value

Returns an object of class Arena containing the organisms and substance conditions in simulation step time.

See Also

[Eval-class](#) and [Arena-class](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
eval <- simEnv(arena,5)
eval_reduce <- redEval(eval,5)
```

reset_screen

Reset plotting screen

Description

The function reset_screen set plotting window to default

Usage

```
reset_screen()
```

rmSubs	<i>Remove substances</i>
--------	--------------------------

Description

The generic function rmSubs removes all amounts of substances available in the arena for given compounds.

Usage

```
rmSubs(object, mediac)
```

```
## S4 method for signature 'Arena'
rmSubs(object, mediac)
```

Arguments

object	An object of class Arena.
mediac	A character vector giving the names of substances, which should be added to the environment (the default takes all possible substances).

selPheno	<i>Function for selecting phenotypes which occurred on the arena from specific iterations and species</i>
----------	---

Description

The generic function selPheno selects phenotypes from specific simulation step in an Eval object.

Usage

```
selPheno(object, time, type, reduce = F)
```

```
## S4 method for signature 'Eval'
selPheno(object, time, type, reduce = F)
```

Arguments

object	An object of class Eval.
time	A numeric vector giving the simulation steps which should be plotted.
type	A names indicating the species of interest in the arena.
reduce	A boolean variable indicating if the resulting matrix should be reduced.

Details

The phenotypes are defined by flux through exchange reactions, which indicate potential differential substrate usages.

Value

Returns a matrix with the substrate usage and the number of individuals using the phenotype.

See Also

[Eval-class](#) and [getPhenoMat](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
eval <- simEnv(arena,5)
selPheno(eval,time=5,type='ecoli_core_model',reduce=TRUE)
```

setKinetics

Function to set Michaelis-Menten kinetics for uptake of a substance

Description

The generic function setKinetics provides kinetics for exchange reactions.

Usage

```
setKinetics(object, exchangeR, Km, vmax)
```

```
## S4 method for signature 'Organism'
setKinetics(object, exchangeR, Km, vmax)
```

Arguments

object	An object of class Organisms.
exchangeR	Name of an exchange reaction
Km	Parameter Michaelis-Menten-Kinetics (in mM)
vmax	Parameter Michaelis-Menten-Kinetics (in mmol/(g*h))

sihumi_test	<i>Multi-species test data set</i>
-------------	------------------------------------

Description

Data contains results of a simulation with 8 gut bacteria for 10 time steps

Usage

```
data(sihumi_test)
```

Format

An object of class eval

simBac	<i>Function for one simulation iteration for objects of Bac class</i>
--------	---

Description

The generic function simBac implements all necessary functions for the individuals to update the complete environment.

Usage

```
simBac(object, arena, j, sublb, bacnum, sec_obj = "none",
        cutoff = 1e-06, pcut = 1e-06, with_shadow = FALSE)
```

```
## S4 method for signature 'Bac'
```

```
simBac(object, arena, j, sublb, bacnum, sec_obj = "none",
        cutoff = 1e-06, pcut = 1e-06, with_shadow = FALSE)
```

Arguments

object	An object of class Bac.
arena	An object of class Arena defining the environment.
j	The index of the organism of interest in orgdat.
sublb	A vector containing the substance concentrations in the current position of the individual of interest.
bacnum	integer indicating the number of bacteria individuals per gridcell
sec_obj	character giving the secondary objective for a bi-level LP if wanted.
cutoff	value used to define numeric accuracy.
pcut	A number giving the cutoff value by which value of objective function is considered greater than 0.
with_shadow	True if shadow cost should be stores (default off).

Details

Bacterial individuals undergo step by step the following procedures: First the individuals are constrained with `constrain` to the substrate environment, then flux balance analysis is computed with `optimizeLP`, after this the substrate concentrations are updated with `consume`, then the bacterial growth is implemented with `growth`, the potential new phenotypes are added with `checkPhen`, finally the additional and conditional functions `lysis`, `move` or `chemotaxis` are performed. In case of many compounds in the vector of `chemotaxis`, the change of the position takes place by the order of the compounds in the vector of `chemotaxis`. Can be used as a wrapper for all important bacterial functions in a function similar to `simEnv`.

Value

Returns the updated environment of the population parameter with all new positions of individuals on the grid and all new substrate concentrations.

See Also

[Bac-class](#), [Arena-class](#), [simEnv](#), `constrain`, `optimizeLP`, `consume`, `growth`, `checkPhen`, `lysis`, `move` and `chemotaxis`

Examples

NULL

simBac_par

Function for one simulation iteration for objects of Bac class

Description

The generic function `simBac_par` implements all necessary functions for the individuals to update the complete environment.

Usage

```
simBac_par(object, arena, j, sublb, bacnum, lproject, sec_obj = "none",
           cutoff = 1e-06, with_shadow = FALSE)
```

```
## S4 method for signature 'Bac'
simBac_par(object, arena, j, sublb, bacnum, lproject,
           sec_obj = "none", cutoff = 1e-06, with_shadow = FALSE)
```

Arguments

<code>object</code>	An object of class <code>Bac</code> .
<code>arena</code>	An object of class <code>Arena</code> defining the environment.
<code>j</code>	The index of the organism of interest in <code>orgdat</code> .

sublb	A vector containing the substance concentrations in the current position of the individual of interest.
bacnum	integer indicating the number of bacteria individuals per gridcell
lpobject	linear programming object (copy of organism@lpobj) that have to be a deep copy in parallel due to pointer use in sybil.
sec_obj	character giving the secondary objective for a bi-level LP if wanted.
cutoff	value used to define numeric accuracy
with_shadow	True if shadow cost should be stores (default off).

Value

Returns the updated environment of the population parameter with all new positions of individuals on the grid and all new substrate concentrations.

simEnv	<i>Main function for simulating all processes in the environment</i>
--------	--

Description

The generic function simEnv for a simple simulation of the environment.

Usage

```
simEnv(object, time, lrw = NULL, continue = FALSE, reduce = FALSE,
       diffusion = TRUE, diff_par = FALSE, cl_size = 2,
       sec_obj = "none", cutoff = 1e-06, pcut = 1e-06,
       with_shadow = TRUE, verbose = TRUE)
```

```
## S4 method for signature 'Arena'
simEnv(object, time, lrw = NULL, continue = FALSE,
       reduce = FALSE, diffusion = TRUE, diff_par = FALSE, cl_size = 2,
       sec_obj = "none", cutoff = 1e-06, pcut = 1e-06,
       with_shadow = TRUE, verbose = TRUE)
```

Arguments

object	An object of class Arena or Eval.
time	A number giving the number of iterations to perform for the simulation
lrw	A numeric value needed by solver to estimate array size (by default lrw is estimated in the simEnv() by the function estimate_lrw())
continue	A boolean indicating whether the simulation should be continued or restarted.
reduce	A boolean indicating if the resulting Eval object should be reduced
diffusion	True if diffusion should be done (default on).
diff_par	True if diffusion should be run in parallel (default off).

cl_size	If diff_par is true then cl_size defines the number of cores to be used in parallelized diffusion.
sec_obj	character giving the secondary objective for a bi-level LP if wanted. Use "mtf" for minimizing total flux, "opt_rxn" for optimizing a random reaction, "opt_ex" for optimizing a random exchange reaction, and "sumex" for optimizing the sum of all exchange fluxes.
cutoff	value used to define numeric accuracy
pcut	A number giving the cutoff value by which value of objective function is considered greater than 0.
with_shadow	True if shadow cost should be stored.
verbose	Set to false if no status messages should be printed.

Details

The returned object itself can be used for a subsequent simulation, due to the inheritance between Eval and Arena. The parameter for sec_obj can be used to optimize a bi-level LP with a secondary objective if wanted. This can be helpful to subselect the solution space and create less alternative optimal solution. The secondary objective can be set to "mtf" to minimize the total flux, to simulate minimal enzyme usage of an organisms. If set to "opt_rxn" or "opt_ex", the secondary objective is picked as a random reaction or exchange reaction respectively everytime a fba is performed. This means that every individual of a population will select a different secondary reaction to optimize. The "sumex" option maximizes the secretion of products.

Value

Returns an object of class Eval which can be used for subsequent analysis steps.

See Also

[Arena-class](#) and [Eval-class](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
          minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
eval <- simEnv(arena,5)
```

simEnv_par	<i>Main function for simulating in parallel all processes in the environment</i>
------------	--

Description

The generic function `simEnv_par` for a simple in parallel all simulation of the environment.

Usage

```
simEnv_par(object, time, lrw = NULL, continue = FALSE,
           reduce = FALSE, cluster_size = NULL, diffusion = TRUE,
           sec_obj = "none", cutoff = 1e-06, with_shadow = FALSE,
           verbose = TRUE)
```

```
## S4 method for signature 'Arena'
simEnv_par(object, time, lrw = NULL,
           continue = FALSE, reduce = FALSE, cluster_size = NULL,
           diffusion = TRUE, sec_obj = "none", cutoff = 1e-06,
           with_shadow = FALSE, verbose = TRUE)
```

Arguments

<code>object</code>	An object of class <code>Arena</code> or <code>Eval</code> .
<code>time</code>	A number giving the number of iterations to perform for the simulation
<code>lrw</code>	A numeric value needed by solver to estimate array size (by default <code>lrw</code> is estimated in the <code>simEnv()</code> by the function <code>estimate_lrw()</code>)
<code>continue</code>	A boolean indicating whether the simulation should be continued or restarted.
<code>reduce</code>	A boolean indicating if the resulting <code>Eval</code> object should be reduced
<code>cluster_size</code>	Number of cpu cores to be used.
<code>diffusion</code>	True if diffusion should be done (default on).
<code>sec_obj</code>	character giving the secondary objective for a bi-level LP if wanted.
<code>cutoff</code>	value used to define numeric accuracy
<code>with_shadow</code>	True if shadow cost should be stores (default off).
<code>verbose</code>	Set to false if no status messages should be printed.

Details

The returned object itself can be used for a subsequent simulation, due to the inheritance between `Eval` and `Arena`.

Value

Returns an object of class `Eval` which can be used for subsequent analysis steps.

See Also

[Arena-class](#) and [Eval-class](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
eval <- simEnv(arena,5)
```

simHum

Function for one simulation iteration for objects of Human class

Description

The generic function `simHum` implements all necessary functions for the individuals to update the complete environment.

Usage

```
simHum(object, arena, j, sublb, bacnum, sec_obj = "none",
        cutoff = 1e-06, pcut = 1e-06, with_shadow = FALSE)
```

```
## S4 method for signature 'Human'
simHum(object, arena, j, sublb, bacnum,
        sec_obj = "none", cutoff = 1e-06, pcut = 1e-06,
        with_shadow = FALSE)
```

Arguments

<code>object</code>	An object of class <code>Human</code> .
<code>arena</code>	An object of class <code>Arena</code> defining the environment.
<code>j</code>	The number of the iteration of interest.
<code>sublb</code>	A vector containing the substance concentrations in the current position of the individual of interest.
<code>bacnum</code>	integer indicating the number of bacteria individuals per gridcell
<code>sec_obj</code>	character giving the secondary objective for a bi-level LP if wanted.
<code>cutoff</code>	value used to define numeric accuracy.
<code>pcut</code>	A number giving the cutoff value by which value of objective function is considered greater than 0.
<code>with_shadow</code>	True if shadow cost should be stores (default off).

Details

Human cell individuals undergo the step by step the following procedures: First the individuals are constrained with `constrain` to the substrate environment, then flux balance analysis is computed with `optimizeLP`, after this the substrate concentrations are updated with `consume`, then the cell growth is implemented with `cellgrowth`, the potential new phenotypes are added with `checkPhen`, finally the conditional function `lysis` is performed. Can be used as a wrapper for all important cell functions in a function similar to `simEnv`.

Value

Returns the updated environment of the arena parameter with all new positions of individuals on the grid and all new substrate concentrations.

See Also

[Human-class](#), [Arena-class](#), [simEnv](#), `constrain`, `optimizeLP`, `consume`, `cellgrowth`, `checkPhen` and `lysis`

Examples

```
NULL
```

statPheno

Function for investigating a specific phenotype of an organism

Description

The generic function `statPheno` provides statistical and visual information about a certain phenotype.

Usage

```
statPheno(object, type_nr = 1, phenotype_nr, dict = NULL)
```

```
## S4 method for signature 'Eval'
statPheno(object, type_nr = 1, phenotype_nr,
          dict = NULL)
```

Arguments

<code>object</code>	An object of class <code>Eval</code> .
<code>type_nr</code>	A number indicating the Organism type of the phenotype to be investigated (from <code>orgdat</code>)
<code>phenotype_nr</code>	A number indicating the phenotype to be investigated (from <code>orgdat</code>)
<code>dict</code>	A character vector of all substance IDs with names that should be used instead of possibly cryptic IDs

Details

The phenotypes are defined by flux through exchange reactions, which indicate potential differential substrate usages.

See Also

[Eval-class](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
           minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
eval <- simEnv(arena,5)
statPheno(eval, type_nr=1, phenotype_nr=2)
```

stirEnv

Function for stirring/mixing the complete environment

Description

The generic function `stirEnv` simulates the event of mixing all substrates and organisms in the environment.

Usage

```
stirEnv(object, sublb)

## S4 method for signature 'Arena'
stirEnv(object, sublb)
```

Arguments

<code>object</code>	An object of class <code>Arena</code> .
<code>sublb</code>	A matrix with the substrate concentration for every individual in the environment based on their x and y position.

Details

The stirring is implemented as a random permutation of organism positions and the equalization of of all substrate concentrations.

Value

Returns the substrate concentration for every individual in the environment with substrates as well as x and y positions as columns and rows for each organism.

See Also

[Arena-class](#) and [getSublb](#)

Examples

```
data(Ec_core, envir = environment()) #get Escherichia coli core metabolic model
bac <- Bac(Ec_core,deathrate=0.05,
          minweight=0.05,growtype="exponential") #initialize a bacterium
arena <- Arena(n=20,m=20) #initialize the environment
arena <- addOrg(arena,bac,amount=10) #add 10 organisms
arena <- addSubs(arena,40) #add all possible substances
sublb <- getSublb(arena)
stirEnv(arena,sublb)
```

Substance-class

Structure of the S4 class "Substance"

Description

Structure of the S4 class Substance representing substances in the environment which can be produced or consumed.

Slots

smax A number representing the start concentration of the substance for each grid cell in the environment.

diffmat A sparse matrix containing all concentrations of the substance in the environment.

name A character vector representing the name of the substance.

id A character vector representing the identifier of the substance.

diffunc A character vector ("pde","cpp" or "r") describing the function for diffusion.

difspeed A number indicating the diffusion rate (given by cm^2/h). Default is set to glucose diffusion in a aqueous solution ($6.7\text{e-}6 \text{ cm}^2/\text{s} * 3600 \text{ s/h} = 0.02412 \text{ cm}^2/\text{h}$).

advspeed A number indicating the advection rate in x direction (given by cm/h).

diffgeometry Diffusion coefficient defined on all grid cells (initially set by constructor).

pde Choose diffusion transport reaction to be used (default is diffusion only)

boundS A number defining the attached amount of substance at the boundary (Warning: boundary-function must be set in pde!)

Substance-constructor *Constructor of the S4 class Substance*

Description

The constructor to get a new object of class Substance

Usage

```
Substance(n, m, smax, gridgeometry, difspeed = 0.02412, advspeed = 0,
  occupyM, Dgrid = NULL, Vgrid = NULL, diffmat = NULL,
  template = FALSE, ...)
```

Arguments

n	A number giving the horizontal size of the environment.
m	A number giving the vertical size of the environment.
smax	A number representing the start concentration of the substance for each grid cell in the environment.
gridgeometry	A list containing grid geometry parameter
difspeed	A number indicating the diffusion speed in x and y direction (given by cm ² /h). For more complex setup define Dgrid.
advspeed	A number indicating the advection speed in x direction (given by cm/h). For more complex setup define Vgrid.
occupyM	A matrix indicating grid cells that are obstacles
Dgrid	A matrix indicating the diffusion speed in x and y direction (given by cm ² /h).
Vgrid	A number indicating the advection speed in x direction (given by cm/h).
diffmat	A matrix with spatial distributed initial concentrations (unit in fmol) (if not set, a homogenous matrix using smax is created)
template	True if diffmat matrix should be used as template only (will be multiplied with smax to obtain concentrations)
...	Arguments of Substance-class

Value

Object of class Substance

unit_conversion	<i>Function for unit conversion</i>
-----------------	-------------------------------------

Description

The generic function unit_conversion converts units for e.g. substance concentrations

Usage

```
unit_conversion(object, unit)
```

```
## S4 method for signature 'Arena'  
unit_conversion(object, unit)
```

Arguments

object	An object of class Arena or Eval.
unit	Unit to be converted to fmol/cell

Value

Conversion factor

usd	<i>Computer standard deviation upper bound</i>
-----	--

Description

Helper function to get upper error bounds in plotting

Usage

```
usd(y)
```

Arguments

y	Vector with numbers
---	---------------------

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