# Package: DstarM (via r-universe)

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**Type** Package

**Title** Analyze Two Choice Reaction Time Data with the D*M Method

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**Description** A collection of functions to estimate parameters of a diffusion model via a D*M analysis. Build in models are: the Ratcliff diffusion model, the RWiener diffusion model, and Linear Ballistic Accumulator models. Custom models functions can be specified as long as they have a density function.

**License** GPL (>= 2)

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---

`chisq` *Calculates the distance between two probability densities.*

---

## Description

Calculates the distance between two probability densities.

## Usage

```
chisq(tt, a, b)

battacharyya(tt, a, b)

hellinger(tt, a, b)
```

## Arguments

| | |
|---|---|
| `tt` | the time grid on which the densities are evaluated. |
| `a` | a vector with values of the first density. |
| `b` | a vector with values of the second density. |

## Value

The distance between densities a and b.

## Examples

```
# Lets simulate a bunch of parameters and compare the three distance measures.

tt = seq(0, 5, .001)
parsMatV = cbind(.8, seq(0, 5, .5), .5, .5, .5) # differ only in drift speed
parsMatA = cbind(seq(.5, 2, .15), 2, .5, .5, .5)# differ only in boundary
# calculate densities for all these parameters
dV = apply(parsMatV, 1, function(x, tt) Voss.density(tt, x, boundary = 'upper'), tt = tt)
dA = apply(parsMatA, 1, function(x, tt) Voss.density(tt, x, boundary = 'upper'), tt = tt)
# make plots of the densities
matplot(tt, dA, xlim = c(0, .6), main = 'Densities with different Boundary',
        col = rainbow(ncol(dA)),type = 'l', lty = 1, las = 1, bty = 'n',
        xlab = 'Time', ylab = 'Density')
legend('topright', lty = 1, bty = 'n', col = rainbow(ncol(dA)),
       legend = paste('a = ', parsMatA[, 1]))
matplot(tt, dV, xlim = c(0, .6), main = 'Densities with different Drift Speed',
        col = rainbow(ncol(dV)), type = 'l', lty = 1, las = 1, bty = 'n',
        xlab = 'Time', ylab = 'Density')
legend('topright', lty = 1, bty = 'n', col = rainbow(ncol(dV)),
       legend = paste('v = ',parsMatV[, 2]))
# empty matrices for data storage
distMatV = matrix(NA, nrow = ncol(dV) - 1, ncol = 3,
                  dimnames = list(NULL, c('Chisq', 'Bhattacharyya', 'Hellinger')))
distMatA = matrix(NA, nrow = ncol(dA) - 1, ncol = 3,
                  dimnames = list(NULL, c('Chisq', 'Bhattacharyya', 'Hellinger')))
# calculate distances between densities in column i and i + 1.
# this is done using three different distance measures
for (i in 1:(ncol(dA) - 1)) {
  distMatV[i, ] = c(chisq(tt, dV[, i], dV[, i + 1]),
                    battacharyya(tt, dV[, i], dV[, i + 1]),
                    hellinger(tt, dV[, i], dV[, i + 1]))
  distMatA[i, ] = c(chisq(tt, dA[, i], dA[, i + 1]),
                    battacharyya(tt, dA[, i], dA[, i + 1]),
                    hellinger(tt, dA[, i], dA[, i + 1]))
}
# The three distance measures correlate highly for differences in Boundary
cor(distMatA)
# The battacharyya distance measures does not correlate with the others
# when calculating differences in drift speed
cor(distMatV)
```

---

`chisqFit` *Calculate model fit*

---

## Description

Calculate model fit

## Usage

```
chisqFit(resObserved, data, DstarM = FALSE, tt = NULL, formula = NULL)
```

## Arguments

| | |
|---|---|
| resObserved | either output from estObserved or a matrix containing custom densities to calculate the fitness for. |
| data | A dataframe containing data. |
| DstarM | Logical. Should the DstarM fit measure be calculated or the traditional fit measure? |
| tt | time grid custom densities where calculated on. Should only be supplied if resOberved is a matrix containing custom densities |
| formula | Optional formula argument, for when columns names in the data are different from those used to obtain the results. |

## Details

This function allows a user to manually calculate a chi-square goodness of fit measure for model densities. This is useful for comparing a traditional analysis and a D*M analysis. For completion, this function can also calculate a D*M fit measure. We do not recommend usage of the D*M measure. While the chi-square fit measure is identical to the value of the optimizer when fitting, the DstarM fit measure is not equal to that of a DstarM analysis. This is because this function calculates the DstarM fit measure on the complete distribution, not on the model distributions, as is done during the optimization.

## Examples

```
tt = seq(0, 5, .1)
pars = c(.8, 2, .5, .5, .5, # condition 1
         .8, 3, .5, .5, .5,  # condition 2
         .8, 4, .5, .5, .5)  # condition 3
pdfND = dbeta(tt, 10, 30)

# simulate data
allDat = simData(n = 3e3, pars = pars, tt = tt, pdfND = pdfND, return.pdf = TRUE)
truePdf = allDat$pdfUnnormalized
dat = allDat$dat
chisqFit(resObserved = truePdf, data = dat, tt = tt)
## Not run:
# estimate it
define restriction matrix
restr = matrix(1:5, 5, 3)
restr[2, 2:3] = 6:7 # allow drift rates to differ
# fix parameters for speed up
fixed = matrix(c('z1', 'a1 / 2', 'sz1', .5, 'sv1', .5), 2, 3)
resD = estDstarM(data = dat, tt = tt, restr = restr, fixed = fixed,
                 Optim = list(parallelType = 1))
resN = estND(resD, Optim = list(parallelType = 1))
```

```
resO = estObserved(resD, resN, data = dat)
resO$fit # proper fit

## End(Not run)
```

---

`Density` *Density function*

---

## Description

Density function

## Usage

```
Density(rt, tt)
```

## Arguments

| | |
|---|---|
| `rt` | vector of reaction times |
| `tt` | grid to evaluate the density on |

## Details

Can be passed to the argument `densityMethod` of `estDstarM`. This function is a minimal example to use as custom smoothing function.

## Value

a vector of `length(tt)`

## Examples

```
x <- rgamma(1e5, 1, 1)
tt <- seq(0, 5, .01)
d <- Density(x, tt)
hist(x, freq = FALSE)
lines(tt, DstarM:::Density(x, tt))
```

estCdf *Estimate cumulative distribution for D*M models*

## Description

Estimate cumulative distribution for D*M models

## Usage

```
estCdf(x)
```

## Arguments

| | |
|---|---|
| x | Any density function to calculate a cumulative distribution for. The code is designed for input of class DstarM but other input is also accepted. Other input can be either a matrix where columns represent densities or a single vector representing a density. |

## Details

Cumulative distributions functions are calculated by: `cumsum(x) / sum(x)`. This method works well enough for our purposes. The example below shows that the ecdf functions seems to work slightly better. However, this estimates a cdf from raw data and does not transform a pdf into a cdf and is therefore not useful for D*M models.

## Value

Cumulative density function(s). If the input was a matrix, a matrix of cumulative density functions is returned.

## Examples

```
x = rnorm(1000)
xx = seq(-5, 5, .1)
approx1 = stats::ecdf(x)(xx)
approx2 = estCdf(dnorm(xx, mean(x), sd(x)))
trueCdf = pnorm(xx)
matplot(xx, cbind(trueCdf, approx1, approx2), type = c('l', 'p', 'p'),
        lty = 1, col = 1:3, pch = 1, bty = 'n', las = 1, ylab = 'Prob')
legend('topleft', legend = c('True Cdf', 'Stats Estatimation', 'DstarM Estimation'),
       col = 1:3, lty = c(1, NA, NA), pch = c(NA, 1, 1), bty = 'n')
```

`estDstarM` *Do a D*M analysis*

## Description

Do a D*M analysis

## Usage

```
estDstarM(
  formula = NULL,
  data,
  tt,
  restr = NULL,
  fixed = list(),
  lower,
  upper,
  Optim = list(),
  DstarM = TRUE,
  SE = 0,
  oscPdf = TRUE,
  splits = rep(0L, (ncondition)),
  forceRestriction = TRUE,
  mg = NULL,
  h = 1,
  pars,
  fun.density = Voss.density,
  args.density = list(),
  fun.dist = chisq,
  args.dist = list(tt = tt),
  verbose = 1L,
  useRcpp = TRUE
)
```

## Arguments

| | |
|---|---|
| `formula` | A formula object of the form: `binary response ~ reaction time + condition1 * condition2 * ... conditionN`. |
| `data` | A dataframe for looking up data specified in formula. For backwards compatibility this can also be with: a column named `rt` containing response times in ms, a column named `response` containing at most 2 response options, and an optional column named `condition` containing a numeric index as to which conditions observations belong. |
| `tt` | A time grid on which the density function will be evaluated. Should be larger than the highest observed reaction time. |

`restr` A restriction matrix where each column depicts one condition. The number of rows should match the number of parameters (and be equal to the length of lower). The contents of `restr` should be numbers, identical numbers means that these parameters (either within or between condition) will be constrained. Different numbers means parameters will not be constrained.

`fixed` A matrix that allows for fixing parameters to certain values.

`lower` Should be a vector containing lower bounds for each parameter. Has a default if `fun.density == Voss.density`.

`upper` Should be a vector containing upper bounds for each parameter. Has a default if `fun.density == Voss.density`.

`Optim` a named list with identical arguments to `DEoptim.control`. In addition, if `verbose == TRUE` `Optim$steptol` can be a vector, i.e. `c(200, 50, 10)` means: Do 200 iterations then check for convergence, do 50 iterations then check for convergence, check every 10 iterations for convergence until itermax is reached. Defaults to `Optim = list(reltol = 1e-6, itermax = 1e3,steptol = 50, CR = .9, trace = 0, parallelType = 0)`.

`DstarM` If TRUE a D*M analysis is done, otherwise the Chi square distance between data and model is minimized.

`SE` positive value, how many standard error to add to the variance to relax the variance restriction a bit.

`oscPdf` Logical, if TRUE check for oscillations in calculated densities and remove densities with oscillations.

`splits` Numeric vector determining which conditions have an equal nondecision density. Identical values in two positions indicate that the conditions corresponding to the indices of those values have an identical nondecision distribution.

`forceRestriction` if TRUE the variance restriction is enforced.

`mg` Supply a data density, useful if a uniform kernel approximation does not suffice. Take care that densities of response categories within conditions are degenerate and therefore integrate to the proportion a category was observed (and not to 1).

`h` bandwidth of a uniform kernel used to generate data based densities.

`pars` Optional parameter vector to supply if one wishes to evaluate the objective function in a given parameter vector. Only used if `itermax` equal zero.

`fun.density` Function used to calculate densities. See details.

`args.density` A names list containing additional arguments to be send to fun.density.

`fun.dist` Function used to calculate distances between densities. Defaults to a chi-square distance.

`args.dist` A named list containing additional arguments to be send to fun.dist.

`verbose` Numeric, should intermediate output be printed? Defaults to 1, higher values result in more progress output. Estimation will speed up if set to 0. If set to TRUE, `Optim$trace` will be forced to 0, hereby disabling the build in printing of `DEoptim`. To enable the printing of `DEoptim`, set `verbose` to 0 and specify `Optim$trace`. `Optim`. If set to 1, ETA refers to the expected maximum time until completion (when the iterations limit is reached).

useRcpp — Logical, setting this to true will make the objective function use an Rcpp implementation of Voss.density with the distance function chisq. This gains speed at the cost of flexibility.

## Details

Response options will be alphabetically sorted and the first response option will be treated as the 'lower' option. This means that if the observed proportion of the first response options is higher, the drift speed will most likely be negative.

fun.density allows a user to specify a custom density function. This function must (at least) take the following arguments: t: a vector specifying at which time points to calculate the density pars: a parameter vector boundary: character 'upper' or 'lower' specifying for which response option the density will be calculated. DstarM: Logical, if TRUE the density should not describe the nondecision density, if FALSE it should describe the nondecision density. Any additional arguments can be passed to fun.density via the argument args.density. If one intends to use a custom density function it is recommended to test the function first with testFun. When specifying a custom density function it is probably also necessary to change the lower and upper bounds of the parameter space.

For purposes of speed, the function can be run in parallel by providing the argument Optim = list(parallelType = 1). See DEoptim.control for details. Also, for Ratcliff models the objective function has been rewritten in Rcpp. This limits some functionality but does result in a faster estimation. Usage of Rcpp can be enabled via useRcpp = TRUE.

When verbose is set to 1, the ETA is an estimated of the time it takes to execute ALL iterations. Convergence can (and is usually) reached before then.

## Value

Returns a list of class DstarM.fitD that contains:

Bestvals — Named numeric vector. Contains the best parameter estimates.

fixed — Numeric vector. Contains the best parameter estimates.

GlobalOptimizer — List. All output from the call to DEoptim

Debug — List. contains the number of DEoptim iterations, the number of function evaluation of the objective function, and the maximum number of iterations.

note — String. A possible note that is used for summary purposes

tt — Numeric vector. Contains the time grid used.

g.hat — Numeric matrix. Named columns represent the (possibly smoothed) densities of the data distribution of each condition-response pair.

modelDist — Numeric matrix. Named columns represent the densities of the model evaluated at grid tt with parameters Bestvals.

ncondition — Numeric scalar. The number of conditions

var.data — Numeric vector. The variance of each condition-response pair. There are as many values as hypothesized nondecision densities.

var.m — Numeric vector. The variance of the model distributions. There are as many values as hypothesized nondecision densities.

| | |
|---|---|
| restr.mat | Numeric matrix. Contains the restrictions used. |
| splits | Numeric vector. Equal to the input argument with the same name. |
| n | Numeric scalar. The total number of observations. |
| DstarM | Logical. Equal to the input argument with the same name. |
| fun.density | Function. Equal to the input argument with the same name. |
| fun.dist | Function. Equal to the input argument with the same name. |
| h | Scalar. Equal to the input argument with the same name. |
| args.density | Named list. Equal to the input argument with the same name. |
| args.dist | Named list. Equal to the input argument with the same name. |

## Examples

```
# simulate data with three stimuli of different difficulty.
# this implies different drift rates across conditions.
# define a time grid. A more reasonable stepsize is .01; this is just for speed.
tt = seq(0, 5, .1)
pars = c(.8, 2, .5, .5, .5, # condition 1
        .8, 3, .5, .5, .5,  # condition 2
        .8, 4, .5, .5, .5)  # condition 3
pdfND = dbeta(tt, 10, 30)
# simulate data
data = simData(n = 3e3, pars = pars, tt = tt, pdfND = pdfND)
# define restriction matrix
restr = matrix(1:5, 5, 3)
restr[2, 2:3] = 6:7 # allow drift rates to differ
# fix variance parameters
fixed = matrix(c('sz1', .5, 'sv1', .5), 2, 2)
## Not run:
# Run D*M analysis
res = estDstarM(data = data, tt = tt, restr = restr, fixed = fixed)
coef(res)
summary(res)

## End(Not run)
```

---

# estND *Estimate nondecision density*

---

## Description

Estimate nondecision density

## Usage

```
estND(
  res,
  tt = NULL,
  data = NULL,
  h = res$h,
  zp = 5,
  upper.bound = 1,
  lower.bound = 0,
  Optim = list(),
  verbose = TRUE,
  dist = NULL,
  NDindex,
  max = 100,
  useRcpp = TRUE
)
```

## Arguments

| | |
|---|---|
| res | an object of class D*M. |
| tt | optional timegrid if the nondecision density is to be estimated at a different grid than the model density. |
| data | if tt is specified then the original dataset must be supplied too. |
| h | Optional smoothing parameter to be used when estimating the nondecision model on a different time grid than the decision model. If omitted, the smoothing parameter of the decision model is used. |
| zp | Zero padding the estimated nondecision density by this amount to avoid numerical artefacts. |
| upper.bound | An upper bound for the nondecision density. Defaults to one. Lowering this bound can increase estimation speed, at the cost of assuming that the density of the nondecision distribution is zero past this value. |
| lower.bound | A lower bound for the nondecision density. Defaults to zero. Increasing this bound can increase estimation speed, at the cost of assuming that the density of the nondecision distribution is zero past this value. |
| Optim | a named list with identical arguments to DEoptim.control. In addition, if verbose == TRUE Optim$steptol can be a vector, i.e. c(200, 50, 10) means: Do 200 iterations then check for convergence, do 50 iterations then check for convergence, check every 10 iterations for convergence until itermax is reached. If there are multiple nondecision distributions to estimate, one can supply different estimation parameters for every nondecision distribution by supplying Optim as a list of lists. Every sublists then corresponds to parameters for one nondecision distribution and should consist of arguments for DEoptim.control. Defaults to Optim = list(reltol = 1e-6, itermax = 1e4, steptol = 200, CR = .9, trace = 0). |
| verbose | Numeric, should intermediate output be printed? Defaults to 1, higher values result in more progress output. Estimation will speed up if set to 0. If |

nonzero, Optim$trace will be forced to 0, hereby disabling the build in printing of DEoptim. To enable the printing of DEoptim, set verbose to 0 and specify Optim$trace.

dist — A matrix where columns represent nondecision distributions. If this argument is supplied then the objective function will be evaluated in these values.

NDindex — A vector containing indices of which nondecision distributions to estimate. If omitted, all nondecision distributions that complement the results in res are estimated.

max — A positive float which indicates the maximum height of the nondecision distribution. If estimated nondecision distributions appear chopped of or have a lot of values at this max value it is recommended to re-estimate the nondecision distributions with a higher max value. Increasing the max value without reason will increase the size of the parameter space and slow the estimation procedure.

useRcpp — Logical, setting this to true will make use of an Rcpp implementation of the objective function. This gains speed at the cost of flexibility.

## Details

When verbose is set to 1, the ETA is an estimated of the time it takes to execute ALL iterations. Convergence can (and is usually) reached before then.

## Examples

```
# simulate data with three stimuli of different difficulty.
# this implies different drift rates across conditions.
# define a time grid. A more reasonable stepsize is .01; this is just for speed.
tt = seq(0, 5, .1)
pars = c(.8, 2, .5, .5, .5, # condition 1
        .8, 3, .5, .5, .5, # condition 2
        .8, 4, .5, .5, .5) # condition 3
pdfND = dbeta(tt, 10, 30)
# simulate data
dat = simData(n = 3e5, pars = pars, tt = tt, pdfND = pdfND)
# define restriction matrix
restr = matrix(1:5, 5, 3)
restr[2, 2:3] = 6:7 # allow drift rates to differ
# fix variance parameters
fixed = matrix(c('sz1', .5, 'sv1', .5), 2, 2)
## Not run:
# Run D*M analysis
res = estDstarM(data = dat, tt = tt, restr = restr, fixed = fixed)
# Estimate nondecision density
resND = estND(res)
plot(resND)
lines(tt, pdfND, type = 'b', col = 2)

## End(Not run)
```

`estObserved` *Estimate observed data density*

## Description

Estimates the density of the observed data by convoluting the estimated decision distributions with the estimated nondecision distributions. If a traditional analysis was run the argument resND can be omitted.

## Usage

```
estObserved(
  resDecision,
  resND = NULL,
  data = NULL,
  interpolateND = FALSE,
  tt = NULL
)
```

## Arguments

| | |
|---|---|
| `resDecision` | output of `estDstarM`. |
| `resND` | output of `estND`. |
| `data` | Optional. If the data used to estimate the decision model is supplied additional fitmeasures are calculated. |
| `interpolateND` | Logical. If the decision model and nondecision model have been estimated on different time grids, should the rougher time grid be interpolated to match the smaller grid? If FALSE (the default) the decision model will be recalculated on the grid of the nondecision model. This tends to produce better fit values. |
| `tt` | Optional time grid to recalculate the model densities on. Unused in case of a DstarM analysis. |

## Value

Returns a list of class `DstarM.fitObs` that contains:

| | |
|---|---|
| `obsNorm` | A matrix containing normalized densities of each condition response pair. |
| `obs` | A matrix containing unnormalized densities of each condition response pair. |
| `tt` | The time grid used. |
| `fit` | A list containing the values of the objective function for the total model ($total), for the decision model ($Decision) and for the nondecision distribution(s) ($ND). |
| `npar` | The number of parameters used in the decision model. |
| `obsIdx` | A numeric vector containing indices of any not observed condition-response pairs. |

## Examples

```
# simulate data with three stimuli of different difficulty.
# this implies different drift rates across conditions.
# define a time grid. A more reasonable stepsize is .01; this is just for speed.
tt = seq(0, 5, .1)
pars = c(.8, 2, .5, .5, .5, # condition 1
        .8, 3, .5, .5, .5, # condition 2
        .8, 4, .5, .5, .5) # condition 3
pdfND = dbeta(tt, 10, 30)
# simulate data
lst = simData(n = 3e5, pars = pars, tt = tt, pdfND = pdfND, return.pdf = TRUE)
dat = lst$dat
# define restriction matrix
restr = matrix(1:5, 5, 3)
restr[2, 2:3] = 6:7 # allow drift rates to differ
# fix variance parameters
fixed = matrix(c('sz1', .5, 'sv1', .5), 2, 2)
## Not run:
# Run D*M analysis
resD = estDstarM(dat = dat, tt = tt, restr = restr, fixed = fixed)
# Estimate nondecision density
resND = estND(resD)
# Estimate observed density
resObs = estObserved(resD, resND)
# plot histograms with overlayed
# densities per condition-response pair
plotObserved(resObserved = resObs, data = dat,
            xlim = c(0, 1))
# plot estimated and true densities
plot(resObs, col = rep(1:3, each = 2), xlim = 0:1)
matlines(tt, lst$pdfNormalized, col = rep(1:3, each = 2), lty = 2)

## End(Not run)
```

---

`estQdf` *Estimate quantiles of distribution*

---

## Description

Estimate quantiles of distribution

## Usage

```
estQdf(p, x, cdf)
```

## Arguments

| | |
|---|---|
| p | A vector of probabilities. |
| x | The x-axis values corresponding to the cumulative distribution function. |
| cdf | A cumulative distributions function, i.e. output of `estCdf`. |

## Details

Quantiles are obtained in the following manner. For p = 0 and p = 1, the minimum and maximum of x is used. For other probabilities the quantiles are obtained via `q[i] = uniroot(x, cdf - p[i])$root`. Y values are interpolated via `approxfun`.

## Value

Quantiles of cumulative distribution function(s). If the input was a matrix of cumulative distributions functions, a matrix of quantiles is returned.

## Examples

```
x = seq(-9, 9, .1) # x-grid
d = dnorm(x) # density functions
p = seq(0, 1, .2) # probabilities of interest
cEst = estCdf(d) # estimate cumulative distribution functions
qEst = estQdf(p = p, x = x, cdf = cEst) # estimate quantiles
plot(x, cEst, bty = 'n', las = 1, type = 'l', ylab = 'Probability') # plot cdf
abline(h = p, v = qEst, col = 1:6, lty = 2) # add lines for p and for obtained quantiles
points(x = qEst, y = p, pch = 18, col = 1:6, cex = 1.75) # add points for intersections
```

---

# getPdfs *(Re)Calculate model densities with given parameters and time grid*

---

## Description

This function is a convenience function for calculating model pdfs for multiple sets of parameters at a specified timegrid. If `resDecision` is supplied, the density function and any additional arguments for the density function will be extracted from that object. If `pars` is missing these will also be extracted from this object. This function is intended to recalculate model densities at a new timegrid.

## Usage

```
getPdfs(
  resDecision,
  tt,
  pars,
  DstarM = TRUE,
  fun.density = Voss.density,
  args.density = list()
)
```

### Arguments

| | |
|---|---|
| resDecision | output of estDstarM. |
| tt | Time grid to calculate the model densities on. |
| pars | Model parameters, can be a matrix where every column is a set of parameters. |
| DstarM | Logical. Do the model pdfs also describe the nondecision distribution? |
| fun.density | density function to calculate pdfs from. |
| args.density | Additional arguments for fun.density |

### Value

A matrix containing model pdfs.

---

## getSter *Estimate variance of nondecision density*

---

### Description

Estimate variance of nondecision density

### Usage

```
getSter(res)
```

### Arguments

| | |
|---|---|
| res | An object of class D*M. |

### Details

The object res can either be output from estDstarM or output from estND. if the former is supplied, getSter attempts to calculate the variance of the nondecision distribution by subtracting the variance of the model distribution from the variance of the data distribution. If the latter is supplied, the variance is calculated by integrating the nondecision distribution.

## getTer *Calculate Mean of the nondecision distribution.*

### Description

Calculate Mean of the nondecision distribution.

### Usage

```
getTer(res, data, formula = NULL)
```

### Arguments

| | |
|---|---|
| res | An object of class D*M. |
| data | The data object used to create res. |
| formula | Optional formula argument, for when columns names in the data are different from those used to obtain the results. |

### Details

The object res can either be output from estDstarM or output from estND. If the former is supplied it is also necessary to supply the data used for the estimation. The mean will then be estimated by subtracting the mean of the model densities from the mean of the data density. If the latter is supplied than this is not required; the mean will be calculated by integrating the nondecision distribution.

### Value

A vector containing estimates for the mean of the nondecision densities.

## normalize *Normalize two pdfs*

### Description

Normalize two pdfs

### Usage

```
normalize(x, tt, props = NULL)
```

### Arguments

| | |
|---|---|
| x | Probability density function(s) evaluated at grid x. Input should be either a vector or matrix. If input is a matrix, each column represents a single pdf. |
| tt | a numeric grid defined in x. |
| props | the value each density should integrate to. |

## Examples

```
tt <- seq(0, 9, length.out = 1e4)
# 2 poper densities
x1 <- cbind(dexp(tt, .5), dexp(tt, 2))
# still 2 poper densities
x2 <- normalize(10*x1, tt)
# 2 densities that integrate to .5
x3 <- normalize(x1, tt, props = c(.5, .5))
# plot the results
matplot(tt, cbind(x1, x2, x3), type = "l", ylab = "density",
        col = rep(1:3, each = 2), lty = rep(1:2, 3), las = 1, bty = "n")
legend("topright", legend = rep(paste0("x", 1:3), each = 2),
       col = rep(1:3, each = 2), lty = rep(1:2, 3), bty = "n")
```

---

`obsQuantiles` *Calculate model fit*

---

## Description

This function is nothing but a wrapper for `quantile`.

## Usage

```
obsQuantiles(data, probs = seq(0, 1, 0.01), what = "cr")
```

## Arguments

| | |
|---|---|
| `data` | A dataframe with: a column named `rt` containing response times in ms, a column named `response` containing at most 2 response options, and an optional column named `condition` containing a numeric index as to which conditions observations belong. |
| `probs` | vector of probabilities for which the corresponding values should be called |
| `what` | Character. `'cr'` if the quantiles are to be calculated per condition-response pair, `'c'` if the quantiles are to be calculated per condition, and `'r'` if the quantiles are to be calculated per response. |

## Examples

```
tt = seq(0, 5, .01)
pars = c(.8, 2, .5, .5, .5, # condition 1
         .8, 3, .5, .5, .5,  # condition 2
         .8, 4, .5, .5, .5)  # condition 3
pdfND = dbeta(tt, 10, 30)
# simulate data
data = simData(n = 3e3, pars = pars, tt = tt, pdfND = pdfND)
probs = seq(0, 1, .01)
q = obsQuantiles(data, probs = probs)
matplot(probs, q, type = 'l', las = 1, bty = 'n')
```

plotObserved *Plot quantiles of data against model implied quantiles.*

## Description

Plots histograms for each condition-response pair/ condition/ response with overlayed estimated densities.

## Usage

```
plotObserved(
  resObserved,
  data,
  what = c("cr", "c", "r"),
  layout = NULL,
  main = NULL,
  linesArgs = list(),
  ggplot = FALSE,
  prob = seq(0, 1, 0.01),
  probType = 3,
  ...
)
```

## Arguments

| | |
|---|---|
| resObserved | output of estObserved. |
| data | The dataset used to estimate the model. |
| what | What to plot. Can be 'cr' for 'condition-response pairs, 'c' for condition, and 'r' for response. |
| layout | An optional layout matrix. |
| main | an optional vector containing names for each plot. |
| linesArgs | A list containing named arguments to be passed to lines. |
| ggplot | Deprecated and ignored. |
| prob | Should a qqplot of observed vs model implied quantiles be plotted? By default, it is seq(0, 1, .01), the probabilities between 0 and 1 to compare the model implied quantiles to the observed quantiles. If this argument is NULL, then a histogram overlayed with model implied densities will be plotted. Internally, estQdf is used for generating quantiles. |
| probType | A numeric value defining several plotting options. 0 does nothing, 1 removes the 0% quantile, 2 removes the 100% quantile and 3 removes both the 0% and 100% quantile. |
| ... | Further arguments to be passed to hist. |

## Details

Keep in mind when using `what = 'c'` or `what = 'r'` pdfs are simply averaged, not weighted to the number of observed responses.

## Value

if ggplot is FALSE `invisible()`, otherwise a list

## Examples

```
# simulate data with three stimuli of different difficulty.
# this implies different drift rates across conditions.
# define a time grid. A more reasonable stepsize is .01; this is just for speed.
tt = seq(0, 5, .1)
pars = c(.8, 2, .5, .5, .5, # condition 1
        .8, 3, .5, .5, .5, # condition 2
        .8, 4, .5, .5, .5) # condition 3
pdfND = dbeta(tt, 10, 30)
# simulate data
lst = simData(n = 3e5, pars = pars, tt = tt, pdfND = pdfND, return.pdf = TRUE)
dat = lst$dat
# define restriction matrix
restr = matrix(1:5, 5, 3)
restr[2, 2:3] = 6:7 # allow drift rates to differ
# fix variance parameters
fixed = matrix(c('sz1', .5, 'sv1', .5), 2, 2)
## Not run:
# Run D*M analysis
resD = estDstarM(dat = dat, tt = tt, restr = restr, fixed = fixed)
# Estimate nondecision density
resND = estND(resD)
# Estimate observed density
resObs = estObserved(resD, resND)
# plot histograms with overlayed
# densities per condition-response pair
plotObserved(resObserved = resObs, data = dat,
            xlim = c(0, 1))
# plot estimated and true densities
plot(resObs, col = rep(1:3, each = 2), xlim = 0:1)
matlines(tt, lst$pdfNormalized, col = rep(1:3, each = 2), lty = 2)
# other uses of plotObserved
plotObserved(resObserved = resObs, data = dat, what = 'cr', xlim = c(0, 1))
plotObserved(resObserved = resObs, data = dat, what = 'c', xlim = c(0, 1))
plotObserved(resObserved = resObs, data = dat, what = 'r', xlim = c(0, 1))

## End(Not run)
```

`rtDescriptives` *Descriptives of reaction time data*

## Description

Descriptives of reaction time data

## Usage

```
rtDescriptives(formula = NULL, data, plot = TRUE, verbose = TRUE)
```

## Arguments

| | |
|---|---|
| `formula` | A formula object of the form: `binary response ~ reaction time + condition1 * condition2` |
| `data` | A dataframe for looking up data specified in formula. For backwards compatibility this can also be with: a column named `rt` containing response times in ms, a column named `response` containing at most 2 response options, and an optional column named `condition` containing a numeric index as to which conditions observations belong. |
| `plot` | Logical, should a density plot of all condition-response pairs be made? |
| `verbose` | Logical, should a table of counts and proportions be printed? |

## Details

This function and `rtHist` are helper functions to inspect raw data.

## Value

Invisibly returns an object of class 'D*M'. It's first element is `table` and contains raw counts and proportions for condition response pairs, conditions, and responses. It's second element `plot` contains a ggplot object.

## Examples

```
tt <- seq(0, 5, .01)
pars <- matrix(.5, 5, 2)
pars[1, ] <- 1
pars[2, ] <- c(0, 2)
dat <- simData(n = 3e3, pars = pars, tt = tt, pdfND = dbeta(tt, 10, 30))
x <- rtDescriptives(data = dat)

print(x$table, what = 'cr')
print(x$table, what = 'c')
print(x$table, what = 'r')
```

`rtHist` *Make histograms of reaction time data*

## Description

Make histograms of reaction time data

## Usage

```
rtHist(data, what = "cr", layout = NULL, nms = NULL, ggplot = FALSE, ...)
```

## Arguments

| | |
|---|---|
| `data` | A reaction time dataset. Must be a dataframe with $rt, $condition and $response. |
| `what` | @param what What to plot. Can be 'cr' for 'condition-response pairs, 'c' for condition, and 'r' for response. |
| `layout` | An optional layout. |
| `nms` | An optional vector of names for each plot. If omitted the names will be based on the contents of `data$condition` and/or `data$response`. |
| `ggplot` | ggplot Logical, should ggplot2 be used instead of base R graphics? If set to TRUE, some arguments from `linesArgs` and `...` will be ignored (but can be added to plots manually). |
| `...` | Arguments to be passed to `hist` |

## Details

This function and `rtDescriptives` are helper functions to inspect raw data.

## Value

invisible()

## Examples

```
tt = seq(0, 5, .01)
dat = simData(n = 3e4, pars = rep(.5, 5), tt = tt, pdfND = dbeta(tt, 10, 30))
rtHist(dat, breaks = tt, xlim = c(0, 1))
```

simData *Simulate data from a given density function via multinomial sampling*

## Description

Simulate data from a given density function via multinomial sampling

## Usage

```
simData(
  n,
  pars,
  tt,
  pdfND,
  fun.density = Voss.density,
  args.density = list(prec = 3),
  npars = 5,
  return.pdf = FALSE,
  normalizePdfs = TRUE
)
```

## Arguments

| | |
|---|---|
| n | Number of observations to be sampled |
| pars | Parameter values for the density function to be evaluated with. length(pars) must be a multiple of npars. |
| tt | time grid on which the density function will be evaluated. Responses not in this time grid cannot appear. |
| pdfND | either a vector of length tt specifying the nondecision density for all condition-response pairs, or a matrix where columns corresponds to the nondecision densities of condition-response pairs. Supplying NULL implies no nondecision distribution. |
| fun.density | Density function to use. |
| args.density | Additional arguments to be passed to fun.density, aside from tt, pars, and a boundary argument ('upper' or 'lower') |
| npars | Number of parameters fun.density must be evaluated with. If length(pars) > npars each npars values in pars will be seen as the parameter values of a condition. |
| return.pdf | Logical, if TRUE genData returns a list containing the probability density function used and the data, if FALSE genData returns a dataframe with simulated data. |
| normalizePdfs | Logical, should the pdf of the nondecision distribution be normalized? |

## Details

Simulate data via multinomial sampling. The response options to sample from should be provided in `tt`. The number of conditions is defined as `length(pars) / npars`.

## Value

A sorted dataframe where rows represent trials. It contains: a column named rt containing reaction times in seconds, a column named response containing either response option lower or upper, and a column named condition indicating which condition a trials belongs to. If `return.pdf` is TRUE it returns a list where the first element is the sorted dataframe, the second through the fifth elements are lists that contain densities used for simulating data.

## Examples

```
tt = seq(0, 5, .01)
pdfND = dbeta(tt, 10, 30)
n = 100
pars = c(1, 2, .5, .5, .5)
dat = simData(n, pars, tt, pdfND)
head(dat)
```

---

# testFun *Test fun.density with lower and upper bounds*

---

## Description

Test fun.density with lower and upper bounds

## Usage

```
testFun(fun.density, lower, upper, args = list())
```

## Arguments

| | |
|---|---|
| `fun.density` | A density function to be evaluated. |
| `lower` | Lower bounds of the parameter space with which `fun.density` can be evaluated. |
| `upper` | Upper bounds of the parameter space with which `fun.density` can be evaluated. |
| `args` | Additional arguments for fun.density. |

## Details

A function that is called whenever a nondefault density function is passed to `DstarM`. It does some rough error checking.

### Value

Returns TRUE if no errors occurred, otherwise returns an error message

### Examples

```
lower = c(.5, -6, .1, 0, 0)
upper = c(2, 6, .99, .99, 10)
args = list(t = seq(0, 5, .01), pars = lower, boundary = 'lower',
DstarM = TRUE)
testFun(fun.density = Voss.density, lower = lower, upper = upper,
args = args)
# TRUE
```

---

## upgradeDstarM *Upgrade a DstarM object for backwards compatibility*

---

### Description

Upgrade a DstarM object for backwards compatibility

### Usage

```
upgradeDstarM(x)
```

### Arguments

| | |
|---|---|
| x | an object of class D*M or DstarM. |

### Value

An object of class DstarM.fitD, DstarM.fitND, or DstarM.fitObs.

---

## Voss.density *Calculate model density for a given set of parameters*

---

### Description

Calculate model density for a given set of parameters

### Usage

```
Voss.density(t, pars, boundary, DstarM = TRUE, prec = 3)

LBA.density(t, pars, boundary, DstarM = TRUE, ...)

Wiener.density(t, pars, boundary, DstarM)
```

## Arguments

| | |
|---|---|
| t | Time grid for density to be calculated on. |
| pars | Parameter vector where (if DstarM == TRUE) the first index contains the boundary parameter, the second contains the drift speed, the third contains the relative starting point, the fourth contains a proportion of the maximum size of the variance on the relative starting point, the fifth contains the standard deviation of the drift speed. if DstarM == FALSE then third index of pars contains the Ter, the fifth the drift speed, the the sixth contains a proportion of the maximum size of the variance on the relative starting point, the fifth contains the standard deviation of the drift speed, and the seventh contains a proportion of the maximum variance of the Ter. |
| boundary | For which response option will the density be calculated? Either 'upper' or 'lower'. |
| DstarM | Logical, see pars. |
| prec | Precision with which the density is calculated. Corresponds roughly to the number of decimals accurately calculated. |
| ... | Other arguments, see dLBA |

## Details

These functions are examples of what fun.density should look like. Voss.density is an adaptation of ddiffusion, LBA.density is an adaptation of dLBA, and wiener.density is an adaptation of dwiener. To improve speed one can remove error handling. Normally error handling is useful, however because differential evolution can result in an incredible number of function evaluations (more than 10.000) it is recommended to omit error handling in custom density functions. estDstarM will apply some internal error checks (see testFun) on the density functions before starting differential evolution. A version of ddifusion without error handling can be found in the source code (commented out to pass R check). Note that for in Voss.density if DstarM == FALSE nondecision parameters are implemented manually and might differ from from how they are implemented in other packages. The parameter t0 specifies the mean of a uniform distribution and st0 specifies the relative size of this uniform distribution. To obtain the lower and upper range of the uniform distribution calculate a = t0 - t0*st0, and b = t0 + t0*st0.

## Value

A numeric vector of length length(t) containing a density.

## Examples

```
t = seq(0, .75, .01)
V.pars = c(1, 2, .5, .5, .5)
L.pars = c(1, .5, 2, 1, 1, 1)
W.pars = V.pars[1:3]
V1 = Voss.density(t = t, pars = V.pars, boundary = 'upper', DstarM = TRUE)
V2 = Voss.density(t = t, pars = V.pars, boundary = 'lower', DstarM = TRUE)
L1 = LBA.density(t = t, pars = L.pars, boundary = 'upper', DstarM = TRUE)
L2 = LBA.density(t = t, pars = L.pars, boundary = 'lower', DstarM = TRUE)
W1 = Wiener.density(t = t, pars = W.pars, boundary = 'upper', DstarM = TRUE)
```

```
W2 = Wiener.density(t = t, pars = W.pars, boundary = 'lower', DstarM = TRUE)
densities = cbind(V1, V2, L1, L2, W1, W2)
matplot(t, densities, type = 'b', ylab = 'Density', lty = 1, las = 1, bty = 'n',
        col = rep(1:3, each = 2), pch = c(0, 15, 1, 16, 2, 17), cex = .8,
        main = 'Model densities')
legend('topright', legend = c('Voss', 'LBA', 'RWiener'), lty = 1,
       pch = 15:17, col = 1:3, bty = 'n')
```

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