

# Package ‘dynConfIR’

October 31, 2025

**Type** Package

**Title** Dynamic Models for Confidence and Response Time Distributions

**Version** 1.1.1

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**Description** Provides density functions for the joint distribution of choice, response time and confidence for discrete confidence judgments as well as functions for parameter fitting, prediction and simulation for various dynamical models of decision confidence. All models are explained in detail by Hellmann et al. (2023; Preprint available at <<https://osf.io/9jfqr/>>, published version: <[doi:10.1037/rev0000411](https://doi.org/10.1037/rev0000411)>). Implemented models are the dynaViTE model, dynWEV model, the 2DSD model (Pleskac & Busemeyer, 2010, <[doi:10.1037/a0019737](https://doi.org/10.1037/a0019737)>), and various race models. C++ code for dynWEV and 2DSD is based on the 'rtdists' package by Henrik Singmann.

**License** GPL (>= 3)

**URL** <https://github.com/SeHellmann/dynConfIR>,  
<https://sehellmann.github.io/dynConfIR/>

**BugReports** <https://github.com/SeHellmann/dynConfIR/issues>

**Depends** R (>= 4.0)

**Imports** dplyr, magrittr, minqa, parallel, progress, Rcpp, rlang, stats

**Suggests** covr, ggplot2, MASS, Hmisc, knitr, logger, rmarkdown,  
testthat (>= 3.0.0), tidy

**LinkingTo** Rcpp

**VignetteBuilder** knitr

**Config/testthat/edition** 3

**Encoding** UTF-8

**LazyData** true

**NeedsCompilation** yes

**Repository** CRAN

**RoxygenNote** 7.3.3

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dynConfiR-package	<i>The dynConfiR Package</i>
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Description

Dynamic Models for Confidence and Response Time Distributions

Details

Package: dynConfiR  
 Type: Package  
 Version: 0.1.0  
 Date: 2023-06-19  
 Depends: R ( $\geq 4.0$ )  
 License: GPL ( $\geq 3$ )  
 URL: <https://github.com/SeHellmann/dynConfiR>

Provides response time and confidence distributions (density/PDF) for following models: dynaViTE, dynWEV, 2DSD, 2DSDT, IRM and PCRM

### Author(s)

Sebastian Hellmann

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ConfidenceOrientation *Confidence and response time data*

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### Description

A data set containing results from an orientation discrimination experiment with confidence judgments. The data set includes results from 16 participants and 3 sessions. The task was to identify the orientation (horizontal or vertical) of a grid that was briefly visible and then covered by a mask in form of a checkerboard pattern.

### Usage

```
data(ConfidenceOrientation)
```

### Format

A data frame with 25920 rows and 12 variables:

**participant** integer values as unique participant identifier

**session** session identifier ranging from 1 to 3

**gender** gender of the participant: "w" for female; "m" for male participants

**age** the age of participants in years

**SOA** stimulus-onset-asynchrony in ms (i.e. time between stimulus and mask onset)

**orientation** orientation of the target stimulus (0: vertical, 90: horizontal)

**stimulus** stimulus identity ("senkrecht": vertical, "waagrecht": horizontal)

**response** response for the discrimination task (see stimulus column)

**correct** 0-1 column indicating whether the discrimination response was correct (1) or not (0)

**rt** response time for the discrimination response in sec

**cont\_rating** confidence rating as registered (continuous values ranging from -1 (unsure) to 1 (sure))

**disc\_rating** confidence rating discretized in 5 steps using equidistant breaks

**Source**

<https://github.com/SeHellmann/SeqSamplingConfidenceModels>

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d2DSD

*Pleskac and Busemeyer's 2DSD Model for Decision Confidence*


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**Description**

Likelihood function and random number generator for a generalization of the 2DSD Model presented by Pleskac & Busemeyer (2010). It includes following parameters: DDM parameters:  $a$  (threshold separation),  $z$  (starting point; relative),  $v$  (drift rate),  $t_0$  (non-decision time/ response time constant),  $d$  (differences in speed of response execution),  $sv$  (inter-trial-variability of drift),  $st_0$  (inter-trial-variability of non-decisional components),  $sz$  (inter-trial-variability of relative starting point),  $s$  (diffusion constant).

**Usage**

```
d2DSD(rt, response = "upper", th1, th2, a, v, t0 = 0, z = 0.5, d = 0,
      sz = 0, sv = 0, st0 = 0, tau = 1, lambda = 0, s = 1,
      simult_conf = FALSE, precision = 6, z_absolute = FALSE,
      stop_on_error = TRUE, stop_on_zero = FALSE)

r2DSD(n, a, v, t0 = 0, z = 0.5, d = 0, sz = 0, sv = 0, st0 = 0,
      tau = 1, lambda = 0, s = 1, delta = 0.01, maxrt = 15,
      simult_conf = FALSE, z_absolute = FALSE, stop_on_error = TRUE)
```

**Arguments**

<code>rt</code>	a vector of RTs. Or for convenience also a <code>data.frame</code> with columns <code>rt</code> and <code>response</code> .
<code>response</code>	character vector, indicating the decision, i.e. which boundary was met first. Possible values are <code>c("upper", "lower")</code> (possibly abbreviated) and <code>"upper"</code> being the default. Alternatively, a numeric vector with values 1=lower and 2=upper or -1=lower and 1=upper, respectively. For convenience, <code>response</code> is converted via <code>as.numeric</code> also allowing factors. Ignored if the first argument is a <code>data.frame</code> .
<code>th1</code>	together with <code>th2</code> : scalars or numerical vectors giving the lower and upper bound of the interval, in which the accumulator should end at the time of the confidence judgment (i.e. at time <code>rt+tau</code> ). Only values with <code>th2&gt;=th1</code> are accepted.
<code>th2</code>	(see <code>th1</code> )
<code>a</code>	threshold separation. Amount of information that is considered for a decision. Large values indicate a conservative decisional style. Typical range: $0.5 < a < 2$

<code>v</code>	drift rate. Average slope of the information accumulation process. The drift gives information about the speed and direction of the accumulation of information. Large (absolute) values of drift indicate a good performance. If received information supports the response linked to the upper threshold the sign will be positive and vice versa. Typical range: $-5 < v < 5$
<code>t0</code>	non-decision time or response time constant (in seconds). Lower bound for the duration of all non-decisional processes (encoding and response execution). Typical range: $0.1 < t0 < 0.5$ . Default is 0.
<code>z</code>	(by default relative) starting point. Indicator of an a priori bias in decision making. When the relative starting point <code>z</code> deviates from 0.5, the amount of information necessary for a decision differs between response alternatives. Default is 0.5 (i.e., no bias).
<code>d</code>	differences in speed of response execution (in seconds). Positive values indicate that response execution is faster for responses linked to the upper threshold than for responses linked to the lower threshold. Typical range: $-0.1 < d < 0.1$ . Default is 0.
<code>sz</code>	inter-trial-variability of starting point. Range of a uniform distribution with mean <code>z</code> describing the distribution of actual starting points from specific trials. Values different from 0 can predict fast errors (but can slow computation considerably). Typical range: $0 < sz < 0.2$ . Default is 0. (Given in relative range i.e. bounded by $2 * \min(z, 1-z)$ )
<code>sv</code>	inter-trial-variability of drift rate. Standard deviation of a normal distribution with mean <code>v</code> describing the distribution of actual drift rates from specific trials. Values different from 0 can predict slow errors. Typical range: $0 < sv < 2$ . Default is 0.
<code>st0</code>	inter-trial-variability of non-decisional components. Range of a uniform distribution with mean $t0 + st0/2$ describing the distribution of actual <code>t0</code> values across trials. Accounts for response times below <code>t0</code> . Reduces skew of predicted RT distributions. Values different from 0 can slow computation considerably. Typical range: $0 < st0 < 0.2$ . Default is 0.
<code>tau</code>	post-decisional accumulation time. The length of the time period after the decision was made until the confidence judgment is made. Range: $tau > 0$ . Default: $tau=1$ .
<code>lambda</code>	power for judgment time in the division of the confidence measure by the judgment time (Default: 0, i.e. no division which is the version of 2DSD proposed by Pleskac and Busemeyer)
<code>s</code>	diffusion constant. Standard deviation of the random noise of the diffusion process (i.e., within-trial variability), scales <code>a</code> , <code>v</code> , <code>sv</code> , and <code>th</code> 's. Needs to be fixed to a constant in most applications. Default is 1. Note that the default used by Ratcliff and in other applications is often 0.1.
<code>simult_conf</code>	logical. Whether in the experiment confidence was reported simultaneously with the decision, as then decision and confidence judgment are assumed to have happened subsequent before response and computations are different, when there is an observable interjudgment time (then <code>simult_conf</code> should be FALSE).

precision	numerical scalar value. Precision of calculation. Determines the the stepsize of integration w.r.t. $z$ and $t_0$ . Represents the number of decimals precisely computed on average. Default is 6.
z_absolute	logical. Determines whether $z$ is treated as absolute start point (TRUE) or relative (FALSE; default) to $a$ .
stop_on_error	Should the diffusion functions return 0 if the parameters values are outside the allowed range (= FALSE) or produce an error in this case (= TRUE).
stop_on_zero	Should the computation of densities stop as soon as a density value of 0 occurs. This may save a lot of time if the function is used for a likelihood function. Default: FALSE
n	integer. The number of samples generated.
delta	numeric. Discretization step size for simulations in the stochastic process
maxrt	numeric. Maximum decision time returned. If the simulation of the stochastic process exceeds a decision time of <code>maxrt</code> , the response will be set to 0 and the <code>maxrt</code> will be returned as <code>rt</code> .

## Details

For confidence:  $\tau$  (post-decisional accumulation time),  $\lambda$  the exponent of judgment time for the division by judgment time in the confidence measure,  $th1$  and  $th2$  (lower and upper thresholds for confidence interval).

**Note that the parameterization or defaults of non-decision time variability  $st_0$  and diffusion constant  $s$  differ from what is often found in the literature.**

The drift diffusion model (DDM; Ratcliff and McKoon, 2008) is a mathematical model for two-choice discrimination tasks. It is based on the assumption that information is accumulated continuously until one of two decision thresholds is hit. For introduction see Ratcliff and McKoon (2008).

The 2DSD is an extension of the DDM to explain confidence judgments based on the preceding decision. It assumes a post decisional period where the process continues the accumulation of information. At the end of the period a confidence judgment (i.e. a judgment of the probability that the decision was correct) is made based on the state of the process. Here, we use a given interval, given by  $th1$  and  $th2$ , assuming that the data is given with discrete judgments and pre-processed, s.t. these discrete ratings are translated to the respective intervals. The 2DSD Model was proposed by Pleskac and Busemeyer (2010).

All functions are fully vectorized across all parameters as well as the response to match the length or `rt` (i.e., the output is always of length equal to `rt`). This allows for trial wise parameters for each model parameter.

For convenience, the function allows that the first argument is a `data.frame` containing the information of the first and second argument in two columns (i.e., `rt` and `response`). Other columns (as well as passing response separately argument) will be ignored.

## Value

`d2DSD` gives the density/likelihood/probability of the diffusion process producing a decision of response at time `rt` and a confidence judgment corresponding to the interval  $[th1, th2]$ . The value will be a numeric vector of the same length as `rt`.

r2DSD returns a `data.frame` with three columns and `n` rows. Column names are `rt` (response time), `response` (-1 (lower) or 1 (upper), indicating which bound was hit), and `conf` (the value of the confidence measure; not discretized!).

The distribution parameters (as well as `response`, `tau`, `th1` and `th2`) are recycled to the length of the result. In other words, the functions are completely vectorized for all parameters and even the response boundary.

## Note

The parameterization of the non-decisional components, `t0` and `st0`, differs from the parameterization sometimes used in the literature. In the present case `t0` is the lower bound of the uniform distribution of length `st0`, but *not* its midpoint. The parameterization employed here is in line with the functions in the `rtdists` package.

The default diffusion constant `s` is 1 and not 0.1 as in most applications of Roger Ratcliff and others. Usually `s` is not specified as the other parameters: `a`, `v`, and `sv`, may be scaled to produce the same distributions (as is done in the code).

The function code is basically an extension of the `ddiffusion` function from the package `rtdists` for the Ratcliff diffusion model.

## Author(s)

For the original `rtdists` package: Underlying C code by Jochen Voss and Andreas Voss. Porting and R wrapping by Matthew Gretton, Andrew Heathcote, Scott Brown, and Henrik Singmann. `qdiffusion` by Henrik Singmann. For the `d2DSD` function the C code was extended by Sebastian Hellmann.

## References

- Pleskac, T. J., & Busemeyer, J. R. (2010). Two-Stage Dynamic Signal Detection: A Theory of Choice, Decision Time, and Confidence, *Psychological Review*, 117(3), 864-901. doi:10.1037/a0019737
- Ratcliff, R., & McKoon, G. (2008). The diffusion decision model: Theory and data for two-choice decision tasks. *Neural Computation*, 20(4), 873-922.

## Examples

```
# Plot rt distribution ignoring confidence
curve(d2DSD(x, "upper", -Inf, Inf, tau=1, a=2, v=0.4, sz=0.2, sv=0.9), xlim=c(0, 2), lty=2)
curve(d2DSD(x, "lower", -Inf, Inf, tau=1, a=2, v=0.4, sz=0.2, sv=0.9), col="red", lty=2, add=TRUE)
curve(d2DSD(x, "upper", -Inf, Inf, tau=1, a=2, v=0.4), add=TRUE)
curve(d2DSD(x, "lower", -Inf, Inf, tau=1, a=2, v=0.4), col="red", add=TRUE)
# Generate a random sample
dfu <- r2DSD(5000, a=2, v=0.5, t0=0, z=0.5, d=0, sz=0, sv=0, st0=0, tau=1, s=1)
# Same RT distribution but upper and lower responses changed
df1 <- r2DSD(50, a=2, v=-0.5, t0=0, z=0.5, d=0, sz=0, sv=0, st0=0, tau=1, s=1)
head(dfu)

d2DSD(dfu, th1=-Inf, th2=Inf, a=2, v=.5)[1:5]
# Scaling diffusion parameters leads do same density values
s <- 2
```

```

d2DSD(dfu, th1=-Inf, th2=Inf, a=2*s, v=.5*s, s=2)[1:5]
if (requireNamespace("ggplot2", quietly = TRUE)) {
  require(ggplot2)
  ggplot(dfu, aes(x=rt, y=conf))+
    stat_density_2d(aes(fill = after_stat(density)), geom = "raster", contour = FALSE, na.rm=TRUE) +
    facet_wrap(~response)
}
boxplot(conf~response, data=dfu)

# Restricting to specific confidence region
dfu <- dfu[dfu$conf >0 & dfu$conf <1,]
d2DSD(dfu, th1=0, th2=1, a=2, v=0.5)[1:5]

# If lower confidence threshold is higher than the upper, the function throws an error,
# except when stop_on_error is FALSE
d2DSD(dfu[1:5,], th1=1, th2=0, a=2, v=0.5, stop_on_error = FALSE)

```

---

dDDConf

*Drift Diffusion Model with time-dependent confidence*


---

## Description

Likelihood function and random number generator for the Drift Diffusion Model with confidence computed as decision time. It includes following parameters: DDM parameters: a (threshold separation), z (starting point; relative), v (drift rate), t0 (non-decision time/ response time constant), d (differences in speed of response execution), sv (inter-trial-variability of drift), st0 (inter-trial-variability of non-decision components), sz (inter-trial-variability of relative starting point), s (diffusion constant).

## Usage

```

dDDConf(rt, response = "upper", th1, th2, a, v, t0 = 0, z = 0.5, d = 0,
  sz = 0, sv = 0, st0 = 1, s = 1, precision = 3,
  z_absolute = FALSE, stop_on_error = TRUE, stop_on_zero = FALSE,
  st0stepsize = 0.001)

rDDConf(n, a, v, t0 = 0, z = 0.5, d = 0, sz = 0, sv = 0, st0 = 2,
  s = 1, delta = 0.01, maxrt = 15, z_absolute = FALSE,
  stop_on_error = TRUE)

```

## Arguments

rt	a vector of RTs. Or for convenience also a data.frame with columns rt and response.
response	character vector, indicating the decision, i.e. which boundary was met first. Possible values are c("upper", "lower") (possibly abbreviated) and "upper"

being the default. Alternatively, a numeric vector with values 1=lower and 2=upper or -1=lower and 1=upper, respectively. For convenience, response is converted via `as.numeric` also allowing factors. Ignored if the first argument is a `data.frame`.

th1	together with th2: scalars or numerical vectors giving the lower and upper bound of the interval, in which the accumulator should end at the time of the confidence judgment (i.e. at time $rt + \tau$ ). Only values with $th2 \geq th1$ are accepted.
th2	(see th1)
a	threshold separation. Amount of information that is considered for a decision. Large values indicate a conservative decisional style. Typical range: $0.5 < a < 2$
v	drift rate. Average slope of the information accumulation process. The drift gives information about the speed and direction of the accumulation of information. Large (absolute) values of drift indicate a good performance. If received information supports the response linked to the upper threshold the sign will be positive and vice versa. Typical range: $-5 < v < 5$
t0	non-decision time or response time constant (in seconds). Lower bound for the duration of all non-decisional processes (encoding and response execution). Typical range: $0.1 < t_0 < 0.5$ . Default is 0.
z	(by default relative) starting point. Indicator of an a priori bias in decision making. When the relative starting point $z$ deviates from 0.5, the amount of information necessary for a decision differs between response alternatives. Default is 0.5 (i.e., no bias).
d	differences in speed of response execution (in seconds). Positive values indicate that response execution is faster for responses linked to the upper threshold than for responses linked to the lower threshold. Typical range: $-0.1 < d < 0.1$ . Default is 0.
sz	inter-trial-variability of starting point. Range of a uniform distribution with mean $z$ describing the distribution of actual starting points from specific trials. Values different from 0 can predict fast errors (but can slow computation considerably). Typical range: $0 < sz < 0.2$ . Default is 0. (Given in relative range i.e. bounded by $2 * \min(z, 1 - z)$ )
sv	inter-trial-variability of drift rate. Standard deviation of a normal distribution with mean $v$ describing the distribution of actual drift rates from specific trials. Values different from 0 can predict slow errors. Typical range: $0 < sv < 2$ . Default is 0.
st0	inter-trial-variability of non-decisional components. Range of a uniform distribution with mean $t_0 + st_0/2$ describing the distribution of actual $t_0$ values across trials. Accounts for response times below $t_0$ . Reduces skew of predicted RT distributions. Values different from 0 can slow computation considerably. Typical range: $0 < st_0 < 0.2$ . Default is 0.
s	diffusion constant. Standard deviation of the random noise of the diffusion process (i.e., within-trial variability), scales $a$ , $v$ , $sv$ , and $th$ 's. Needs to be fixed to a constant in most applications. Default is 1. Note that the default used by Ratcliff and in other applications is often 0.1.

precision	numerical scalar value. Precision of calculation. Corresponds to the stepsize of integration w.r.t. $z$ . Default is $1e-5$ .
z_absolute	logical. Determines whether $z$ is treated as absolute start point (TRUE) or relative (FALSE; default) to $a$ .
stop_on_error	Should the diffusion functions return 0 if the parameters values are outside the allowed range (= FALSE) or produce an error in this case (= TRUE).
stop_on_zero	Should the computation of densities stop as soon as a density value of 0 occurs. This may save a lot of time if the function is used for a likelihood function. Default: FALSE
st0stepsize	numerical scalar value. Stepsize for integration over $t0$ .
n	integer. The number of samples generated.
delta	numeric. Discretization step size for simulations in the stochastic process
maxrt	numeric. Maximum decision time returned. If the simulation of the stochastic process exceeds a decision time of <code>maxrt</code> , the response will be set to 0 and the <code>maxrt</code> will be returned as <code>rt</code> .

## Details

For the confidence part: `th1` and `th2` (lower and upper thresholds for decision time interval).

**Note that the parameterization or defaults of non-decision time variability `st0` and diffusion constant `s` differ from what is often found in the literature.**

The Ratcliff diffusion model (Ratcliff and McKoon, 2008) is a mathematical model for two-choice discrimination tasks. It is based on the assumption that information is accumulated continuously until one of two decision thresholds is hit. For introduction see Ratcliff and McKoon (2008).

This model incorporates the idea, that the decision time  $T$  is informative for stimulus difficulty and thus confidence is computed as a monotone function of  $\frac{1}{\sqrt{T}}$ . In this implementation, confidence is the decision time, directly. Here, we use an interval, given by `th1` and `th2`, assuming that the data is given with discrete judgments and pre-processed, s.t. these discrete ratings are translated to the respective intervals.

All functions are fully vectorized across all parameters as well as the response to match the length or `rt` (i.e., the output is always of length equal to `rt`). This allows for trial wise parameters for each model parameter.

For convenience, the function allows that the first argument is a `data.frame` containing the information of the first and second argument in two columns (i.e., `rt` and `response`). Other columns (as well as passing response separately argument) will be ignored.

## Value

`dDDConf` gives the density/likelihood/probability of the diffusion process producing a decision of response at time `rt` and a confidence judgment corresponding to the interval  $[th1, th2]$ . The value will be a numeric vector of the same length as `rt`.

`rDDConf` returns a `data.frame` with three columns and `n` rows. Column names are `rt` (response time), `response` (-1 (lower) or 1 (upper), indicating which bound was hit), `conf` for the decision time (without non-decision time component; not discretized!).

The distribution parameters (as well as response, th1 and th2) are recycled to the length of the result. In other words, the functions are completely vectorized for all parameters and even the response boundary.

### Note

The parameterization of the non-decisional components,  $t_0$  and  $st_0$ , differs from the parameterization sometimes used in the literature. In the present case  $t_0$  is the lower bound of the uniform distribution of length  $st_0$ , but *not* its midpoint. The parameterization employed here is in line with the functions in the `rtdists` package.

The default diffusion constant  $s$  is 1 and not 0.1 as in most applications of Roger Ratcliff and others. Usually  $s$  is not specified as the other parameters:  $a$ ,  $v$ , and  $sv$ , may be scaled to produce the same distributions (as is done in the code).

The function code is basically an extension of the `ddiffusion` function from the package `rtdists` for the Ratcliff diffusion model.

### Author(s)

For the original `rtdists` package: Underlying C code by Jochen Voss and Andreas Voss. Porting and R wrapping by Matthew Gretton, Andrew Heathcote, Scott Brown, and Henrik Singmann. `qdiffusion` by Henrik Singmann. For the `dDDConf` function the C code was extended by Sebastian Hellmann.

### References

- Ratcliff, R., & McKoon, G. (2008). The diffusion decision model: Theory and data for two-choice decision tasks. *Neural Computation*, 20(4), 873-922.
- Hellmann, S., Zehetleitner, M., & Rausch, M. (2023). Simultaneous modeling of choice, confidence and response time in visual perception. *Psychological Review* 2023 Mar 13. doi: 10.1037/rev0000411. Epub ahead of print. PMID: 36913292.

### Examples

```
# Plot rt distribution ignoring confidence
curve(dDDConf(x, "upper", 0, Inf, a=2, v=0.4, sz=0.1, sv=0.9), xlim=c(0, 2), lty=2, n=20)
curve(dDDConf(x, "lower", 0, Inf, a=2, v=0.4, sz=0.1, sv=0.9), col="red", lty=2, add=TRUE, n=20)
curve(dDDConf(x, "upper", 0, Inf, a=2, v=0.4), add=TRUE, n=20)
curve(dDDConf(x, "lower", 0, Inf, a=2, v=0.4), col="red", add=TRUE, n=20)
# Generate a random sample
dfu <- rDDConf(1500, a=2, v=0.5, t0=0, z=0.5, d=0, sz=0, sv=0, st0=0.2, s=1)
# Same RT distribution but upper and lower responses changed
df1 <- rDDConf(50, a=2, v=-0.5, t0=0, z=0.5, d=0, sz=0, sv=0, st0=0.2, s=1)
head(dfu)

dDDConf(dfu, th1=0.5, th2=2.5, a=2, v=.5, st0=0.2)[1:5]
# Scaling diffusion parameters leads do same density values
s <- 2
dDDConf(dfu, th1=0.5, th2=2.5, a=2*s, v=.5*s, s=2, st0=0.2)[1:5]
if (requireNamespace("ggplot2", quietly = TRUE)) {
  require(ggplot2)
```

```

ggplot(dfu, aes(x=rt, y=conf))+
  stat_density_2d(aes(fill = after_stat(density)), geom = "raster", contour = FALSE, na.rm=TRUE) +
  facet_wrap(~response)
}
boxplot(conf~response, data=dfu)

# Restricting to specific confidence region
dfu <- dfu[dfu$conf >0 & dfu$conf <1,]
dDDConf(dfu, th1=0, th2=1, a=2, v=0.5, st0=0.1)[1:5]

# If lower confidence threshold is higher than the upper, the function throws an error,
# except when stop_on_error is FALSE
dDDConf(dfu[1:5,], th1=1, th2=0, a=2, v=0.5, stop_on_error = FALSE)

```

---

dynaViTE

---

*Dynamical visibility, time, and evidence model (dynaViTE) and Dynamical weighted evidence and visibility model (dynWEV)*


---

## Description

Likelihood function and random number generator for the dynaViTE and dynWEV model (Hellmann et al., 2023). It includes following parameters from the drift diffusion model:  $a$  (threshold separation),  $z$  (starting point; relative),  $v$  (drift rate),  $t_0$  (non-decision time/response time constant),  $d$  (differences in speed of response execution),  $sv$  (inter-trial-variability of drift),  $st_0$  (inter-trial-variability of non-decisional components),  $sz$  (inter-trial-variability of relative starting point) and  $s$  (diffusion constant). For the computation of confidence following parameters were added:  $\tau$  (post-decisional accumulation time),  $w$  (weight on the decision evidence (weight on visibility is  $(1-w)$ )),  $\mu_{vis}$  (mean drift rate of visibility process),  $s_{vis}$  (diffusion constant of visibility process),  $\sigma_{vis}$  (variability in drift rate of visibility accumulator),  $th_1$  and  $th_2$  (lower and upper thresholds for confidence interval).  $\lambda$  for dynaViTE only, the exponent of judgment time for the division by judgment time in the confidence measure, and **Note that the parametrization or defaults of non-decision time variability  $st_0$  and diffusion constant  $s$  differ from what is often found in the literature.**

Likelihood function and random number generator for the dynaViTE and dynWEV model (Hellmann et al., 2023). It includes following parameters from the drift diffusion model:  $a$  (threshold separation),  $z$  (starting point; relative),  $v$  (drift rate),  $t_0$  (non-decision time/response time constant),  $d$  (differences in speed of response execution),  $sv$  (inter-trial-variability of drift),  $st_0$  (inter-trial-variability of non-decisional components),  $sz$  (inter-trial-variability of relative starting point) and  $s$  (diffusion constant). For the computation of confidence following parameters were added:  $\tau$  (post-decisional accumulation time),  $w$  (weight on the decision evidence (weight on visibility is  $(1-w)$ )),  $\mu_{vis}$  (mean drift rate of visibility process),  $s_{vis}$  (diffusion constant of visibility process),  $\sigma_{vis}$  (variability in drift rate of visibility accumulator),  $th_1$  and  $th_2$  (lower and upper thresholds for confidence interval).  $\lambda$  for dynaViTE only, the exponent of judgment time for the division by judgment time in the confidence measure, and **Note that the parametrization or defaults of non-decision time variability  $st_0$  and diffusion constant  $s$  differ from what is often found in the literature.**

**Usage**

```
dWEV(rt, response = "upper", th1, th2, a, v, t0 = 0, z = 0.5, d = 0,
      sz = 0, sv = 0, st0 = 0, tau = 1, w = 0.5, muvis = NULL,
      sigvis = 0, svis = 1, lambda = 0, s = 1, simult_conf = FALSE,
      precision = 6, z_absolute = FALSE, stop_on_error = TRUE,
      stop_on_zero = FALSE)
```

```
ddynaViTE(rt, response = "upper", th1, th2, a, v, t0 = 0, z = 0.5,
           d = 0, sz = 0, sv = 0, st0 = 0, tau = 1, w = 0.5, muvis = NULL,
           sigvis = 0, svis = 1, lambda = 0, s = 1, simult_conf = FALSE,
           precision = 6, z_absolute = FALSE, stop_on_error = TRUE,
           stop_on_zero = FALSE)
```

```
rdynaViTE(n, a, v, t0 = 0, z = 0.5, d = 0, sz = 0, sv = 0, st0 = 0,
           tau = 1, w = 0.5, muvis = NULL, sigvis = 0, svis = 1, lambda = 0,
           s = 1, delta = 0.01, maxrt = 15, simult_conf = FALSE,
           z_absolute = FALSE, stop_on_error = TRUE, process_results = FALSE)
```

```
dWEV(rt, response = "upper", th1, th2, a, v, t0 = 0, z = 0.5, d = 0,
      sz = 0, sv = 0, st0 = 0, tau = 1, w = 0.5, muvis = NULL,
      sigvis = 0, svis = 1, lambda = 0, s = 1, simult_conf = FALSE,
      precision = 6, z_absolute = FALSE, stop_on_error = TRUE,
      stop_on_zero = FALSE)
```

```
rWEV(n, a, v, t0 = 0, z = 0.5, d = 0, sz = 0, sv = 0, st0 = 0,
      tau = 1, w = 0.5, muvis = NULL, sigvis = 0, svis = 1, lambda = 0,
      s = 1, delta = 0.01, maxrt = 15, simult_conf = FALSE,
      z_absolute = FALSE, stop_on_error = TRUE, process_results = FALSE)
```

```
ddynaViTE(rt, response = "upper", th1, th2, a, v, t0 = 0, z = 0.5,
           d = 0, sz = 0, sv = 0, st0 = 0, tau = 1, w = 0.5, muvis = NULL,
           sigvis = 0, svis = 1, lambda = 0, s = 1, simult_conf = FALSE,
           precision = 6, z_absolute = FALSE, stop_on_error = TRUE,
           stop_on_zero = FALSE)
```

```
rdynaViTE(n, a, v, t0 = 0, z = 0.5, d = 0, sz = 0, sv = 0, st0 = 0,
           tau = 1, w = 0.5, muvis = NULL, sigvis = 0, svis = 1, lambda = 0,
           s = 1, delta = 0.01, maxrt = 15, simult_conf = FALSE,
           z_absolute = FALSE, stop_on_error = TRUE, process_results = FALSE)
```

**Arguments**

<code>rt</code>	a vector of RTs. Or for convenience also a data.frame with columns <code>rt</code> and <code>response</code> .
<code>response</code>	character vector, indicating the decision, i.e. which boundary was met first. Possible values are <code>c("upper", "lower")</code> (possibly abbreviated) and <code>"upper"</code> being the default. Alternatively, a numeric vector with values <code>1=lower</code> and <code>2=upper</code> or <code>-1=lower</code> and <code>1=upper</code> , respectively. For convenience, <code>response</code> is con-

	<p>verted via <code>as.numeric</code> also allowing factors. Ignored if the first argument is a <code>data.frame</code>.</p>
th1	<p>together with th2: scalars or numerical vectors giving the lower and upper bound of the interval of the confidence measure (see Details). Only values with <math>th2 \geq th1</math> are accepted.</p>
th2	<p>(see th1)</p>
a	<p>threshold separation. Amount of information that is considered for a decision. Large values indicate a conservative decisional style. Typical range: <math>0.5 &lt; a &lt; 2</math></p>
v	<p>drift rate of decision process. Average slope of the information accumulation process. The drift gives information about the speed and direction of the accumulation of information. Large (absolute) values of drift indicate a good performance. If received information supports the response linked to the upper threshold the sign will be positive and vice versa. Typical range: <math>-5 &lt; v &lt; 5</math></p>
t0	<p>non-decision time or response time constant (in seconds). Lower bound for the duration of all non-decisional processes (encoding and response execution). Typical range: <math>0.1 &lt; t0 &lt; 0.5</math>. Default is 0.</p>
z	<p>(by default relative) starting point of decision process. Indicator of an a priori bias in decision making. When the relative starting point <math>z</math> deviates from 0.5, the amount of information necessary for a decision differs between response alternatives. Default is 0.5 (i.e., no bias).</p>
d	<p>differences in speed of response execution (in seconds). Positive values indicate that response execution is faster for responses linked to the upper threshold than for responses linked to the lower threshold. Typical range: <math>-0.1 &lt; d &lt; 0.1</math>. Default is 0.</p>
sz	<p>inter-trial-variability of starting point. Range of a uniform distribution with mean <math>z</math> describing the distribution of actual starting points from specific trials. Values different from 0 can predict fast errors (but can slow computation considerably). Typical range: <math>0 &lt; sz &lt; 0.2</math>. Default is 0. (Given in relative range i.e. bounded by <math>2 * \min(z, 1-z)</math>)</p>
sv	<p>inter-trial-variability of drift rate of decision process. Standard deviation of a normal distribution with mean <math>v</math> describing the distribution of actual drift rates from specific trials. Values different from 0 can predict slow errors. Typical range: <math>0 &lt; sv &lt; 2</math>. Default is 0.</p>
st0	<p>inter-trial-variability of non-decisional components. Range of a uniform distribution with mean <math>t0 + st0/2</math> describing the distribution of actual <math>t0</math> values across trials. Accounts for response times below <math>t0</math>. Reduces skew of predicted RT distributions. Values different from 0 can slow computation considerably. Typical range: <math>0 &lt; st0 &lt; 0.2</math>. Default is 0.</p>
tau	<p>post-decisional accumulation time; the length of the time period after the decision was made until the confidence judgment is made. Range: <math>tau &gt; 0</math>. Default: <math>tau=1</math>.</p>
w	<p>weight put on the final state of the decision accumulator for confidence computation. <math>1-w</math> is the weight on the visibility accumulator. Range: <math>0 &lt; w &lt; 1</math>. Default: <math>w=0.5</math>.</p>

<code>muvis</code>	mean drift of visibility process. If NULL (default), <code>muvis</code> will be set to the absolute value of <code>v</code> .
<code>sigvis</code>	the variability in drift rate of the visibility process (which varies independently from the drift rate in decision process). Range: <code>sigvis</code> ≥ 0. Default: <code>sigvis</code> = 0.
<code>svis</code>	diffusion constant of visibility process. Range: <code>svis</code> > 0. Default: <code>svis</code> = 1.
<code>lambda</code>	power for judgment time in the division of the confidence measure by the judgment time (Default: 0, i.e. no division which is the version of <code>dynWEV</code> proposed by Hellmann et al., 2023)
<code>s</code>	diffusion constant of decision process; standard deviation of the random noise of the diffusion process (i.e., within-trial variability), scales other parameters (see Note). Needs to be fixed to a constant in most applications. Default is 1. Note that the default used by Ratcliff and in other applications is often 0.1.
<code>simult_conf</code>	logical. Whether in the experiment confidence was reported simultaneously with the decision. If that is the case decision and confidence judgment are assumed to have happened subsequent before the response. Therefore <code>tau</code> is included in the response time. If the decision was reported before the confidence report, <code>simul_conf</code> should be FALSE.
<code>precision</code>	numerical scalar value. Precision of calculation. Determines the step size of integration w.r.t. <code>z</code> and <code>t0</code> . Represents the number of decimals precisely computed on average. Default is 6.
<code>z_absolute</code>	logical. Determines whether <code>z</code> is treated as absolute start point (TRUE) or relative (FALSE; default) to <code>a</code> .
<code>stop_on_error</code>	Should the diffusion functions return 0 if the parameters values are outside the allowed range (= FALSE) or produce an error in this case (= TRUE).
<code>stop_on_zero</code>	Should the computation of densities stop as soon as a density value of 0 occurs. This may save a lot of time if the function is used for a likelihood function. Default: FALSE
<code>n</code>	integer. The number of samples generated.
<code>delta</code>	numeric. Discretization step size for simulations in the stochastic process
<code>maxrt</code>	numeric. Maximum decision time returned. If the simulation of the stochastic process exceeds a decision time of <code>maxrt</code> , the response will be set to 0 and the <code>maxrt</code> will be returned as <code>rt</code> .
<code>process_results</code>	logical. Whether the output simulations should contain the final state of the decision (and visibility) process as additional column. Default is FALSE, meaning that no additional columns for the final process states are returned.

## Details

The function `dWEV` was renamed to `ddynaViTE` in version 0.1.0 of the package. It is still here for reasons of backwards compatibility. The function just calls the `ddynaViTE` function (and produces a deprecation warning).

The dynamical visibility, time, and evidence (dynaViTE) model and the weighted evidence and visibility model (dynWEV) are extensions of the 2DSD model for decision confidence (see [d2DSD](#)). It assumes that the decision follows a drift diffusion model with two additional assumptions to

account for confidence. First, there is a post-decisional period of further evidence accumulation  $\tau$ . Second, another accumulation process accrues information about stimulus reliability (the visibility process) including also evidence about decision irrelevant features. See Hellmann et al. (2023) for more information. The measure for confidence is then a weighted sum of the final state of the decision process  $X$  and the visibility process  $V$  over a power-function of total accumulation time, i.e. for a decision time  $T$  (which is not the response time), the confidence variable is

$$conf = \frac{wX(T + \tau) + (1 - w)V(T + \tau)}{(T + \tau)^\lambda}.$$

The dynWEV model is a special case of dynaViTE, with the parameter  $\lambda=0$ .

All functions are fully vectorized across all parameters as well as the response to match the length or `rt` (i.e., the output is always of length equal to `rt`). This allows for trial wise parameters for each model parameter.

For convenience, the function allows that the first argument is a `data.frame` containing the information of the first and second argument in two columns (i.e., `rt` and `response`). Other columns (as well as passing response separately argument) will be ignored.

The functions `dWEV` and `rWEV` were renamed to `ddynaViTE` and `rdynaViTE`, respectively in version 0.1.0 of the package. They are still here for reasons of backwards compatibility. The functions just calls their counterpart `ddynaViTE` and `rdynaViTE` (and produce a deprecation warning).

The dynamical visibility, time, and evidence (dynaViTE) model and the weighted evidence and visibility model (dynWEV) are extensions of the 2DSD model for decision confidence (see [d2DSD](#)). It assumes that the decision follows a drift diffusion model with two additional assumptions to account for confidence. First, there is a post-decisional period of further evidence accumulation  $\tau$ . Second, another accumulation process accrues information about stimulus reliability (the visibility process) including also evidence about decision irrelevant features. See Hellmann et al. (2023) for more information. The measure for confidence is then a weighted sum of the final state of the decision process  $X$  and the visibility process  $V$  over a power-function of total accumulation time, i.e. for a decision time  $T$  (which is not the response time), the confidence variable is

$$conf = \frac{wX(T + \tau) + (1 - w)V(T + \tau)}{(T + \tau)^\lambda}.$$

The dynWEV model is a special case of dynaViTE, with the parameter  $\lambda=0$ .

All functions are fully vectorized across all parameters as well as the response to match the length or `rt` (i.e., the output is always of length equal to `rt`). This allows for trial wise parameters for each model parameter.

For convenience, the function allows that the first argument is a `data.frame` containing the information of the first and second argument in two columns (i.e., `rt` and `response`). Other columns (as well as passing response separately argument) will be ignored.

## Value

`ddynaViTE` gives the density/likelihood/probability of the diffusion process producing a decision of response at time `rt` and a confidence judgment corresponding to the interval  $[th1, th2]$ . The value will be a numeric vector of the same length as `rt`.

`rdynaViTE` returns a `data.frame` with three columns and `n` rows. Column names are `rt` (response time), `response` (-1 (lower) or 1 (upper), indicating which bound was hit), and `conf` (the value of the confidence measure; not discretized!).

The distribution parameters (as well as response, tau, th1 and th2, w and sig) are recycled to the length of the result. In other words, the functions are completely vectorized for all parameters and even the response boundary.

ddynaViTE gives the density/likelihood/probability of the diffusion process producing a decision of response at time `rt` and a confidence judgment corresponding to the interval [ `th1`, `th2`]. The value will be a numeric vector of the same length as `rt`.

rdynaViTE returns a `data.frame` with three columns and `n` rows. Column names are `rt` (response time), `response` (-1 (lower) or 1 (upper), indicating which bound was hit), and `conf` (the value of the confidence measure; not discretized!).

The distribution parameters (as well as response, tau, th1 and th2, w and sig) are recycled to the length of the result. In other words, the functions are completely vectorized for all parameters and even the response boundary.

### Note

The parameterization of the non-decisional components, `t0` and `st0`, differs from the parameterization sometimes used in the literature. In the present case `t0` is the lower bound of the uniform distribution of length `st0`, but *not* its midpoint. The parameterization employed here is in line with the functions in the `rtdists` package.

The default diffusion constant `s` is 1 and not 0.1 as in most applications of Roger Ratcliff and others. Usually `s` is not specified as the other parameters: `a`, `v`, `sv`, `muvis`, `sigvis`, and `svis` respectively, may be scaled to produce the same distributions (as is done in the code).

The function code is basically an extension of the `ddiffusion` function from the package `rtdists` for the Ratcliff diffusion model.

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The function code is basically an extension of the `ddiffusion` function from the package `rtdists` for the Ratcliff diffusion model.

### Author(s)

Sebastian Hellmann

### References

- Hellmann, S., Zehetleitner, M., & Rausch, M. (2023). Simultaneous modeling of choice, confidence and response time in visual perception. *Psychological Review* 2023 Mar 13. doi: 10.1037/rev0000411. Epub ahead of print. PMID: 36913292.
- Hellmann, S., Zehetleitner, M., & Rausch, M. (2023). Simultaneous modeling of choice, confidence and response time in visual perception. *Psychological Review* 2023 Mar 13. doi: 10.1037/rev0000411. Epub ahead of print. PMID: 36913292.

## Examples

```
# Plot rt distribution ignoring confidence
curve(ddynaViTE(x, "upper", -Inf, Inf, tau=1, a=2, v=0.4, sz=0.2, sv=0.9), xlim=c(0, 2), lty=2)
curve(ddynaViTE(x, "lower", -Inf, Inf, tau=1, a=2, v=0.4, sz=0.2, sv=0.9))
curve(ddynaViTE(x, "upper", -Inf, Inf, tau=1, a=2, v=0.4), add=TRUE)
curve(ddynaViTE(x, "lower", -Inf, Inf, tau=1, a=2, v=0.4), col="red", add=TRUE)
# Generate a random sample
df1 <- rdynaViTE(5000, a=2, v=0.5, t0=0, z=0.5, d=0, sz=0, sv=0, st0=0, tau=1, s=1, w=0.9)
# Same RT and response distribution but different confidence distribution
df2 <- rdynaViTE(5000, a=2, v=0.5, t0=0, z=0.5, d=0, sz=0, sv=0, st0=0, tau=1, s=1, w=0.1)
head(df1)

# Scaling diffusion parameters leads do same density values
ddynaViTE(df1[1:5,], th1=-Inf, th2=Inf, a=2, v=.5)[1:5]
s <- 2
ddynaViTE(df1[1:5,], th1=-Inf, th2=Inf, a=2*s, v=.5*s, s=2)[1:5]

# Diffusion constant also scales confidence parameters
ddynaViTE(df1[1:5,], th1=0.2, th2=1, a=2, v=.5, sv=0.2, w=0.5, sigvis = 0.2, svis = 1)[1:5]
s <- 2
ddynaViTE(df1[1:5,], th1=0.2*s, th2=1*s, a=2*s, v=.5*s, s=2,
          sv=0.2*s, w=0.5, sigvis=0.2*s, svis=1*s)[1:5]

two_samples <- rbind(cbind(df1, w="high"),
                    cbind(df2, w="low"))
# no difference in RT distributions
boxplot(rt~w+response, data=two_samples)
# but different confidence distributions
boxplot(conf~w+response, data=two_samples)
if (requireNamespace("ggplot2", quietly = TRUE)) {
  require(ggplot2)
  ggplot(two_samples, aes(x=rt, y=conf))+
    stat_density_2d(aes(fill = after_stat(density)), geom = "raster", contour = FALSE, na.rm=TRUE) +
    xlim(c(0, 2))+ ylim(c(-1.5, 4))+
    facet_grid(cols=vars(w), rows=vars(response), labeller = "label_both")
}

# Restricting to specific confidence region
df1 <- df1[df1$conf >0 & df1$conf <1,]
ddynaViTE(df1[1:5,], th1=0, th2=1, a=2, v=0.5)[1:5]

# If lower confidence threshold is higher than the upper, the function throws an error,
# except when stop_on_error is FALSE
ddynaViTE(df1[1:5,], th1=1, th2=0, a=2, v=0.5, stop_on_error = FALSE)

# Plot rt distribution ignoring confidence
curve(ddynaViTE(x, "upper", -Inf, Inf, tau=1, a=2, v=0.4,
              sz=0.2, sv=0.9),
      xlim=c(0, 2), lty=2)
curve(ddynaViTE(x, "lower", -Inf, Inf, tau=1, a=2, v=0.4,
              sz=0.2, sv=0.9),
```

```

      col="red", lty=2, add=TRUE)
curve(ddynaViTE(x, "upper", -Inf, Inf, tau=1, a=2, v=0.4), add=TRUE)
curve(ddynaViTE(x, "lower", -Inf, Inf, tau=1, a=2, v=0.4), add=TRUE)
# Generate a random sample
df1 <- rdynaViTE(5000, a=2, v=0.5, t0=0, z=0.5, d=0, sz=0, sv=0, st0=0, tau=1, s=1, w=0.9)
# Same RT and response distribution but different confidence distribution
df2 <- rdynaViTE(5000, a=2, v=0.5, t0=0, z=0.5, d=0, sz=0, sv=0, st0=0, tau=1, s=1, w=0.1)
head(df1)

# Scaling diffusion parameters leads do same density values
ddynaViTE(df1[1:5,], th1=-Inf, th2=Inf, a=2, v=.5)[1:5]
s <- 2
ddynaViTE(df1[1:5,], th1=-Inf, th2=Inf, a=2*s, v=.5*s, s=2)[1:5]

# Diffusion constant also scales confidence parameters
ddynaViTE(df1[1:5,], th1=0.2, th2=1, a=2, v=.5, sv=0.2, w=0.5, sigvis = 0.2, svis = 1)[1:5]
s <- 2
ddynaViTE(df1[1:5,], th1=0.2*s, th2=1*s, a=2*s, v=.5*s, s=2,
          sv=0.2*s, w=0.5, sigvis=0.2*s, svis=1*s)[1:5]

two_samples <- rbind(cbind(df1, w="high"),
                    cbind(df2, w="low"))
# no difference in RT distributions
boxplot(rt~w+response, data=two_samples)
# but different confidence distributions
boxplot(conf~w+response, data=two_samples)
if (requireNamespace("ggplot2", quietly = TRUE)) {
  require(ggplot2)
  ggplot(two_samples, aes(x=rt, y=conf))+
    stat_density_2d(aes(fill = after_stat(density)), geom = "raster", contour = FALSE, na.rm=TRUE) +
    xlim(c(0, 2))+ ylim(c(-1.5, 4))+
    facet_grid(cols=vars(w), rows=vars(response), labeller = "label_both")
}

# Restricting to specific confidence region
df1 <- df1[df1$conf >0 & df1$conf <1,]
ddynaViTE(df1[1:5,], th1=0, th2=1, a=2, v=0.5)[1:5]

# If lower confidence threshold is higher than the upper, the function throws an error,
# except when stop_on_error is FALSE
ddynaViTE(df1[1:5,], th1=1, th2=0, a=2, v=0.5, stop_on_error = FALSE)

```

---

fitRTConf

---

*Function for fitting sequential sampling confidence models*


---

## Description

Fits the parameters of different models of response time and confidence, including the 2DSD model (Pleskac & Busemeyer, 2010), dynWEV, DDConf, and various flavors of race models (Hellmann

et al., 2023). Which model to fit is specified by the argument `model`. Only a ML method is implemented. See [ddynaViTE](#), [d2DSD](#), [dRM](#), and [dMTLNR](#) for more information about the parameters and Details for not-fitted parameters.

## Usage

```
fitRTConf(data, model = "dynWEV", fixed = list(sym_thetas = FALSE),
  init_grid = NULL, grid_search = TRUE, data_names = list(),
  nRatings = NULL, restr_tau = Inf, precision = 3, logging = FALSE,
  opts = list(), optim_method = "bobyqa", useparallel = FALSE,
  n.cores = NULL, ...)
```

## Arguments

<code>data</code>	<p>a <code>data.frame</code> where each row is one trial, containing following variables (column names can be changed by passing additional arguments of the form <code>condition="contrast"</code>):</p> <ul style="list-style-type: none"> <li>• <code>condition</code> (not necessary; for different levels of stimulus quality, will be transformed to a factor),</li> <li>• <code>rating</code> (discrete confidence judgments, should be given as integer vector; otherwise will be transformed to integer),</li> <li>• <code>rt</code> (giving the reaction times for the decision task),</li> <li>• either 2 of the following (see details for more information about the accepted formats):             <ul style="list-style-type: none"> <li>– <code>stimulus</code> (encoding the stimulus category in a binary choice task),</li> <li>– <code>response</code> (encoding the decision response),</li> <li>– <code>correct</code> (encoding whether the decision was correct; values in 0, 1)</li> </ul> </li> <li>• <code>sbj</code> or <code>participant</code> (optional; giving the subject ID; only relevant if <code>logging = TRUE</code>; if unique the ID is used in saved files with interim results and logging messages; if non-unique or missing and <code>logging = TRUE</code>, 999 will be used then)</li> </ul>
<code>model</code>	character scalar. One of "dynWEV", "2DSD", "IRM", "PCRM", "IRMt", "PCRMt", "MTLNR", or "DDConf" for the model to be fit.
<code>fixed</code>	list. List with parameter-value pairs for parameters that should not be fitted. See Details.
<code>init_grid</code>	<code>data.frame</code> or <code>NULL</code> . Grid for the initial parameter search. Each row is one parameter constellation. See details for more information. If <code>NULL</code> a default grid will be used.
<code>grid_search</code>	logical. If <code>FALSE</code> , the grid search before the optimization algorithm is omitted. The fitting is then started with a mean parameter set from the default grid (if <code>init_grid=NULL</code> ) or directly with the rows from <code>init_grid</code> , if not <code>NULL</code> . (Default: <code>TRUE</code> )
<code>data_names</code>	named list (e.g. <code>c(rating="confidence")</code> ). Alternative possibility of giving other column names for the variables in the data. By default column names are identical to the ones given in the data argument description.

nRatings	integer. Number of rating categories. If NULL, the maximum of rating and length(unique(rating)) is used. This argument is especially important for data sets where not the whole range of rating categories is realized. If given, ratings has to be given as factor or integer.
restr_tau	numerical or Inf or "simult_conf". For 2DSD and dynWEV only. Upper bound for tau. Fits will be in the interval (0,restr_tau). If FALSE tau will be unbound. For "simult_conf", see the documentation of <a href="#">d2DSD</a> and <a href="#">ddynaViTE</a>
precision	numeric. Precision of calculation for the density functions (see <a href="#">ddynaViTE</a> and <a href="#">dPCRM</a> for more information).
logging	logical. If TRUE, a folder 'autosave/fitmodel' is created and messages about the process are printed in a logging file and to console (depending on OS). Additionally intermediate results are saved in a .RData file with the participant ID in the name.
opts	list. A list for more control options in the optimization routines (depending on the optim_method). See details for more information.
optim_method	character. Determines which optimization function is used for the parameter estimation. Either "bobyqa" (default), "L-BFGS-B" or "Nelder-Mead". "bobyqa" uses a box-constrained optimization with quadratic interpolation. (See <a href="#">bobyqa</a> for more information.) The first two use a box-constraint optimization. For Nelder-Mead a transfinite function rescaling is used (i.e. the constrained arguments are suitably transformed to the whole real line).
useparallel	logical. If TRUE the grid search in the beginning is done with a parallel back-end, using the parallel package.
n.cores	integer or NULL. Number of cores used for parallelization. If NULL (default) the number of available cores -1 is used.
...	Possibility of giving alternative variable names in data frame (in the form condition = "SOA", or response="pressedKey").

## Details

The fitting involves a first grid search through computation of the likelihood on an initial grid with possible sets of parameters to start the optimization routine. Then the best nAttempts parameter sets are chosen for an optimization, which is done with an algorithm, depending on the argument optim-method. The Nelder-Mead algorithm uses the R function [optim](#). The optimization routine is restarted nRestarts times with the starting parameter set equal to the best parameters from the previous routine.

**stimulus, response and correct.** Two of these columns must be given in data. If all three are given, correct will have no effect (and will be not checked!). stimulus can always be given in numerical format with values -1 and 1. response can always be given as a character vector with "lower" and "upper" as values. Correct must always be given as a 0-1-vector. If the stimulus column is given together with a response column and they both do not match the above format, they need to have the same values/levels (if factor). In the case that only stimulus/response is given in any other format together with correct, the unique values will be sorted increasingly and the first value will be encoded as "lower"/-1 and the second as "upper"/+1.

**fixed.** Parameters that should not be fitted but kept constant. These will be dropped from the initial grid search but will be present in the output, to keep all parameters for prediction in the result. Includes the possibility for symmetric confidence thresholds for both alternative (sym\_thetas=logical).

Other examples are  $z = .5$ ,  $sv=0$ ,  $st0=0$ ,  $sz=0$ . For race models, the possibility of setting  $a='b'$  (or vice versa) leads to identical upper bounds on the decision processes, which is the equivalence for  $z=.5$  in a diffusion process.

**Parameters not fitted.** The models get developed continuously and not all changes are adopted in the fitting function instantly. Following parameters are currently not included in the fitting routine:

- in race models:  $sza$ ,  $szb$ ,  $smu1$ , and  $smu2$

`init_grid`. Each row should be one parameter set to check. The column names should include the parameters of the desired model, which are the following for 2DSD:  $a$ ,  $vmin$  and  $vmax$  (will be equidistantly spanned across conditions),  $sv$ ,  $z$  (as the relative starting point between 0 and  $a$ ),  $sz$  (also in relative terms),  $t0$ ,  $st0$ ,  $theta0$  (minimal threshold),  $thetamax$  (maximal threshold; the others will be equidistantly spanned symmetrically for both decisions), and  $tau$ . For dynWEV, additionally  $w$ ,  $svis$ , and  $sigvis$  are required.

For the race models the parameters are:  $vmin$ ,  $vmax$  (will be equidistantly spanned across conditions),  $a$  and  $b$  (decision thresholds),  $t0$ ,  $st0$ ,  $theta0$  (minimal threshold),  $thetamax$  (maximal threshold; the others will be equidistantly spanned symmetrically for both decisions), and for time-dependent confidence race models additionally  $wrt$  and  $wint$  (as weights compared to  $wx=1$ ).

For the multiple-threshold log-normal race model (MTLNR) the parameters are:  $vmin$ ,  $vmax$  (will be equidistantly spanned across conditions),  $\mu_d1$  and  $\mu_d2$  (mean pars for boundary distance),  $s_v1$ ,  $s_v2$ ,  $s_d1$ ,  $s_d2$  (variance parameters for the accumulation rates ( $v$ ) and boundary distances ( $d$ )),  $\rho_v$  and  $\rho_d$  (correlation parameters for the correlation of accumulation rates and boundary distances between accumulators),  $t0$ ,  $st0$ ,  $theta0$  (minimal threshold),  $thetamax$  (maximal threshold; the others will be equidistantly spanned symmetrically for both decisions).

**opts.** A list with numerical values. Possible options are listed below (together with the optimization method they are used for).

- `nAttempts` (all) number of best performing initial parameter sets used for optimization; default 5, if `grid_search` is TRUE. If `grid_search` is FALSE and `init_grid` is NULL, then `nAttempts` will be set to 1 (and any input will be ignored). If `grid_search` is FALSE and `init_grid` is not NULL, the rows of `init_grid` will be used from top to bottom (since no initial grid search is done) with not more than `nAttempts` rows used.
- `nRestarts` (all) number of successive optim routines for each of the starting parameter sets; default 5,
- `maxfun` ('bobyqa') maximum number of function evaluations; default: 5000,
- `maxit` ('Nelder-Mead' and 'L-BFGS-B') maximum iterations; default: 2000,
- `reltol` ('Nelder-Mead') relative tolerance; default:  $1e-6$ ,
- `fctr` ('L-BFGS-B') tolerance in terms of reduction factor of the objective, default:  $1e-10$

## Value

Gives a one-row data frame with columns for the different parameters as fitted result as well as additional information about the fit (`negLogLik` (for final parameters), `k` (number of parameters), `N` (number of data rows), `BIC`, `AICc` and `AIC`) and the column `fixed`, which includes all information about fixed and not fitted parameters.

**Author(s)**

Sebastian Hellmann.

**References**

Hellmann, S., Zehetleitner, M., & Rausch, M. (2023). Simultaneous modeling of choice, confidence and response time in visual perception. *Psychological Review* 2023 Mar 13. doi: 10.1037/rev0000411. Epub ahead of print. PMID: 36913292.

<https://nashjc.wordpress.com/2016/11/10/why-optim-is-out-of-date/>

Powell, M.J. (2009). The BOBYQA algorithm for bound constrained optimization without derivatives. Report DAMTP 2009/NA06

**Examples**

```
# We use one of the implemented models, "dynWEV"
# 1. Generate data
# data with positive drift (stimulus = "upper")
data <- rdynaViTE(20, a=2,v=0.5,t0=0.2,z=0.5, sz=0.1,sv=0.1, st0=0, tau=4, s=1, w=0.3)
data$stimulus <- "upper"
# data with negative drift (stimulus = "lower") but same intensity
data2 <- rdynaViTE(100, a=2,v=-0.5,t0=0.2,z=0.5,sz=0.1,sv=0.1, st0=0, tau=4, s=1, w=0.3)
data2$stimulus <- "lower"
data <- rbind(data, data2)
# Transfer response column and add dummy condition column
data$response <- ifelse(data$response==1, "upper", "lower")
data$condition <- 1
# Take some confidence thresholds for discrete ratings
threshs <- c(-Inf, 1, 2, Inf)
data$rating <- as.numeric(cut(data$conf, breaks = threshs, include.lowest = TRUE))
head(data)

# 2. Use fitting function
# Fitting the model with these opts results in a pretty bad fit
# (especially because of omitting the grid_search)

fitRTConf(data, "dynWEV", fixed=list(sym_thetas=TRUE, z=0.5, st0=0),
           grid_search = FALSE, logging=FALSE,
           opts = list(nAttempts=1, nRestarts=2, maxfun=2000))
```

## Description

This function is a wrapper of the function `fitConfModel` (see there for more information). It calls the function for every possible combination of model and participant/subject in model and data respectively. Also, see `ddynaViTE`, `d2DSD`, `dddConf`, `drm`, and `dMTLNR` for more information about the parameters.

## Usage

```
fitRTConfModels(data, models = c("dynaViTE", "2DSD", "PCRMt"),
  nRatings = NULL, fixed = list(sym_thetas = FALSE), restr_tau = Inf,
  grid_search = TRUE, opts = list(), optim_method = "bobyqa",
  logging = FALSE, precision = 3, parallel = TRUE, n.cores = NULL, ...)
```

## Arguments

<code>data</code>	<p>a data.frame where each row is one trial, containing following variables (column names can be changed by passing additional arguments of the form <code>condition="contrast"</code>):</p> <ul style="list-style-type: none"> <li>• <code>condition</code> (not necessary; for different levels of stimulus quality, will be transformed to a factor),</li> <li>• <code>rating</code> (discrete confidence judgments, should be given as integer vector; otherwise will be transformed to integer),</li> <li>• <code>rt</code> (giving the reaction times for the decision task),</li> <li>• either 2 of the following (see details for more information about the accepted formats):             <ul style="list-style-type: none"> <li>– <code>stimulus</code> (encoding the stimulus category in a binary choice task),</li> <li>– <code>response</code> (encoding the decision response),</li> <li>– <code>correct</code> (encoding whether the decision was correct; values in 0, 1)</li> </ul> </li> <li>• <code>sbj</code> alternatively subject or participant (giving the subject ID; the models given in the second argument are fitted for each subject individually. (Furthermore, if <code>logging = TRUE</code>, the ID is used in files saved with interim results and logging messages.) The output data frame reused the name of the column in the input (i.e. the output contains a subject column, if the input contains subject instead of sbj).)</li> </ul>
<code>models</code>	character vector with following possible elements "dynWEV", "2DSD", "IRM", "PCRM", "IRMt", "PCRMt", and "MTLNR" for the models to be fit.
<code>nRatings</code>	integer. Number of rating categories. If NULL, the maximum of rating and <code>length(unique(rating))</code> is used. This argument is especially important for data sets where not the whole range of rating categories is realized. If given, ratings has to be given as factor or integer.
<code>fixed</code>	list. List with parameter value pairs for parameters that should not be fitted. (see Details).
<code>restr_tau</code>	numerical or Inf or "simult_conf". Used for 2DSD and dynWEV only. Upper bound for tau. Fits will be in the interval (0, restr_tau). If FALSE tau will be unbound. For "simult_conf", see the documentation of <code>d2DSD</code> and <code>ddynaViTE</code>

grid_search	logical. If FALSE, the grid search before the optimization algorithm is omitted. The fitting is then started with a mean parameter set from the default grid. (Default: TRUE)
opts	list. A list for more control options in the optimization routines (depending on the optim_method). See details for more information.
optim_method	character. Determines which optimization function is used for the parameter estimation. Either "bobyqa" (default), "L-BFGS-B" or "Nelder-Mead". "bobyqa" uses a box-constrained optimization with quadratic interpolation. (See <a href="#">bobyqa</a> for more information.) The first two use a box-constraint optimization. For Nelder-Mead a transfinite function rescaling is used (i.e. the constrained arguments are suitably transformed to the whole real line).
logging	logical. If TRUE, a folder 'autosave/fitmodel' is created and messages about the process are printed in a logging file and to console (depending on OS). Additionally intermediate results are saved in a .RData file with the participant/subject ID in the name.
precision	numerical numeric. Precision of calculation for the density functions (see <a href="#">ddynaViTE</a> and <a href="#">dPCRm</a> for more information).
parallel	"models", "single", "both" or FALSE. If FALSE no parallelization is used in the fitting process. If "models" the fitting process is parallelized over participants and models (i.e. over the calls for fitting functions). If "single" parallelization is used within the fitting processes (over initial grid search and optimization processes for different start points, but see <a href="#">fitRTConf</a> ). If "both", parallelization is done hierarchical. For small number of models and participants "single" or "both" is preferable. Otherwise, you may use "models".
n.cores	integer vector or NULL. If parallel is "models" or "single", a single integer for the number of cores used for parallelization is required. If parallel is "both", two values are required. The first for the number of parallel model-participant combinations and the second for the parallel processes within the fitting procedures (this may be specified to match the nAttempts-Value in the opts argument. If NULL (default) the number of available cores -1 is used. If NULL and parallel is "both", the cores will be used for model-participant-parallelization, only.
...	Possibility of giving alternative variable names in data frame (in the form condition = "SOA", or response="pressedKey").

## Details

The fitting involves a first grid search through an initial grid. Then the best nAttempts parameter sets are chosen for an optimization, which is done with an algorithm, depending on the argument optim-method. The Nelder-Mead algorithm uses the R function [optim](#). The optimization routine is restarted nRestarts times with the starting parameter set equal to the best parameters from the previous routine.

**stimulus, response and correct.** Two of these columns must be given in data. If all three are given, correct will have no effect (and will be not checked!). stimulus can always be given in numerical format with values -1 and 1. response can always be given as a character vector with "lower" and "upper" as values. Correct must always be given as a 0-1-vector. If stimulus is given together with response and they both do not match the above format, they need to have the same values/levels (if

factor). In the case that only stimulus/response is given in any other format together with correct, the unique values will be sorted increasingly and the first value will be encoded as "lower"/-1 and the second as "upper"/+1.

**fixed.** Parameters that should not be fitted but kept constant. These will be dropped from the initial grid search but will be present in the output, to keep all parameters for prediction in the result. Includes the possibility for symmetric confidence thresholds for both alternative (`sym_thetas=logical`). Other examples are `z = .5`, `sv=0`, `st0=0`, `sz=0`. For race models, the possibility of setting `a='b'` (or vice versa) leads to identical upper bounds on the decision processes, which is the equivalence for `z=.5` in a diffusion process

**opts.** A list with numerical values. Possible options are listed below (together with the optimization method they are used for).

- `nAttempts` (all) number of best performing initial parameter sets used for optimization; default 5
- `nRestarts` (all) number of successive optim routines for each of the starting parameter sets; default 5,
- `maxfun` ('bobyqa') maximum number of function evaluations; default: 5000,
- `maxit` ('Nelder-Mead' and 'L-BFGS-B') maximum iterations; default: 2000,
- `reltol` ('Nelder-Mead') relative tolerance; default: 1e-6),
- `factr` ('L-BFGS-B') tolerance in terms of reduction factor of the objective, default: 1e-10)

## Value

Gives data frame with rows for each model-participant combination and columns for the different parameters as fitted result as well as additional information about the fit (`negLogLik` (for final parameters), `k` (number of parameters), `N` (number of data rows), `BIC`, `AICc` and `AIC`)

## Author(s)

Sebastian Hellmann, <sebastian.hellmann@ku.de>

## References

Hellmann, S., Zehetleitner, M., & Rausch, M. (2023). Simultaneous modeling of choice, confidence and response time in visual perception. *Psychological Review* 2023 Mar 13. doi: 10.1037/rev0000411. Epub ahead of print. PMID: 36913292.

## Examples

```
# 1. Generate data from two artificial participants
# Get random drift direction (i.e. stimulus category) and
# stimulus discriminability (two steps: hard, easy)
stimulus <- sample(c(-1, 1), 400, replace=TRUE)
discriminability <- sample(c(1, 2), 400, replace=TRUE)

# generate data for participant 1
data <- rdynaViTE(400, a=2, v=stimulus*discriminability*0.5,
                 t0=0.2, z=0.5, sz=0.1, sv=0.1, st0=0, tau=4, s=1, w=0.3)
```

```

# discretize confidence ratings (only 2 steps: unsure vs. sure)
data$rating <- as.numeric(cut(data$conf, breaks = c(-Inf, 1, Inf), include.lowest = TRUE))
data$participant = 1
data$stimulus <- stimulus
data$discriminability <- discriminability
# generate data for participant 2
data2 <- rdynaViTE(400, a=2.5, v=stimulus*discriminability*0.7,
                  t0=0.1, z=0.7, sz=0, sv=0.2, st0=0, tau=2, s=1, w=0.5)
data2$rating <- as.numeric(cut(data$conf, breaks = c(-Inf, 0.3, Inf), include.lowest = TRUE))
data2$participant = 2
data2$stimulus <- stimulus
data2$discriminability <- discriminability

# bind data from participants
data <- rbind(data, data2)
data <- data[data$response!=0, ] # drop not finished decision processes
data <- data[,-3] # drop conf measure (unobservable variable)
head(data)

# 2. Use fitting function
## Not run:
# Fitting takes very long to run and uses multiple (6) cores with this
# call:
fitRTConfModels(data, models=c("dynWEV", "PCRM"), nRatings = 2,
                 logging=FALSE, parallel="both",
                 n.cores = c(2,3), # fit two participant-model combination in parallel
                 condition="discriminability")# tell which column is "condition"

## End(Not run)

```

---

LogLikMTLNR

---

*Log-Likelihood functions for the multiple-threshold log-normal race model*


---

## Description

Computes the Log-likelihood for given data and parameters in the MTLNR. It is a wrapped version of the respective density [dMTLNR](#), where one can find more information about the parameters. . The function is mainly used inside [fitRTConf](#) for the MTLNR but exported for individual usage in other contexts.

## Usage

```
LogLikMTLNR(data, paramDf, precision = 6, data_names = list(), ...)
```

## Arguments

data	<p>a dataframe where each row is one trial. Containing following variables:</p> <ul style="list-style-type: none"> <li>• condition (not necessary; convertible to integer (e.g. factor); for different levels of stimulus quality),</li> <li>• rating (convertible to integer (e.g. factor); discrete confidence judgments),</li> <li>• rt (numeric; giving reaction times for decision task),</li> <li>• stimulus (values at least convertible to c(1,2), i.e. integer or factor; stimulus category (index of accumulator with higher drift))</li> <li>• response (values at least convertible to c(1,2); direction of decision; (index of accumulator reaching the boundary first))</li> </ul>
paramDf	<p>a list or data frame with one row. Column names should match the following names (see <a href="#">dMTLNR</a>): For different stimulus quality/mean drift rates, names should be v1, v2, v3,... (corresponding to the mean parameter for the accumulation rate for the stimulus-corresponding accumulator, therefore mu_v1 and mu_v2 are not used in this context but replaced by the parameter v); mu_d1 and mu_d2 correspond to the mean parameters for boundary distance of the two accumulators; s1 and s2 correspond to the variance parameters of the first and second boundary hitting time; rho corresponds to the correlation of boundary hitting times. Note that s_v1,s_v2,rho_v,s_d1,s_d2, and rho_d are not used in this context, although the accumulation rate-related parameters can be used to replace the above-mentioned variance parameters. Additionally, the confidence thresholds should be given by names with thetaUpper1, thetaUpper2,..., thetaLower1,... or, for symmetric thresholds only by theta1, theta2,... (see Details for the correspondence to the data)</p>
precision	numerical scalar. Precision of calculation for integration over t0.
data_names	list. Possibility of giving alternative column names for the variables in the data. By default column names are identical to the ones given in the data argument description.
...	Another possibility of giving alternative variable names in data frame (in the form condition = "SOA").

## Details

Note, that the requirements on the format of the columns for the likelihood functions are much stricter, than in [fitRTConf](#). This is because the function is very frequently called in the optimization routines of the fitting process and the preprocessing steps are therefore included in the other function.

**rating, condition.** If integer, values should range from 1 to number of possible ratings/conditions. If factor, the number of levels should be equal to number of possible ratings/conditions. This should be consistent with the parameter vector. The confidence thresholds should be named as thetaUpper1, thetaLower1,... (or theta1,... for symmetric thresholds), with the number of ratings -1 and the mean drift rates (and possibly the standard deviation in drift rates) should be denoted as v1, v2,... If only one condition is used v will be accepted as well as v1.

**stimulus, response.** stimulus and response should always be given in numerical format with values 1 and 2. Stimulus determines which of two accumulators has positive drift. The other has negative

drift with the same absolute value. Response gives the index of the accumulator that reaches the boundary first.

### Value

Numeric scalar. The summed Log-likelihood of the data given the parameters in the respective model. If one or more row-wise probabilities is  $\leq 0$ , the function returns  $-1e+12$ .

### Author(s)

Sebastian Hellmann.

### References

Reynolds, A., Kvam, P. D., Osth, A. F., & Heathcote, A. (2020). Correlated racing evidence accumulator models. *Journal of Mathematical Psychology*, 96, 102331. doi: 10.1016/j.jmp.2020.102331

### Examples

```
# 1. Generate data from an artificial participants
# Get random index for accumulator with positive
# drift (i.e. stimulus category) and
# stimulus discriminability (two steps: hard, easy)
stimulus <- sample(c(1, 2), 200, replace=TRUE)
discriminability <- sample(c(1, 2), 200, replace=TRUE)
# generate data for participant 1
data <- rMTLNR(200,
  mu_v1 = as.numeric(stimulus==1)*discriminability*0.5,
  mu_v2 = as.numeric(stimulus==2)*discriminability*0.5,
  mu_d1=1, mu_d2=1, t0=0.1)
# discretize confidence ratings (only 2 steps: unsure vs. sure)
data$rating <- as.numeric(cut(data$conf, breaks = c(0, 3, Inf), include.lowest = TRUE))
data$stimulus <- stimulus
data$discriminability <- discriminability
data <- data[, -c(3,4)] # drop Tdec and conf measure (unobservable variable)
head(data)

# 2. Define some parameter set in a data.frame
paramDf <- data.frame(v1=0.5, v2=1.0, t0=0.1, st0=0,
  mu_d1=1, mu_d2=1,
  s1=0.5, s2=0.5,
  rho=0.2, theta1=2.5)

# 3. Compute log likelihood for parameter and data
LogLikMTLNR(data, paramDf, condition="discriminability")
```

LogLikRM

*Log-Likelihood functions for the independent and partially anti-correlated race models of confidence*

## Description

Computes the Log-likelihood for given data and parameters in the IRM and PCRM with or without time-scaled confidence measure. It is a wrapped version of the respective densities [dIRM](#) and [dPCRM](#), where one can find more information about the parameters. It restricts the rates of accumulation to be the negative of each other, though (a common assumption in perceptual decision tasks). The function is mainly used inside [fitRTConf](#) for race models but exported for individual usage in other contexts.

## Usage

```
LogLikRM(data, paramDf, model = "IRM", time_scaled = FALSE,
         precision = 6, data_names = list(), ...)
```

## Arguments

data	<p>a dataframe where each row is one trial. Containing following variables:</p> <ul style="list-style-type: none"> <li>• condition (not necessary; convertible to integer (e.g. factor); for different levels of stimulus quality),</li> <li>• rating (convertible to integer (e.g. factor); discrete confidence judgments),</li> <li>• rt (numeric; giving reaction times for decision task),</li> <li>• stimulus (values at least convertible to c(1,2), i.e. integer or factor; stimulus category (index of accumulator with higher drift))</li> <li>• response (values at least convertible to c(1,2); direction of decision; (index of accumulator reaching the boundary first))</li> </ul>
paramDf	<p>a list or data frame with one row. Column names should match the names of <a href="#">RaceModels</a> parameter names (only mu1 and mu2 are not used in this context but replaced by the parameter v). For different stimulus quality/mean drift rates, names should be v1, v2, v3,... Different s parameters are possible with s1, s2, s3,... with equally many steps as for drift rates. Additionally, the confidence thresholds should be given by names with thetaUpper1, thetaUpper2,..., thetaLower1,... or, for symmetric thresholds only by theta1, theta2,... (see Details for the correspondence to the data)</p>
model	<p>character scalar. One of "IRM" or "PCRM". ("IRMt" and "PCRMt" will also be accepted. In that case, time_scaled is set to TRUE.)</p>
time_scaled	<p>logical. Whether the confidence measure should be scaled by 1/sqrt(rt). Default: TRUE.</p>
precision	<p>numerical scalar. Precision of calculation for integration over t0.</p>
data_names	<p>list. Possibility of giving alternative column names for the variables in the data. By default column names are identical to the ones given in the data argument description.</p>

... Another possibility of giving alternative variable names in data frame (in the form condition = "SOA").

## Details

Note, that the requirements on the format of the columns for the likelihood functions are much stricter, than in `fitRTConf`. This is because the function is very frequently called in the optimization routines of the fitting process and the preprocessing steps are therefore included in the other function.

**rating, condition.** If integer, values should range from 1 to number of possible ratings/conditions. If factor, the number of levels should be equal to number of possible ratings/conditions. This should be consistent with the parameter vector. The confidence thresholds should be named as `thetaUpper1, thetaLower1,...` (or `theta1,...` for symmetric thresholds), with the number of ratings -1 and the mean drift rates (and possibly the standard deviation in drift rates) should be denoted as `v1, v2,...` (and `s1, s2,...`) with the number equal to the number of conditions. If only one condition is used `v` will be accepted as well as `v1`.

**stimulus, response.** stimulus and response should always be given in numerical format with values 1 and 2. Stimulus determines which of two accumulators has positive drift. The other has negative drift with the same absolute value. Response gives the index of the accumulator that reaches the boundary first.

## Value

Numeric scalar. The summed Log-likelihood of the data given the parameters in the respective model. If one or more row-wise probabilities is  $\leq 0$ , the function returns  $-1e+12$ .

## Author(s)

Sebastian Hellmann.

## Examples

```
# 1. Generate data from an artificial participants
# Get random index for accumulator with positive
# drift (i.e. stimulus category) and
# stimulus discriminability (two steps: hard, easy)
stimulus <- sample(c(1, 2), 200, replace=TRUE)
discriminability <- sample(c(1, 2), 200, replace=TRUE)
# generate data for participant 1
data <- rPCRM(200, mu1=ifelse(stimulus==1, 1, -1)*discriminability*0.5,
  mu2=ifelse(stimulus==1, -1, 1)*discriminability*0.5,
  a=2, b=1.8, t0=0.2, st0=0, wx=0.7, wint=0.3, wrt=0)
# discretize confidence ratings (only 2 steps: unsure vs. sure)
data$rating <- as.numeric(cut(data$conf, breaks = c(0, 3, Inf), include.lowest = TRUE))
data$stimulus <- stimulus
data$discriminability <- discriminability
data <- data[data$response!=0, ] # drop not finished decision processes
data <- data[,-c(3,4)] # drop x1 and conf measure (unobservable variable)
head(data)
```

```
# 2. Define some parameter set in a data.frame
paramDf <- data.frame(a=2,b=2, v1=0.5, v2=1, t0=0.1,st0=0,
                      wx=0.6, wint=0.2, wrt=0.2,
                      theta1=4)

# 3. Compute log likelihood for parameter and data
LogLikRM(data, paramDf, model="PCRMt", condition="discriminability")
# same result
LogLikRM(data, paramDf, model="PCRM", time_scaled=TRUE,condition="discriminability")
# different
LogLikRM(data, paramDf, model="PCRM", condition="discriminability")

# same parameters used for IRM model
LogLikRM(data, paramDf, model="IRMt", condition="discriminability")
```

---

LogLikWEV

---

*Log-Likelihood functions for the dynWEV and 2DSD models of confidence*


---

## Description

Computes the Log-likelihood for given data and parameters in the dynWEV model (Hellmann et al., 2023) and the 2DSD model (Pleskac & Busmeyer, 2010). It is a wrapped version of the respective densities [ddynaViTE](#) and [d2DSD](#), where one can find more information about the parameters (z is always given relatively, in the likelihood). The function is mainly used in [fitRTConf](#) but exported for individual usage in other contexts.

## Usage

```
LogLikWEV(data, paramDf, model = "dynaViTE", simult_conf = FALSE,
          precision = 6, stop_on_error = TRUE, data_names = list(), ...)
```

## Arguments

**data** a dataframe where each row is one trial. Containing following variables:

- condition (not necessary; convertible to integer (e.g. factor); for different levels of stimulus quality),
- rating (convertible to integer (e.g. factor); discrete confidence judgments),
- rt (numeric; giving reaction times for decision task),
- stimulus (values at least convertible to c(-1,1); stimulus category (direction of evidence accumulation))
- response (characters in "upper", "lower" (or convertible to); direction of decision; correct answers are "lower" for stimulus=-1; and "upper" for stimulus=1),

paramDf	list or data.frame with one row. Names should match the names of <a href="#">dynaViTE</a> and <a href="#">2DSD</a> model specific parameter names. For different stimulus quality/mean drift rates, names should be v1, v2, v3,... Different sv and/or s parameters are possible with sv1, sv2, sv3... (s1, s2, s3,... respectively) with equally many steps as for drift rates. Additionally, the confidence thresholds should be given by names with thetaUpper1, thetaUpper2,..., thetaLower1,... or, for symmetric thresholds only by theta1, theta2,... (see Details for the correspondence to the data)
model	character scalar. One of "dynWEV" or "2DSD" for the model to fit.
simult_conf	logical. Whether in the experiment confidence was reported simultaneously with the decision, as then decision and confidence judgment are assumed to have happened subsequent before response and computations are different, when there is an observable interjudgment time (then simult_conf should be FALSE).
precision	numerical scalar. Precision of calculation for integration over z and t0.
stop_on_error	logical. If TRUE an error in the function will be returned in case of invalid parameters. Otherwise, the output will be 0 without error.
data_names	list. Possibility of giving alternative column names for the variables in the data. By default column names are identical to the ones given in the data argument description.
...	Possibility of giving alternative variable names in data frame (in the form condition = "SOA").

## Details

Note, that the requirements on the format of the columns for the likelihood functions are much stricter, than in [fitRTConf](#). This is because the function is very frequently calls in the optimization routines of the fitting process and the preprocessing steps are therefore included in that function.

**rating, condition.** If integer, values should range from 1 to number of possible ratings/conditions. If a factor, the number of levels should be equal to number of possible ratings/conditions. This should be consistent with the parameter vector. The confidence thresholds should be named as thetaUpper1, thetaLower1,... (or theta1,... for symmetric thresholds), with the number of ratings -1 and the mean drift rates (and possibly the standard deviation in drift rates) should be denoted as v1, v2,... (and sv1, sv2,.../s1, s2, ...) with the number equal to the number of conditions. If only one condition is used v will be accepted as well as v1.

**stimulus, response.** stimulus should always be given in numerical format with values -1 and 1. response should always be given as a character vector with "lower" and "upper" as values. This corresponds to the situation of Ratcliff's diffusion model (Ratcliff, 1978), where stimulus is the sign of the mean drift direction and the response is the "upper" or "lower" boundary that is first hit by the evidence accumulation. A correct decision is therefore "lower", if stimulus is -1, and "upper", if stimulus is 1.

## Value

Numeric scalar. The summed Log-likelihood of the data given the parameters in the respective model. If one or more row-wise probabilities is  $\leq 0$ , the function returns  $-1e+12$ .

**Author(s)**

Sebastian Hellmann.

**References**

Hellmann, S., Zehetleitner, M., & Rausch, M. (2023). Simultaneous modeling of choice, confidence and response time in visual perception. *Psychological Review* 2023 Mar 13. doi: 10.1037/rev0000411. Epub ahead of print. PMID: 36913292.

**Examples**

```
# 1. Generate data from an artificial participants
# Get random drift direction (i.e. stimulus category) and
# stimulus discriminability (two steps: hard, easy)
stimulus <- sample(c(-1, 1), 200, replace=TRUE)
discriminability <- sample(c(1, 2), 200, replace=TRUE)
# generate data for participant 1
data <- rdynaViTE(200, a=2, v=stimulus*discriminability*0.5,
                 t0=0.2, z=0.5, sz=0.1, sv=0.1, st0=0, tau=4, s=1, w=0.3)
# discretize confidence ratings (only 2 steps: unsure vs. sure)
data$rating <- as.numeric(cut(data$conf, breaks = c(-Inf, 1, Inf), include.lowest = TRUE))
data$stimulus <- stimulus
data$discriminability <- discriminability
data <- data[data$response!=0, ] # drop not finished decision processes
data <- data[, -3] # drop conf measure (unobservable variable)
head(data)

# 2. Define some parameter set in a data.frame
paramDf <- data.frame(a=2.5, v1=0.5, v2=1, t0=0.1, z=0.7,
                     sz=0, sv=0.2, st0=0, tau=3, w=0.3,
                     theta1=0.8, svis=0.5, sigvis=0.8)

# 3. Compute log likelihood for parameter and data
LogLikWEV(data, paramDf, model="dynWEV", condition="discriminability")
# adding the hypothetical interjudgment time to response times
# results in the same log likelihood as before when simult_conf=TRUE
data$rt <- data$rt + paramDf$tau
LogLikWEV(data, paramDf, model="dynWEV", condition="discriminability", simult_conf=TRUE)

# the same function for "2DSD" model
paramDf <- data.frame(a=2.5, v1=0.5, v2=1, t0=0.1, z=0.7,
                     sz=0, sv=0.2, st0=0, tau=3, theta1=0.8)
LogLikWEV(data, paramDf, model="2DSD", condition="discriminability", simult_conf=TRUE)
# this results in the same log likelihood as before
```

**Description**

The function `MLE_dirichlet` performs a maximum-likelihood estimation of the  $\alpha$  parameter of a Dirichlet distribution for a given sample of probability vectors.

**Usage**

```
MLE_dirichlet(probs, alpha0 = rep(1, ncol(probs)))
```

**Arguments**

<code>probs</code>	a matrix with $N$ rows representing observations of probability vectors and $K$ columns representing the classes. Therefore, values of each row should sum to 1.
<code>alpha0</code>	vector of $K=\text{ncol}(\text{probs})$ values as starting parameter for the optimization. Values have to be greater 0.

**Details**

The density of the Dirichlet distribution for  $\alpha = (\alpha_1, \dots, \alpha_K)$  and  $\alpha_i > 0 \forall i = 1, \dots, K$  is given by

$$f(p|\alpha) = \frac{1}{B(\alpha)} \prod_{i=1}^K K p_i^{\alpha_i-1},$$

if  $0 \leq p_i \leq 1 \forall i = 1, \dots, K$  and  $\sum_{i=1}^K p_i = 1$ , and  $f(p|\alpha) = 0$ , else.

The function optimizes the log-likelihood of a sample of probability vectors given in `probs` using the function `optim` and a Nelder-Mead algorithm.

**Value**

Returns a numeric vector of length  $K=\text{ncol}(\text{probs})$  representing the  $\alpha$  of the Dirichlet distribution.

**Author(s)**

Sebastian Hellmann.

**Examples**

```
probs <- matrix(c(0.2, 0.4, 0.2, 0.4, 0, 0.2, 0.4, 0.4, 0.6, 0.2, 0.2,
                  0.4, 0.4, 0.2, 0.2, 0.4, 0.8, 0.4), ncol=3)
MLE_dirichlet(probs)
```

MTLNR

*Correlated Multiple-threshold Log-normal Race Model for Decision Confidence***Description**

Probability densities and random number generators for response times, decisions and confidence judgments in the multi-threshold correlated log-normal race model (MTLNR; Reynolds et al., 2020), i.e. the probability of a given response (response: winning accumulator (1 or 2)) at a given time (rt) and the confidence measure in the interval between th1 and th2. The confidence measure is defined as the log-ratio between the time the losing accumulator would hit its boundary and the decision time (which means it is in the interval (0, Inf). The parameters for the model are:  $\mu_{v1}/\mu_{v2}$  and  $s_{v1}/s_{v2}$  the mean and standard deviation parameters for the log-normally distributed accumulation rates of the two accumulators,  $\mu_{d1}/\mu_{d2}$  and  $s_{d1}/s_{d2}$  the mean and standard deviation parameters for the log-normally distributed boundary distances of the two accumulators,  $\rho_v$  and  $\rho_d$  the correlation coefficients for the accumulation rates and the boundary distance, giving the correlation of these parameters between the two accumulators,  $t_0$  and  $st_0$  for the minimum and range of uniformly distributed non-decision times (including encoding and motor time).

**Usage**

```
dMTLNR(rt, response = 1, th1, th2, mu_v1, mu_v2, s_v1 = 1, s_v2 = 1,
       rho_v = 0, mu_d1 = 0, mu_d2 = 0, s_d1 = 1, s_d2 = 1, rho_d = 0,
       t0 = 0, st0 = 0, precision = 6, step_width = NULL)
```

```
dMTLNR_multiple_ratings(rt, response = 1, rating = 1, thresholds, mu_v1,
                        mu_v2, s_v1 = 1, s_v2 = 1, rho_v = 0, mu_d1 = 0, mu_d2 = 0,
                        s_d1 = 1, s_d2 = 1, rho_d = 0, t0 = 0, st0 = 0, precision = 6,
                        step_width = NULL)
```

```
rMTLNR(n, thresholds = NULL, mu_v1, mu_v2, s_v1 = 1, s_v2 = 1,
       rho_v = 0, mu_d1 = 0, mu_d2 = 0, s_d1 = 1, s_d2 = 1, rho_d = 0,
       t0 = 0, st0 = 0)
```

**Arguments**

rt	a numeric vector of RTs. For convenience also a data.frame with columns rt and response is possible.
response	numeric vector with values in c(1, 2), giving the accumulator that hit its boundary first.
th1	numeric. Lower bound of interval range for the confidence measure.
th2	numeric. Upper bound of interval range for the confidence measure.
mu_v1	numeric. Mean parameter of log-normally distributed accumulation rate of first accumulator
mu_v2	numeric. Mean parameter of log-normally distributed accumulation rate of second accumulator

s_v1	numeric. Standard deviation of log-normally distributed accumulation rate of first accumulator
s_v2	numeric. Standard deviation of log-normally distributed accumulation rate of second accumulator
rho_v	numeric. Correlation parameter for accumulation rates
mu_d1	numeric. Mean parameter of log-normally distributed boundary distance of first accumulator
mu_d2	numeric. Mean parameter of log-normally distributed boundary distance of second accumulator
s_d1	numeric. Standard deviation of log-normally distributed boundary distance of first accumulator
s_d2	numeric. Standard deviation of log-normally distributed boundary distance of second accumulator
rho_d	numeric. Correlation parameter for boundary distances
t0	numeric. Lower bound of non-decision time component in observable response times. Range: $t0 \geq 0$ . Default: 0.
st0	numeric. Range of a uniform distribution for non-decision time. Range: $st0 \geq 0$ . Default: 0.
precision	numerical scalar value. Precision of calculation. Determines the step size of integration w.r.t. $t0$ . Represents the number of decimals precisely computed on average. Default is 6.
step_width	numeric. Alternative way to define the precision of integration w.r.t. $t0$ by directly providing the step size for the integration.
rating	numeric vector with integer values from 1 to the number of confidence levels
thresholds	numeric vector of length $2 \times (\text{number of confidence levels})$ . Confidence thresholds, which are used to compare the confidence variable against for producing discrete confidence judgments. The first half of the entries should be increasing and represent the confidence thresholds for a response of 1; the second half of the entries should also be increasing and are the thresholds for response=2.
n	integer. The number of samples generated.

### Details

For the computation of confidence judgments, the parameters for the confidence thresholds  $th1$  and  $th2$  for the lower and upper bound of the interval for the confidence measure, or thresholds for a vector of thresholds (see Details).

The model assumes that each of the two accumulators has a log-normally distributed boundary distance  $D$  with mean parameter  $\mu_D$  and standard deviation parameter  $s_D^2$  and a log-normally distributed accumulation rate  $V$  with respective parameters. The accumulation takes the form of a linear ballistic accumulation without any noise, such that the boundary crossing times  $T = D/V$  are log-normally distributed with mean parameter  $\mu_D - \mu_V$  and variance parameter  $s_D^2 + s_V^2$ . In addition, the boundary distances for the two accumulators are correlated with the correlation determined by  $\rho_D$ . Similarly, the accumulation rates share a correlation with parameter  $\rho_V$ . Confidence

is determined by the log-ratio of the losing over the winning boundary crossing time, i.e., if the first accumulator hit its boundary first, confidence is determined by

$$conf = \log(T_2/T_1).$$

This confidence measure is then compared to the set of thresholds to produce discrete confidence judgments. This is equivalent to a confident computation based on the Balance of Evidence at decision time, although symmetry conditions for the thresholds may differ depending on the interpretation (see Reynolds et al., 2020 for more detail).

For convenience, the likelihood function allows that the first argument is a `data.frame` containing the information of the first and second argument in the columns (i.e., `rt` and `response` (and `rating` if relevant)). Other columns (as well as passing response separately as argument) will be ignored.

*Difference between dMTLNR and dMTLNR\_multiple\_ratings* The function `dMTLNR` allows to compute the probability of a `rt` and `response` with the confidence variable being within an interval given by two thresholds, `th1` and `th2`, similar to the definitions of the other density functions (like `ddynaViTE`). The function `dMTLNR_multiple_ratings` takes a vector for the discrete confidence judgments, `rating`, and a vector `thresholds`, such that the confidence interval can vary from observation to observation. The correct interval limits are chosen by the function depending on the entry in `rating`.

### Value

`dMTLNR` and `dMTLNR_multiple_ratings` return the numerical value of the probability density in a numerical vector of the same length as `rt`.

`rMTLNR` returns a `data.frame` with five columns and `n` rows. Column names are `rt` (response time), `response` (1 or 2, indicating which accumulator hit its boundary first), `Tdec` (the actual decision time (without non-decision time)), `conf` (the log of the ratio of boundary hitting times (the confidence variable)), and `rating` (the discrete confidence judgment).

The race parameters (as well as `response` (and `rating`), `th1`, and `th2`) are recycled to the length of the result (either `rt` or `n`). In other words, the functions are completely vectorized for all parameters and even the response.

### Note

The model is highly over-parametrized because the mean parameters for the boundary distances and accumulation rates trade off. Similarly, the variance parameters and correlation parameters trade off. For this reason, one may only use the first set of parameters for the accumulation rates (`mu_v1,...`).

### Author(s)

Sebastian Hellmann

### References

Reynolds, A., Kvam, P. D., Osth, A. F., & Heathcote, A. (2020). Correlated racing evidence accumulator models. *Journal of Mathematical Psychology*, 96, 102331. doi: 10.1016/j.jmp.2020.102331

**Examples**

```

# Plot rt distribution ignoring confidence
curve(dMTLNR(x, 1, th1=-Inf, th2=Inf, mu_v1=0.5, mu_v2=-0.5,
            mu_d1=1, mu_d2=1, t0=0.1), xlim=c(0,2.5), ylim=c(0,2))
curve(dMTLNR(x, 2, th1=-Inf, th2=Inf, mu_v1=0.5, mu_v2=-0.5,
            mu_d1=1, mu_d2=1, t0=0.1), col="red", add=TRUE)
# t0 indicates minimal response time possible
abline(v=0.1)

# Generate a random sample
df1 <- rMTLNR(5000, mu_v1=0.2, mu_v2=-0.2, mu_d1=1, mu_d2=1, t0=0.1)
head(df1)

# Compute density with rt and response as separate arguments
dMTLNR(seq(0, 2, by =0.4), response=2, th1=0.5, th2=2,
       mu_v1=0.2, mu_v2=-0.2, mu_d1=1, mu_d2=1, t0=0.1)

# Compute density with rt and response in data.frame argument
df1 <- subset(df1, response !=0) # drop trials where no accumulation hit its boundary
dMTLNR(df1[1:5,], th1=0, th2=Inf, mu_v1=0.2, mu_v2=-0.2,
       mu_d1=1, mu_d2=1, t0=0.1)

# Example with correlation parameters
dMTLNR(df1[1:5,], th1=0, th2=Inf, mu_v1=0.2, mu_v2=-0.2,
       mu_d1=1, mu_d2=1, rho_v=0.3, rho_d=0.2, t0=0.1)

# Example with multiple confidence ratings using dMTLNR_multiple_ratings
thresholds <- c(0.5, 1.5, 2.5, # for response=1 (increasing)
               0.3, 1.2, 2.0) # for response=2 (increasing)

# Create some sample data with ratings
sample_data <- data.frame(
  rt = c(0.8, 1.2, 0.9, 1.5, 1.1),
  response = c(1, 2, 1, 2, 1),
  rating = c(1, 2, 3, 1, 2)
)

dMTLNR_multiple_ratings(sample_data, thresholds=thresholds,
                        mu_v1=0.2, mu_v2=-0.2, mu_d1=1, mu_d2=1, t0=0.1)

# Compare RT and confidence distributions for different parameter settings
df_low_var <- rMTLNR(2000, thresholds = thresholds,
                   mu_v1=0.3, mu_v2=-0.3, mu_d1=1, mu_d2=1,
                   s_v1=0.5, s_v2=0.5, s_d1=0.5, s_d2=0.5, t0=0.1)
df_high_var <- rMTLNR(2000, thresholds=thresholds,
                    mu_v1=0.3, mu_v2=-0.3, mu_d1=1, mu_d2=1,
                    s_v1=1.5, s_v2=1.5, s_d1=1.5, s_d2=1.5, t0=0.1)

two_samples <- rbind(cbind(df_low_var, variance="low"),
                    cbind(df_high_var, variance="high"))
two_samples <- two_samples[two_samples$response != 0, ]

```

```
# Compare RT distributions
boxplot(log(rt) ~ variance + response, data = two_samples)

# Compare confidence distributions
boxplot(conf ~ variance + response, data = two_samples)
boxplot(rating ~ variance + response, data = two_samples)
if (requireNamespace("ggplot2", quietly = TRUE)) {
  require(ggplot2)
  ggplot(two_samples, aes(x = rt, y = conf)) +
    stat_density_2d(aes(fill = after_stat(density)), geom = "raster", contour = FALSE, na.rm=TRUE) +
    facet_grid(cols = vars(variance), rows = vars(response),
               labeller = "label_both") +
    xlim(c(0.2, 2.0)) + ylim(c(-2, 4))
}
```

---

PDFtoQuantiles

---

*Get Quantiles from vectors of PDF or CDF values*


---

## Description

CDFtoQuantiles computes quantiles for a given CDF. PDFtoQuantiles computes the quantiles for given PDF values within groups of other variables, if available.

## Usage

```
PDFtoQuantiles(pdf_df, p = c(0.1, 0.3, 0.5, 0.7, 0.9), agg_over = NULL,
               scaled = FALSE)
```

```
CDFtoQuantiles(cdf, x = NULL, p)
```

## Arguments

pdf_df	dataframe. Should have at least two columns: <ul style="list-style-type: none"> <li>• rt (for reaction times) or x for the support values of the pdf</li> <li>• dens or pdf for the pdf values</li> <li>• All other columns will be used as grouping factors, for which separate quantiles will be returned.</li> </ul>
p	numeric vector. Probabilities for returned quantiles. Default: c(.1, .3, .5, .7, .9).
agg_over	character. Names of columns in pdf_df to aggregate over (using the mean of densities, which is valid only, if groups occur with equal probabilities) before computing the quantiles.
scaled	logical. Indicating whether the pdf values are from a proper probability distribution. Non-scaled pdfs will scaled to 1. If scaled is TRUE, this may cause problems with high probabilities. In any case we strongly recommend to cover the most probability mass with the values in the support vector.

cdf	numeric. A increasing vector of the same length as x giving the CDF for respective x-Values. Dataframe inputs are accepted. If a column x is available there, this will be used as support values.
x	numeric. A increasing vector of same length as cdf. Can also be specified as column of cdf.

## Details

For a reasonable accuracy the number of steps in the support column (rt/x) should be high, i.e. the distance between values small. We recommend, to ensure that the support vector in the input to be equidistant, i.e. the difference between consecutive support values should be constant, though this is not required. If both column names x and rt are present in pdf\_df, rt will be preferred. Attention should be given to the columns of pdf\_df other than rt/x and dens/pdf.

The column for the pdf may be scaled to integrate to 1 but do not have to.

### Quantile computation in the dynConfIR package:

As argument pdf\_df, the outputs of predictRT and predictRTModels from the dynConfIR package can be used. In the context of confidence models grouping factors often used are conditions, correct/incorrect answers and confidence ratings.

## Value

PDFtoQuantiles returns a tibble with columns p and q indicating probabilities and respective quantiles. Furthermore, the output has grouping columns identical to the additional columns in the input (without rt/x, dens and densscaled), but without the ones in the agg\_over argument. CDFtoQuantiles returns only a data.frame with columns p and q.

## Author(s)

Sebastian Hellmann.

## Examples

```
## Demonstrate PDFtoQuantiles
pred <- expand.grid(model = c("dynWEV", "PCRMT"),
                   rt = seq(0, 15, length.out=1200),
                   condition = c(1,2,3),
                   rating = c(1,2))
pred$dens <- dchisq(pred$rt, 3) # pdf may also be used as column name
head(pred)
res <- PDFtoQuantiles(pred, p=c(0.3, 0.5, 0.7))
head(res)
nrow(res) #= 3(quantiles)*2(models)*3(conditions)*2(rating)
# Compare to true quantiles of Chi-square distribution
qchisq(p=c(0.3, 0.5, 0.7), 3)
res$q[1:3]

res2 <- PDFtoQuantiles(pred, p=c(0.3, 0.5, 0.7), agg_over = "model")
nrow(res2) #=18 because res aggregated over models
```

```

pred$pdf <- dchisq(pred$rt, 3)
head(pred)
# following call throws a warning, because both columns pdf and dens are present
PDFtoQuantiles(pred, p=c(0.3, 0.5, 0.7), agg_over = "model")

pred2 <- data.frame(rt=seq(0, 7, length.out=100))
pred2$dens <- dchisq(pred2$rt, 5)
# following call throws a warning, because density is assumed to be scaled (scaled=TRUE), i.e.
# integrate to 1, but the .95 quantile is not reached in the rt column
PDFtoQuantiles(pred2, p=c(0.3, 0.5, 0.95), scaled=TRUE) # Gives a warning

## Demonstrate CDFtoQuantiles
X <- seq(-2, 2, length.out=300)
pdf_values <- pnorm(X)
CDFtoQuantiles(pdf_values, X, p=c(0.2, 0.5, 0.8))
qnorm(c(0.2, 0.5, 0.8))

```

---

predictDDConf	<i>Prediction of Confidence Rating and Reaction Time Distribution in the drift diffusion confidence model</i>
---------------	---

---

## Description

predictDDConf\_Conf predicts the categorical response distribution of decision and confidence ratings, predictDDConf\_RT computes the RT distribution (density) in the drift diffusion confidence model (Hellmann et al., 2023), given specific parameter constellations. See [dDDConf](#) for more information about the model and parameters.

## Usage

```

predictDDConf_Conf(paramDf, maxrt = Inf, subdivisions = 100L,
  stop.on.error = FALSE, .progress = TRUE)

predictDDConf_RT(paramDf, maxrt = 9, subdivisions = 100L, minrt = NULL,
  scaled = FALSE, DistConf = NULL, .progress = TRUE)

```

## Arguments

paramDf	a list or data frame with one row. Column names should match the names of <a href="#">DD-Conf</a> model parameter names. For different stimulus quality/mean drift rates, names should be v1, v2, v3,... Different sv and/or s parameters are possible with sv1, sv2, sv3... (s1, s2, s3,... respectively) with equally many steps as for drift rates. Additionally, the confidence thresholds should be given by names with thetaUpper1, thetaUpper2,..., thetaLower1,... or, for symmetric thresholds only by theta1, theta2,....
---------	---

maxrt	numeric. The maximum RT for the integration/density computation. Default: 15 (for predictDDConf_Conf (integration)), 9 (for predictDDConf_RT).
subdivisions	integer (default: 100). For predictDDConf_Conf it is used as argument for the inner integral routine. For predictDDConf_RT it is the number of points for which the density is computed.
stop.on.error	logical. Argument directly passed on to integrate. Default is FALSE, since the densities invoked may lead to slow convergence of the integrals (which are still quite accurate) which causes R to throw an error.
.progress	logical. If TRUE (default) a progress bar is drawn to the console.
minrt	numeric or NULL(default). The minimum rt for the density computation.
scaled	logical. For predictDDConf_RT. Whether the computed density should be scaled to integrate to one (additional column densscaled). Otherwise the output contains only the defective density (i.e. its integral is equal to the probability of a response and not 1). If TRUE, the argument DistConf should be given, if available. Default: FALSE.
DistConf	NULL or data.frame. A data.frame or matrix with column names, giving the distribution of response and rating choices for different conditions and stimulus categories in the form of the output of predictDDConf_Conf. It is only necessary, if scaled=TRUE, because these probabilities are used for scaling. If scaled=TRUE and DistConf=NULL, it will be computed with the function predictDDConf_Conf, which takes some time and the function will throw a message. Default: NULL

## Details

The function predictDDConf\_Conf consists merely of an integration of the response time density, [dDDConf](#), over the response time in a reasonable interval (0 to maxrt). The function predictDDConf\_RT wraps these density functions to a parameter set input and a data.frame output. For the argument paramDf, the output of the fitting function [fitRTConf](#) with the DDConf model may be used.

## Value

predictDDConf\_Conf returns a data.frame/tibble with columns: condition, stimulus, response, rating, correct, p, info, err. p is the predicted probability of a response and rating, given the stimulus category and condition. info and err refer to the respective outputs of the integration routine used for the computation. predictDDConf\_RT returns a data.frame/tibble with columns: condition, stimulus, response, rating, correct, rt and dens (and densscaled, if scaled=TRUE).

## Note

Different parameters for different conditions are only allowed for drift rate  $v$ , drift rate variability  $sv$ , and process variability  $s$ . Otherwise,  $s$  is not required in paramDf but set to 1 by default. All other parameters are used for all conditions.

## Author(s)

Sebastian Hellmann.

## References

Hellmann, S., Zehetleitner, M., & Rausch, M. (2023). Simultaneous modeling of choice, confidence and response time in visual perception. *Psychological Review* 2023 Mar 13. doi: 10.1037/rev0000411. Epub ahead of print. PMID: 36913292.

## Examples

```
# 1. Define some parameter set in a data.frame
paramDf <- data.frame(a=2, v1=0.5, v2=1, t0=0.1, z=0.55,
                      sz=0, sv=0.2, st0=0, theta1=0.8)

# 2. Predict discrete Choice x Confidence distribution:
preds_Conf <- predictDDConf_Conf(paramDf, maxrt = 15)
head(preds_Conf)

# 3. Compute RT density
preds_RT <- predictDDConf_RT(paramDf, maxrt=4, subdivisions=200) #(scaled=FALSE)
# same output with scaled density column:
preds_RT <- predictDDConf_RT(paramDf, maxrt=4, subdivisions=200,
                             scaled=TRUE, DistConf = preds_Conf)
head(preds_RT)

# Example of visualization
library(ggplot2)
preds_Conf$rating <- factor(preds_Conf$rating, labels=c("unsure", "sure"))
preds_RT$rating <- factor(preds_RT$rating, labels=c("unsure", "sure"))
ggplot(preds_Conf, aes(x=interaction(rating, response), y=p))+
  geom_bar(stat="identity")+
  facet_grid(cols=vars(stimulus), rows=vars(condition), labeller = "label_both")
ggplot(preds_RT, aes(x=rt, color=interaction(rating, response), y=dens))+
  geom_line(stat="identity")+
  facet_grid(cols=vars(stimulus), rows=vars(condition), labeller = "label_both")+
  theme(legend.position = "bottom")
ggplot(aggregate(densscaled~rt+correct+rating+condition, preds_RT, mean),
       aes(x=rt, color=rating, y=densscaled))+
  geom_line(stat="identity")+
  facet_grid(cols=vars(condition), rows=vars(correct), labeller = "label_both")+
  theme(legend.position = "bottom")

# Use PDFtoQuantiles to get predicted RT quantiles
head(PDFtoQuantiles(preds_RT, scaled = FALSE))
```

## Description

predictMTLNR\_Conf predicts the categorical response distribution of decision and confidence ratings, predictMTLNR\_RT computes the RT distribution (density) in the multiple-threshold log-normal noise race model (Reynolds et al., 2020), given specific parameter constellations. See [dMTLNR](#) for more information about the models and parameters.

## Usage

```
predictMTLNR_Conf(paramDf, maxrt = Inf, subdivisions = 100L,
  stop.on.error = FALSE, .progress = TRUE)

predictMTLNR_RT(paramDf, maxrt = 9, subdivisions = 100L, minrt = NULL,
  scaled = FALSE, DistConf = NULL, .progress = TRUE)
```

## Arguments

paramDf	a list or data frame with one row. Column names should match the following names (see <a href="#">dMTLNR</a> ): For different stimulus quality/mean drift rates, names should be v1, v2, v3,... (corresponding to the mean parameter for the accumulation rate for the stimulus-corresponding accumulator, therefore mu_v1 and mu_v2 are not used in this context but replaced by the parameter v); mu_d1 and mu_d2 correspond to the mean parameters for boundary distance of the two accumulators; s1 and s2 correspond to the variance parameters of the first and second boundary hitting time; rho corresponds to the correlation of boundary hitting times. Note that s_v1,s_v2,rho_v,s_d1,s_d2, and rho_d are not used in this context, although the accumulation rate-related parameters can be used to replace the above-mentioned variance parameters. Additionally, the confidence thresholds should be given by names with thetaUpper1, thetaUpper2,..., thetaLower1,... or, for symmetric thresholds only by theta1, theta2,... (see Details for the correspondence to the data)
maxrt	numeric. The maximum RT for the integration/density computation. Default: 15 (for predictMTLNR_Conf (integration)), 9 (for predictMTLNR_RT).
subdivisions	integer (default: 100). For predictMTLNR_Conf it is used as argument for the inner integral routine. For predictMTLNR_RT it is the number of points for which the density is computed.
stop.on.error	logical. Argument directly passed on to integrate. Default is FALSE, since the densities invoked may lead to slow convergence of the integrals (which are still quite accurate) which causes R to throw an error.
.progress	logical. If TRUE (default) a progress bar is drawn to the console.
minrt	numeric or NULL(default). The minimum rt for the density computation.
scaled	logical. For predictMTLNR_RT. Whether the computed density should be scaled to integrate to one (additional column densscaled). Otherwise the output contains only the defective density (i.e. its integral is equal to the probability of a response and not 1). If TRUE, the argument DistConf should be given, if available. Default: FALSE.

**DistConf** NULL or data.frame. A data.frame or matrix with column names, giving the distribution of response and rating choices for different conditions and stimulus categories in the form of the output of predictMTLNR\_Conf. It is only necessary, if scaled=TRUE, because these probabilities are used for scaling. If scaled=TRUE and DistConf=NULL, it will be computed with the function predictMTLNR\_Conf, which takes some time and the function will throw a message. Default: NULL

## Details

The function predictMTLNR\_Conf consists merely of an integration of the response time density, [dMTLNR](#), over the response time in a reasonable interval (0 to maxrt). The function predictMTLNR\_RT wraps these density functions to a parameter set input and a data.frame output. For the argument paramDf, the output of the fitting function [fitRTConf](#) with the respective model may be used.

The names of the accumulation rate parameters differ from those used in [dMTLNR](#) because the accumulation rates for the two options depend on stimulus and condition. Here, the mean parameter for the accumulation rate of the correct accumulator is  $v$  ( $v_1$ ,  $v_2$ ,... respectively) in paramDf and the other one has a mean parameter of 0.

## Value

predictMTLNR\_Conf returns a data.frame/tibble with columns: condition, stimulus, response, rating, correct, p, info, err. p is the predicted probability of a response and rating, given the stimulus category and condition. info and err refer to the respective outputs of the integration routine used for the computation. predictMTLNR\_RT returns a data.frame/tibble with columns: condition, stimulus, response, rating, correct, rt and dens (and densscaled, if scaled=TRUE).

## Note

Different parameters for different conditions are only allowed for drift rate,  $v$ . All other parameters are used for all conditions.

## Author(s)

Sebastian Hellmann.

## References

Reynolds, A., Kvam, P. D., Osth, A. F., & Heathcote, A. (2020). Correlated racing evidence accumulator models. *Journal of Mathematical Psychology*, 96, 102331. doi: 10.1016/j.jmp.2020.102331

## Examples

```
# 1. Define some parameter set in a data.frame
paramDf <- data.frame(v1=0.5, v2=1.0, t0=0.1, st0=0,
                      mu_d1=1, mu_d2=1,
                      s1=0.5, s2=0.5, rho=0.2,
                      theta1=0.8, theta2=1.5)
```

```

# 2. Predict discrete Choice x Confidence distribution:
preds_Conf <- predictMTLNR_Conf(paramDf, maxrt=7, subdivisions=50)
head(preds_Conf)

# 3. Compute RT density
preds_RT <- predictMTLNR_RT(paramDf, maxrt=7, subdivisions=50)
# same output with scaled density column:
preds_RT <- predictMTLNR_RT(paramDf, maxrt=7, subdivisions=50,
                             scaled=TRUE, DistConf = preds_Conf)
head(preds_RT)

# produces a warning, if scaled=TRUE and DistConf missing
preds_RT <- predictMTLNR_RT(paramDf, maxrt=7, subdivisions=50,
                             scaled=TRUE)

# Example of visualization
library(ggplot2)
preds_Conf$rating <- factor(preds_Conf$rating, labels=c("unsure", "medium", "sure"))
preds_RT$rating <- factor(preds_RT$rating, labels=c("unsure", "medium", "sure"))
ggplot(preds_Conf, aes(x=interaction(rating, response), y=p))+
  geom_bar(stat="identity")+
  facet_grid(cols=vars(stimulus), rows=vars(condition), labeller = "label_both")
ggplot(preds_RT, aes(x=rt, color=interaction(rating, response), y=dens))+
  geom_line(stat="identity")+
  facet_grid(cols=vars(stimulus), rows=vars(condition), labeller = "label_both")+
  theme(legend.position = "bottom")
ggplot(aggregate(densscaled~rt+correct+rating+condition, preds_RT, mean),
       aes(x=rt, color=rating, y=densscaled))+
  geom_line(stat="identity")+
  facet_grid(cols=vars(condition), rows=vars(correct), labeller = "label_both")+
  theme(legend.position = "bottom")

# Use PDFtoQuantiles to get predicted RT quantiles
# (produces warning because of few rt steps (--> inaccurate calculations))
PDFtoQuantiles(preds_RT, scaled = FALSE)

# Example with asymmetric confidence thresholds
paramDf_asym <- data.frame(v1=0.5, v2=1.0, t0=0.1, st0=0,
                           mu_d1=1, mu_d2=1,
                           s1=0.5, s2=0.5, rho=0.2,
                           thetaLower1=0.5, thetaLower2=1.2,
                           thetaUpper1=0.7, thetaUpper2=1.8)

preds_Conf_asym <- predictMTLNR_Conf(paramDf_asym, maxrt=7, subdivisions=50)
head(preds_Conf_asym)

# Example with multiple conditions
paramDf_multi <- data.frame(v1=0.3, v2=0.6, v3=1.2, t0=0.1, st0=0,
                            mu_d1=1, mu_d2=1,

```

```

s1=0.5, s2=0.5, rho=0.2,
theta1=0.8, theta2=1.5)

preds_Conf_multi <- predictMTLNR_Conf(paramDf_multi, maxrt=7, subdivisions=50)
head(preds_Conf_multi)

```

---

predictRM	<i>Prediction of Confidence Rating and Reaction Time Distribution in race models of confidence</i>
-----------	--

---

## Description

predictRM\_Conf predicts the categorical response distribution of decision and confidence ratings, predictRM\_RT computes the RT distribution (density) in the independent and partially anti-correlated race models (Hellmann et al., 2023), given specific parameter constellations. See [Race-Models](#) for more information about the models and parameters.

## Usage

```

predictRM_Conf(paramDf, model = "IRM", time_scaled = FALSE, maxrt = Inf,
  subdivisions = 100L, stop.on.error = FALSE, .progress = TRUE)

predictRM_RT(paramDf, model = "IRM", time_scaled = FALSE, maxrt = 9,
  subdivisions = 100L, minrt = NULL, scaled = FALSE, DistConf = NULL,
  .progress = TRUE)

```

## Arguments

paramDf	a list or data frame with one row. Column names should match the names of <a href="#">RaceModels</a> parameter names (only mu1 and mu2 are not used in this context but replaced by the parameter v). For different stimulus quality/mean drift rates, names should be v1, v2, v3,... Different s parameters are possible with s1, s2, s3,... with equally many steps as for drift rates. Additionally, the confidence thresholds should be given by names with thetaUpper1, thetaUpper2,..., thetaLower1,... or, for symmetric thresholds only by theta1, theta2,....
model	character scalar. One of "IRM" or "PCRM". ("IRMt" and "PCRMt" will also be accepted. In that case, time_scaled is set to TRUE.)
time_scaled	logical. Whether the confidence measure should be scaled by 1/sqrt(rt). Default: FALSE. (It is set to TRUE, if model is "IRMt" or "PCRMt")
maxrt	numeric. The maximum RT for the integration/density computation. Default: 15 (for predictRM_Conf (integration)), 9 (for predictRM_RT).
subdivisions	integer (default: 100). For predictRM_Conf it is used as argument for the inner integral routine. For predictRM_RT it is the number of points for which the density is computed.

<code>stop.on.error</code>	logical. Argument directly passed on to <code>integrate</code> . Default is <code>FALSE</code> , since the densities invoked may lead to slow convergence of the integrals (which are still quite accurate) which causes R to throw an error.
<code>.progress</code>	logical. If <code>TRUE</code> (default) a progress bar is drawn to the console.
<code>minrt</code>	numeric or <code>NULL</code> (default). The minimum <code>rt</code> for the density computation.
<code>scaled</code>	logical. For <code>predictRM_RT</code> . Whether the computed density should be scaled to integrate to one (additional column <code>densscaled</code> ). Otherwise the output contains only the defective density (i.e. its integral is equal to the probability of a response and not 1). If <code>TRUE</code> , the argument <code>DistConf</code> should be given, if available. Default: <code>FALSE</code> .
<code>DistConf</code>	<code>NULL</code> or <code>data.frame</code> . A <code>data.frame</code> or <code>matrix</code> with column names, giving the distribution of response and rating choices for different conditions and stimulus categories in the form of the output of <code>predictRM_Conf</code> . It is only necessary, if <code>scaled=TRUE</code> , because these probabilities are used for scaling. If <code>scaled=TRUE</code> and <code>DistConf=NULL</code> , it will be computed with the function <code>predictRM_Conf</code> , which takes some time and the function will throw a message. Default: <code>NULL</code>

## Details

The function `predictRM_Conf` consists merely of an integration of the response time density, [dIRM](#) and [dPCRM](#), over the response time in a reasonable interval (0 to `maxrt`). The function `predictRM_RT` wraps these density functions to a parameter set input and a `data.frame` output. For the argument `paramDf`, the output of the fitting function [fitRTConf](#) with the respective model may be used.

The drift rate parameters differ from those used in [dIRM/dPCRM](#) since in many perceptual decision experiments the drift on one accumulator is assumed to be the negative of the other. The drift rate of the correct accumulator is  $v$  ( $v_1, v_2, \dots$  respectively) in `paramDf`.

## Value

`predictRM_Conf` returns a `data.frame/tibble` with columns: `condition`, `stimulus`, `response`, `rating`, `correct`, `p`, `info`, `err`. `p` is the predicted probability of a response and rating, given the stimulus category and condition. `info` and `err` refer to the respective outputs of the integration routine used for the computation. `predictRM_RT` returns a `data.frame/tibble` with columns: `condition`, `stimulus`, `response`, `rating`, `correct`, `rt` and `dens` (and `densscaled`, if `scaled=TRUE`).

## Note

Different parameters for different conditions are only allowed for drift rate,  $v$ , and process variability  $s$ . All other parameters are used for all conditions.

## Author(s)

Sebastian Hellmann.

## References

Hellmann, S., Zehetleitner, M., & Rausch, M. (2023). Simultaneous modeling of choice, confidence and response time in visual perception. *Psychological Review* 2023 Mar 13. doi: 10.1037/rev0000411. Epub ahead of print. PMID: 36913292.

## Examples

```
# Examples for "PCRM" model (equivalent applicable for "IRM" model)
# 1. Define some parameter set in a data.frame
paramDf <- data.frame(a=2,b=2, v1=0.5, v2=1, t0=0.1,st0=0,
                      wx=0.6, wint=0.2, wrt=0.2,
                      theta1=4)

# 2. Predict discrete Choice x Confidence distribution:
preds_Conf <- predictRM_Conf(paramDf, "PCRM", time_scaled=TRUE)
# equivalent:
preds_Conf <- predictRM_Conf(paramDf, "PCRMt")
head(preds_Conf)

# 3. Compute RT density
preds_RT <- predictRM_RT(paramDf, "PCRMt", maxrt=7, subdivisions=50)
# same output with scaled density column:
preds_RT <- predictRM_RT(paramDf, "PCRMt", maxrt=7, subdivisions=50,
                        scaled=TRUE, DistConf = preds_Conf)
head(preds_RT)

# produces a warning, if scaled=TRUE and DistConf missing
preds_RT <- predictRM_RT(paramDf, "PCRMt", maxrt=7, subdivisions=50,
                        scaled=TRUE)

# Example of visualization
library(ggplot2)
preds_Conf$rating <- factor(preds_Conf$rating, labels=c("unsure", "sure"))
preds_RT$rating <- factor(preds_RT$rating, labels=c("unsure", "sure"))
ggplot(preds_Conf, aes(x=interaction(rating, response), y=p))+
  geom_bar(stat="identity")+
  facet_grid(cols=vars(stimulus), rows=vars(condition), labeller = "label_both")
ggplot(preds_RT, aes(x=rt, color=interaction(rating, response), y=dens))+
  geom_line(stat="identity")+
  facet_grid(cols=vars(stimulus), rows=vars(condition), labeller = "label_both")+
  theme(legend.position = "bottom")
ggplot(aggregate(densscaled~rt+correct+rating+condition, preds_RT, mean),
      aes(x=rt, color=rating, y=densscaled))+
  geom_line(stat="identity")+
  facet_grid(cols=vars(condition), rows=vars(correct), labeller = "label_both")+
  theme(legend.position = "bottom")

# Use PDFtoQuantiles to get predicted RT quantiles
# (produces warning because of few rt steps (--> inaccurate calculations))
PDFtoQuantiles(preds_RT, scaled = FALSE)
```

---

predictRTConf	<i>Prediction of confidence rating and response time distribution for sequential sampling confidence models</i>
---------------	---

---

## Description

predictConf predicts the categorical response distribution of decision and confidence ratings, predictRT computes the predicted RT distribution (density) for the sequential sampling confidence model specified by the argument model, given specific parameter constellations. This function calls the respective functions for diffusion based models (dynWEV and 2DSD: [predictWEV](#)) and race models (IRM, PCRM, IRMt, and PCRMt: [predictRM](#); and MTLNR: [predictMTLNR](#)).

## Usage

```
predictConf(paramDf, model = NULL, maxrt = Inf, subdivisions = 100L,
  simult_conf = FALSE, stop.on.error = FALSE, .progress = TRUE)

predictRT(paramDf, model = NULL, maxrt = 9, subdivisions = 100L,
  minrt = NULL, simult_conf = FALSE, scaled = FALSE, DistConf = NULL,
  .progress = TRUE)
```

## Arguments

paramDf	a list or dataframe with one row. Column names should match the names of the respective model parameters. For different stimulus quality/mean drift rates, names should be v1, v2, v3,... Different s parameters are possible with s1, s2, s3... with equally many steps as for drift rates (same for sv parameter in dynWEV and 2DSD), except for MTLNR. Additionally, the confidence thresholds should be given by names with thetaUpper1, thetaUpper2,..., thetaLower1,... or, for symmetric thresholds only by theta1, theta2,....
model	character scalar. One of "2DSD", "dynWEV", "IRM", "PCRM", "IRMt", "PCRMt", or "MTLNR".
maxrt	numeric. The maximum RT for the integration/density computation. Default: 15 (for predictConf (integration)), 9 (for predictRT).
subdivisions	integer (default: 100). For predictConf it is used as argument for the inner integral routine. For predictRT it is the number of points for which the density is computed.
simult_conf	logical, only relevant for dynWEV and 2DSD. Whether in the experiment confidence was reported simultaneously with the decision, as then decision and confidence judgment are assumed to have happened subsequent before response and computations are different, when there is an observable interjudgment time (then simult_conf should be FALSE).
stop.on.error	logical. Argument directly passed on to integrate. Default is FALSE, since the densities invoked may lead to slow convergence of the integrals (which are still quite accurate) which causes R to throw an error.

<code>.progress</code>	logical. If TRUE (default) a progress bar is drawn to the console.
<code>minrt</code>	numeric or NULL(default). The minimum rt for the density computation.
<code>scaled</code>	logical. For predictRT. Whether the computed density should be scaled to integrate to one (additional column <code>densscaled</code> ). Otherwise the output contains only the defective density (i.e. its integral is equal to the probability of a response and not 1). If TRUE, the argument <code>DistConf</code> should be given, if available. Default: FALSE.
<code>DistConf</code>	NULL or data.frame. A data.frame or matrix with column names, giving the distribution of response and rating choices for different conditions and stimulus categories in the form of the output of <code>predictConf</code> . It is only necessary, if <code>scaled=TRUE</code> , because these probabilities are used for scaling. If <code>scaled=TRUE</code> and <code>DistConf=NULL</code> , it will be computed with the function <code>predictRM_Conf</code> , which takes some time and the function will throw a message. Default: NULL

## Details

The function `predictConf` consists merely of an integration of the reaction time density of the given model, `{d*model*}`, over the response time in a reasonable interval (0 to `maxrt`). The function `predictRT` wraps these density functions to a parameter set input and a data.frame output. For the argument `paramDf`, the output of the fitting function `fitRTConf` with the respective model may be used.

## Value

`predictConf` returns a data.frame/tibble with columns: `condition`, `stimulus`, `response`, `rating`, `correct`, `p`, `info`, `err`. `p` is the predicted probability of a response and rating, given the stimulus category and condition. `info` and `err` refer to the respective outputs of the integration routine used for the computation. `predictRT` returns a data.frame/tibble with columns: `condition`, `stimulus`, `response`, `rating`, `correct`, `rt` and `dens` (and `densscaled`, if `scaled=TRUE`).

## Note

Different parameters for different conditions are only allowed for drift rate, `v`, drift rate variability, `sv` (in `dynWEV` and `2DSD`), and process variability `s` (not for `MTLNR`). All other parameters are used for all conditions.

## Author(s)

Sebastian Hellmann.

## Examples

```
# Examples for "dynWEV" model (equivalent applicable for
# all other models (with different parameters!))

# 1. Define some parameter set in a data.frame
paramDf <- data.frame(a=1.5, v1=0.2, v2=1, t0=0.1, z=0.52,
                      sz=0.3, sv=0.4, st0=0, tau=3, w=0.5,
                      theta1=1, svis=0.5, sigvis=0.8)
```

```

# 2. Predict discrete Choice x Confidence distribution:
preds_Conf <- predictConf(paramDf, "dynWEV", maxrt = 25, simult_conf=TRUE)
head(preds_Conf)

# 3. Compute RT density
preds_RT <- predictRT(paramDf, "dynWEV") #(scaled=FALSE)
# same output with default rt-grid and without scaled density column:
preds_RT <- predictRT(paramDf, "dynWEV", maxrt=5, subdivisions=200,
                      minrt=paramDf$tau+paramDf$t0, simult_conf = TRUE,
                      scaled=TRUE, DistConf = preds_Conf)
head(preds_RT)

# produces a warning, if scaled=TRUE and DistConf missing
preds_RT <- predictRT(paramDf, "dynWEV",
                      scaled=TRUE)

# Example of visualization
library(ggplot2)
preds_Conf$rating <- factor(preds_Conf$rating, labels=c("unsure", "sure"))
preds_RT$rating <- factor(preds_RT$rating, labels=c("unsure", "sure"))
ggplot(preds_Conf, aes(x=interaction(rating, response), y=p))+
  geom_bar(stat="identity")+
  facet_grid(cols=vars(stimulus), rows=vars(condition), labeller = "label_both")
ggplot(preds_RT, aes(x=rt, color=interaction(rating, response), y=densscaled))+
  geom_line(stat="identity")+
  facet_grid(cols=vars(stimulus), rows=vars(condition), labeller = "label_both")+
  theme(legend.position = "bottom")+ ggtitle("Scaled Densities")
ggplot(aggregate(dens~rt+correct+rating+condition, preds_RT, mean),
      aes(x=rt, color=rating, y=dens))+
  geom_line(stat="identity")+
  facet_grid(cols=vars(condition), rows=vars(correct), labeller = "label_both")+
  theme(legend.position = "bottom")+ ggtitle("Non-Scaled Densities")

# Use PDFtoQuantiles to get predicted RT quantiles
head(PDFtoQuantiles(preds_RT, scaled = FALSE))

```

---

predictRTConfModels	<i>Prediction of confidence and RT distributions for several sequential sampling confidence models and parameter constellations in parallel</i>
---------------------	---

---

## Description

This function is a wrapper around the functions [predictRTConf](#) (see there for more information). It calls the respective function for predicting the response distribution (discrete decision and rating outcomes) and the rt density (density for decision, rating and response time) for every model and

participant/subject combination in paramDf. Also, see [ddynaViTE](#), [d2DSD](#), [dMTLNR](#), and [dRM](#) for more information about the parameters.

## Usage

```
predictConfModels(paramDf, maxrt = Inf, subdivisions = 100L,
  simult_conf = FALSE, stop.on.error = FALSE, .progress = TRUE,
  parallel = FALSE, n.cores = NULL)

predictRTModels(paramDf, maxrt = 9, subdivisions = 100L, minrt = NULL,
  simult_conf = FALSE, scaled = FALSE, DistConf = NULL,
  .progress = TRUE, parallel = FALSE, n.cores = NULL)
```

## Arguments

paramDf	a dataframe with one row per combination of model and participant/parameter set. Columns may include a participant (sbj, or subject) column, and must include a model column and the names of the model parameters. For different stimulus quality/mean drift rates, names should be v1, v2, v3,... Different s parameters are possible with s1, s2, s3... with equally many steps as for drift rates (same for sv parameter in dynWEV and 2DSD), except for MTLNR. Additionally, the confidence thresholds should be given by names with thetaUpper1, thetaUpper2,..., thetaLower1,... or, for symmetric thresholds only by theta1, theta2,....
maxrt	numeric. The maximum RT for the integration/density computation. Default: 15 (for predictConfModels (integration)) and 9 (for predictRTModels).
subdivisions	integer (default: 100). For predictConfModels it is used as argument for the inner integral routine. For predictRTModels it is the number of points for which the density is computed.
simult_conf	logical, only relevant for dynWEV and 2DSD. Whether in the experiment confidence was reported simultaneously with the decision, as then decision and confidence judgment are assumed to have happened subsequent before response and computations are different, when there is an observable interjudgment time (then simult_conf should be FALSE).
stop.on.error	logical. Argument directly passed on to integrate. Default is FALSE, since the densities invoked may lead to slow convergence of the integrals (which are still quite accurate) which causes R to throw an error.
.progress	logical. If TRUE (default) a progress bar is drawn to the console. (Works for some OS only when parallel=FALSE.)
parallel	logical. If TRUE, prediction is parallelized over participants and models (i.e. over the calls for the respective <a href="#">predictRTConf</a> functions).
n.cores	integer. If parallel is TRUE, the number of cores used for parallelization is required. If NULL (default) the number of available cores -1 is used.
minrt	numeric or NULL(default). The minimum rt for the density computation. If NULL, the minimal possible response time possible with given parameters will be used (min(t0)).

scaled	logical. Whether the computed density should be scaled to integrate to one (additional column densscaled). Otherwise the output contains only the defective density (i.e. its integral is equal to the probability of a response and not 1). If TRUE, the argument DistConf should be given, if available. Default: FALSE.
DistConf	NULL or data.frame. A data.frame with participant and model columns and columns, giving the distribution of response and rating choices for different conditions and stimulus categories in the form of the output of predictConfModels. It is only necessary if scaled=TRUE, because these probabilities are used for scaling. If scaled=TRUE and DistConf=NULL, it will be computed with the function predictConfModels, which takes some time and the function will throw a message. Default: NULL

## Details

These functions merely split the input data frame by model participants combinations, call the equivalent `predictRTConf` functions for the individual parameter sets and bind the outputs together. They are included for convenience and the easy parallelization, which facilitates speeding up computations considerably. For the argument `paramDf`, the output of the fitting function `fitRTConfModels` with the respective models and participants may be used.

The function `predictConf` (called by `predictConfModels`) consists merely of an integration of the reaction time density or the given model, `{d*model*}`, over the reaction time in a reasonable interval (0 to `maxrt`). The function `predictRT` (called by `predictRTModels`) wraps these density functions to a parameter set input and a data.frame output. ' Note, that the encoding for stimulus identity is different between diffusion based models (2DSD, dynWEV) and race models (IRM(t), PCRM(t), and MTLNR). Therefore, in the columns stimulus and response there will be a mix of encodings: -1/1 for diffusion based models and 1/2 for race models. This, usually is not important, since for further aggregation models will not be mixed.

## Value

`predictConfModels` returns a data.frame/tibble with columns: participant (or sbj, subject depending on the input), model, condition, stimulus, response, rating, correct, p, info, err. p is the predicted probability of a response and rating, given the stimulus category and condition. info and err refer to the respective outputs of the integration routine used for the computation. `predictRTModels` returns a data.frame/tibble with columns: participant (or sbj/subject depending on the input), model, condition, stimulus, response, rating, correct, rt and dens (and densscaled, if scaled=TRUE).

## Note

Different parameters for different conditions are only allowed for drift rate  $v$ , drift rate variability  $sv$  (only dynWEV and 2DSD), and process variability  $s$  (not for MTLNR). All other parameters are used for all conditions.

## Author(s)

Sebastian Hellmann.

## Examples

```
# First example for 2 participant and the "dynWEV" model
# (equivalent applicable for
# all other models (with different parameters!))
# 1. Define two parameter sets from different participants
paramDf <- data.frame(participant = c(1,2), model="dynWEV",
                      a=c(1.5, 2),v1=c(0.2,0.1), v2=c(1, 1.5),
                      t0=c(0.1, 0.2),z=c(0.52,0.45),
                      sz=c(0.0,0.3),sv=c(0.4,0.7), st0=c(0,0.01),
                      tau=c(2,3), w=c(0.5,0.2),
                      theta1=c(1,1.5), svis=c(0.5,0.1), sigvis=c(0.8, 1.2))

paramDf
# 2. Predict discrete Choice x Confidence distribution:
# model is not an extra argument but must be a column of paramDf
preds_Conf <- predictConfModels(paramDf, maxrt = 15, simult_conf=TRUE,
                               .progress=TRUE, parallel = FALSE)

# 3. Compute RT density
preds_RT <- predictRTModels(paramDf, maxrt=6, subdivisions=100,
                           scaled=TRUE, DistConf = preds_Conf,
                           parallel=FALSE, .progress = TRUE)

head(preds_RT)

# produces a warning, if scaled=TRUE and DistConf missing
preds_RT <- predictRTModels(paramDf, scaled=TRUE)

# Use PDFtoQuantiles to get predicted RT quantiles
head(PDFtoQuantiles(preds_RT, scaled = FALSE))

# Second Example: only one parameter set but for two different models

paramDf1 <- data.frame(model="dynWEV", a=1.5,v1=0.2, v2=1, t0=0.1,z=0.52,
                      sz=0.3,sv=0.4, st0=0, tau=3, w=0.5,
                      theta1=1, svis=0.5, sigvis=0.8)
paramDf2 <- data.frame(model="PCRMt", a=2,b=2, v1=0.5, v2=1, t0=0.1,st0=0,
                      wx=0.6, wint=0.2, wrt=0.2, theta1=4)
paramDf <- dplyr::full_join(paramDf1, paramDf2)
paramDf # each model parameters sets hat its relevant parameters
predictConfModels(paramDf, parallel=FALSE, .progress=TRUE)
```

---

predictWEV

*Prediction of Confidence Rating and Response Time Distribution in  
dynaViTE, dynWEV, and 2DSD confidence models*

---

## Description

predictWEV\_Conf predicts the categorical response distribution of decision and confidence ratings, predictWEV\_RT computes the predicted RT distribution (density) in the 2DSD Model (Pleskac & Busemeyer, 2010) and the dynWEV model (Hellmann et al., 2023), given specific parameter constellations. See [ddynaViTE](#) and [d2DSD](#) for more information about parameters.

**Usage**

```

predictWEV_Conf(paramDf, model = "dynaViTE", maxrt = Inf,
  subdivisions = 100L, simult_conf = FALSE, stop.on.error = FALSE,
  precision = 3, .progress = TRUE)

predictWEV_RT(paramDf, model = NULL, maxrt = 9, subdivisions = 100L,
  minrt = NULL, simult_conf = FALSE, scaled = FALSE, DistConf = NULL,
  precision = 3, .progress = TRUE)

```

**Arguments**

paramDf	a list or dataframe with one row. Column names should match the names of <a href="#">dynaViTE</a> and <a href="#">2DSD</a> model specific parameter names. For different stimulus quality/mean drift rates, names should be v1, v2, v3,... Different sv and/or s parameters are possible with sv1, sv2, sv3... (s1, s2, s3,... respectively) with equally many steps as for drift rates. Additionally, the confidence thresholds should be given by names with thetaUpper1, thetaUpper2,..., thetaLower1,... or, for symmetric thresholds only by theta1, theta2,....
model	character scalar. One of "dynaViTE", "dynWEV", or "2DSD".
maxrt	numeric. The maximum RT for the integration/density computation. Default: 15 (for predictWEV_Conf (integration)), 9 (for predictWEV_RT).
subdivisions	integer (default: 100). For predictWEV_Conf it is used as argument for the inner integral routine. For predictWEV_RT it is the number of points for which the density is computed.
simult_conf	logical. Whether in the experiment confidence was reported simultaneously with the decision, as then decision and confidence judgment are assumed to have happened subsequent before response and computations are different, when there is an observable interjudgment time (then simult_conf should be FALSE).
stop.on.error	logical. Argument directly passed on to integrate. Default is FALSE, since the densities invoked may lead to slow convergence of the integrals (which are still quite accurate) which causes R to throw an error.
precision	numerical scalar value. Precision of calculation. Corresponds to the step size of integration w.r.t. z and t0. Default is 1e-5.
.progress	logical. if TRUE (default) a progress bar is drawn to the console.
minrt	numeric or NULL(default). The minimum rt for the density computation.
scaled	logical. For predictWEV_RT. Whether the computed density should be scaled to integrate to one (additional column densscaled). Otherwise the output contains only the defective density (i.e. its integral is equal to the probability of a response and not 1). If TRUE, the argument DistConf should be given, if available. Default: FALSE.
DistConf	NULL or data.frame. A data.frame or matrix with column names, giving the distribution of response and rating choices for different conditions and stimulus categories in the form of the output of predictWEV_Conf. It is only necessary, if scaled=TRUE, because these probabilities are used for scaling. If scaled=TRUE and DistConf=NULL, it will be computed with the function predictWEV_Conf, which takes some time and the function will throw a message. Default: NULL

## Details

The function `predictWEV_Conf` consists merely of an integration of the response time density, `ddynaViTE` and `d2DSD`, over the response time in a reasonable interval (`t0` to `maxrt`). The function `predictWEV_RT` wraps these density functions to a parameter set input and a `data.frame` output. For the argument `paramDf`, the output of the fitting function `fitRTConf` with the respective model may be used.

## Value

`predictWEV_Conf` returns a `data.frame/tibble` with columns: `condition`, `stimulus`, `response`, `rating`, `correct`, `p`, `info`, `err`. `p` is the predicted probability of a response and rating, given the stimulus category and condition. `info` and `err` refer to the respective outputs of the integration routine used for the computation. `predictWEV_RT` returns a `data.frame/tibble` with columns: `condition`, `stimulus`, `response`, `rating`, `correct`, `rt` and `dens` (and `densscaled`, if `scaled=TRUE`).

## Note

Different parameters for different conditions are only allowed for drift rate `v`, drift rate variability `sv`, and process variability `s`. Otherwise, `s` is not required in `paramDf` but set to 1 by default. All other parameters are used for all conditions.

## Author(s)

Sebastian Hellmann.

## References

- Hellmann, S., Zehetleitner, M., & Rausch, M. (2023). Simultaneous modeling of choice, confidence and response time in visual perception. *Psychological Review* 2023 Mar 13. doi: 10.1037/rev0000411. Epub ahead of print. PMID: 36913292.
- Pleskac, T. J., & Busemeyer, J. R. (2010). Two-Stage Dynamic Signal Detection: A Theory of Choice, Decision Time, and Confidence, *Psychological Review*, 117(3), 864-901. doi:10.1037/a0019737

## Examples

```
# Examples for "dynWEV" model (equivalent applicable for "2DSD" model (with less parameters))
# 1. Define some parameter set in a data.frame
paramDf <- data.frame(a=2.5, v1=0.5, v2=1, t0=0.1, z=0.55,
                      sz=0, sv=0.2, st0=0, tau=3, w=0.3,
                      theta1=0.8, svis=0.5, sigvis=0.8)

# 2. Predict discrete Choice x Confidence distribution:
preds_Conf <- predictWEV_Conf(paramDf, "dynWEV", maxrt = 15)
head(preds_Conf)

# To set simult_conf=TRUE makes a minor difference in the discrete distribution,
# because we integrate over response times (we just adapt maxrt for comparison)
preds_Conf2 <- predictWEV_Conf(paramDf, "dynWEV", simult_conf = TRUE, maxrt = 15+paramDf$tau)
summary(preds_Conf$p-preds_Conf2$p) # difference in predicted probabilities
```

```

# 3. Compute RT density
preds_RT <- predictWEV_RT(paramDf, "dynWEV", maxrt=4, subdivisions=200) #(scaled=FALSE)
# same output with scaled density column:
preds_RT <- predictWEV_RT(paramDf, "dynWEV", maxrt=4, subdivisions=200,
                          scaled=TRUE, DistConf = preds_Conf)

head(preds_RT)

# produces a warning, if scaled=TRUE and DistConf missing
preds_RT <- predictWEV_RT(paramDf, "dynWEV", maxrt=4, subdivisions=200,
                          scaled=TRUE)

# Example of visualization
library(ggplot2)
preds_Conf$rating <- factor(preds_Conf$rating, labels=c("unsure", "sure"))
preds_RT$rating <- factor(preds_RT$rating, labels=c("unsure", "sure"))
ggplot(preds_Conf, aes(x=interaction(rating, response), y=p))+
  geom_bar(stat="identity")+
  facet_grid(cols=vars(stimulus), rows=vars(condition), labeller = "label_both")
ggplot(preds_RT, aes(x=rt, color=interaction(rating, response), y=dens))+
  geom_line(stat="identity")+
  facet_grid(cols=vars(stimulus), rows=vars(condition), labeller = "label_both")+
  theme(legend.position = "bottom")
ggplot(aggregate(densscaled~rt+correct+rating+condition, preds_RT, mean),
      aes(x=rt, color=rating, y=densscaled))+
  geom_line(stat="identity")+
  facet_grid(cols=vars(condition), rows=vars(correct), labeller = "label_both")+
  theme(legend.position = "bottom")

# Use PDFtoQuantiles to get predicted RT quantiles
head(PDFtoQuantiles(preds_RT, scaled = FALSE))

```

## Description

`subject_modelweights` computes the model weights (as probabilities) of individual subjects based on an information criterion (BIC, AIC, or AICc). `group_BMS` performs a Bayesian model comparison based on marginal likelihoods (alias model evidence), given for different models across different subject on a group level using a fixed effects model and a random effects model on the distribution of model probabilities (see Rigoux et al., 2014 and Details section). `group_BMS_fits` is a wrapper for `group_BMS` that can be used with the output of `fitRTConfModels`, i.e. a data frame with information criteria for different models and subjects, using an information criterion to approximate the model evidence.

**Usage**

```

subject_modelweights(fits, measure = "BIC")

group_BMS_fits(fits, measure = "BIC", opts = list(), alpha0 = NULL)

group_BMS(mlp, opts = list(), alpha0 = NULL)

```

**Arguments**

<code>fits</code>	a data frame as returned by <code>fitRTConfModels</code> . Should contain a column <code>model</code> indicating the model name, a column <code>subject</code> (alternatively <code>sbj</code> or <code>participant</code> ) indicating the grouping structure of the data, and a column with the name given by the <code>measure</code> argument containing the values of the information criterion that should be used to approximate model evidence.
<code>measure</code>	the name of the column indicating the information criterion to approximate model evidence. For outputs of <code>fitRTConfModels</code> , the available measures are 'BIC', 'AIC', and 'AICc'. Any other approximation for the model evidence may be used, the measure is transferred to log model evidence by taking $-\text{measure}/2$ .
<code>opts</code>	a list with options for the iteration algorithm to estimate the parameter of the Dirichlet distribution. Following values may be provided: <ul style="list-style-type: none"> <li>• <code>maxiter</code> the maximum number of iterations (Default: 200)</li> <li>• <code>tol</code> the tolerance for changes in the free energy approximation to stop the algorithm, if <math> \text{FE}(i+1) - \text{FE}(i)  &lt; \text{tol}</math> the algorithm is stopped (Default: <math>1e-4</math>)</li> <li>• <code>eps</code> The number to substitute values of 0 in calls to log (Default: <math>1e-32</math>)</li> </ul>
<code>alpha0</code>	a positive numeric vector representing the parameter of a Dirichlet distribution used as prior over model probabilities. The length should be equal to <code>nrow(mlp)</code> for <code>group_BMS</code> , and equal to the number of unique names in the <code>model</code> column of <code>fits</code> for <code>group_BMS_fits</code> .
<code>mlp</code>	a matrix containing the logarithm of marginal probabilities (i.e. log model evidence) with N columns representing individuals (or any other grouping structure) and K rows representing the models.

**Details**

This set of function can be used for model comparisons on a group level when the models were not fitted hierarchical but by fitting the models independently to different subgroups (e.g. data from different subjects).

The function `subject_modelweights` computes the model weights for each subject separately to inspect predominant models but also heterogeneity within the population. The functions `group_BMS` and `group_BMS_fits` can be used for a Bayesian model selection on the group level following the approach of Rigoux et al. (2014). The approach compares three different models for the generative structure of the data and gives estimates for model probabilities for the fixed and random effects models. The **fixed effects model** assumes that there is a single model that generated the data of all subjects. Thus, model weights may be computed directly by multiplying the model weights computed on a subject-level. This model is formulated in a Bayesian way using a Multinomial distribution over the models as prior with some prior parameter `alpha0` giving the prior model weights.

This is updated according to the marginal model likelihoods resulting in a single posterior vector of model probabilities, reported in the column `fx_prob` of the `model_weights` data frame. The random effects model assumes that there is a vector of model probabilities and each subject may be generated by a different model, each drawn from a Multinomial distribution. The Bayesian prior on this vector of model probabilities is given by a Dirichlet distribution with some parameter  $\alpha_0$ . The function uses a variational technique to approximate the alpha parameter of the posterior Dirichlet distribution. Within this framework several statistics may be used for model selection. The `model_weights` data frame reports the posterior alpha parameter, as well as the posterior mean  $r$  of the corresponding dirichlet distribution. The exceedance probability  $ep$  represents the probability that given a random sample from the Dirichlet distribution the probability of the model is greater than all other probabilities. Finally, the protected exceedance probability ( $pep$ ) is a scaled version of the  $ep$  multiplying the  $ep$  by one minus the Bayesian omnibus risk (BOR). The Bayesian omnibus risk is the posterior probability of the **null model** against the random effects model. The **null model** assumes that all models are generating the subjects' data with equal probability and results from taking the limit of  $\alpha_0$  towards infinity. The Bayesian omnibus risk is reported in the `summary_stats` together with the free energy approximation of the null, the fixed effects, and the random effects models.

### Value

`subject_modelweights` returns a data frame of subject-wise model probabilities with rows for each subject and columns for the models given by name and one column for the subject ID as given in the input. `group_BMS` and `group_BMS_fits` return a list with two entries:

- `model_weights`: a matrix with rows for each model (row names indicate the model names for `group_BMS_fits` and for `group_BMS` if row names are available in `mlp`), and following columns:  $\alpha$  (the alpha parameter of the Dirichlet posterior over model probabilities in the population),  $r$  (the mean probabilities of each model in the population),  $ep$  and  $pep$  (exceedance and protected exceedance probabilities for each model), and `fx_prob` (the posterior model probabilities if a fixed true model is assumed in the population).
- `summary_stats`: a vector giving statistics for the Bayesian model comparison that may be used for other analyses: Bayesian omnibus risks: `bor` (random effects model against the null model), `bor_fixed` (fixed effects model against the null model), and `bor_re_fixed` (random effects model against the fixed effects model) (higher values indicating a simpler model), and estimations of the Free Energy of the Dirichlet distribution `FE` (random effects model), `FE0` (null model), and `FEfixed` (fixed effects model)

### Author(s)

Sebastian Hellmann.

### References

Rigoux, L., Stephan, K. E., Friston, K. J., & Daunizeau, J. (2014). Bayesian model selection for group studies - revisited. *NeuroImage*, 84, 971–985. doi: 10.1016/j.neuroimage.2013.08.065

### Examples

```
# Define a data frame with information criteria from model fits
# (this is a sub-data.frame from an output of fitRTConfModels with
```

```

# 8 subjects, three models and rounded information criteria)
fits <- data.frame(
  participant = rep(1:8, each=3),
  model = rep(c("dynaViTE", "2DSD", "PCRMt"), 8),
  BIC = c(5318, 5665, 1659, 3856, 5508, 3982, 3950, 3998,
          4114, 4216, 4314, 4419, 3170, 3489, 3256, 1950,
          1934, 2051, 3194, 3317, 3359, 9656, 10161, 4024),
  AIC = c(5211, 5577, 1577, 3750, 5420, 3899, 3843, 3911,
          4031, 4109, 4226, 4337, 3063, 3401, 3173, 1844,
          1847, 1969, 3087, 3229, 3277, 9549, 10074, 3942),
  AICc = c(5212, 5578, 1577, 3751, 5421, 3900, 3844, 3911,
           4032, 4110, 4227, 4337, 3064, 3402, 3174, 1845,
           1848, 1970, 3088, 3230, 3277, 9550, 10074, 3942))
# Compute subject-wise model probabilities based on different ICs
subject_modelweights(fits, measure = "BIC")
subject_modelweights(fits, measure = "AIC")
subject_modelweights(fits, measure = "AICc")
# Conduct group-level Bayesian model selection based on BIC
group_BMS_fits(fits, measure="BIC")

## General group-level Bayesian model selection based on any marginal log-probabilities
# Compute marginal log-likelihood based on BIC from fits
mlp <- matrix(NA, ncol=8, nrow=3)
for (i in 1:8) mlp[,i] <- fits[(i-1)*3 + 1:3, "BIC"]
mlp <- - mlp/(2)
rownames(mlp) <- c("dynaViTE", "2DSD", "PCRMt")
# conduct group BMS:
group_BMS(mlp)

```

---

RaceModels

---

*Independent and partially anti-correlated Race Model for Decision Confidence*


---

## Description

Probability densities and random number generators for response times, decisions and confidence judgments in the independent Race Model (dIRM/rIRM) or partially (anti-)correlated Race Model (dPCRM/rPCRM), i.e. the probability of a given response (response: winning accumulator (1 or 2)) at a given time (rt) and the confidence measure in the interval between th1 and th2 (Hellmann et al., 2023). The definition of the confidence measure depends on the argument `time_scaled` (see Details). The computations are based on Moreno-Bote (2010). The parameters for the models are  $\mu_1$  and  $\mu_2$  for the drift rates,  $a$ ,  $b$  for the upper thresholds of the two accumulators and  $s$  for the incremental standard deviation of the processes and  $t_0$  and  $st_0$  for the minimum and range of uniformly distributed non-decision times (including encoding and motor time). For the computation of confidence judgments, the parameters  $th_1$  and  $th_2$  for the lower and upper bound of the interval for confidence measure and if `time_scaled` is TRUE the weight parameters  $w_x$ ,  $w_{rt}$ ,  $w_{int}$  for the computation of the confidence measure are required (see Details).

**Usage**

```

dIRM(rt, response = 1, mu1, mu2, a, b, th1, th2, wx = 1, wrt = 0,
     wint = 0, t0 = 0, st0 = 0, s1 = 1, s2 = 1, s = NULL,
     time_scaled = TRUE, precision = 6, step_width = NULL)

dIRM2(rt, response = 1, mu1, mu2, a, b, th1, th2, wx = 1, wrt = 0,
     wint = 0, t0 = 0, st0 = 0, s1 = 1, s2 = 1, smu1 = 0, smu2 = 0,
     sza = 0, szb = 0, s = NULL, time_scaled = TRUE, precision = 6,
     step_width = NULL)

dIRM3(rt, response = 1, mu1, mu2, a, b, th1, th2, wx = 1, wrt = 0,
     wint = 0, t0 = 0, st0 = 0, s1 = 1, s2 = 1, smu1 = 0, smu2 = 0,
     s = NULL, time_scaled = TRUE, precision = 6, step_width = NULL)

dPCRM(rt, response = 1, mu1, mu2, a, b, th1, th2, wx = 1, wrt = 0,
     wint = 0, t0 = 0, st0 = 0, s1 = 1, s2 = 1, s = NULL,
     time_scaled = TRUE, precision = 6, step_width = NULL)

rIRM(n, mu1, mu2, a, b, wx = 1, wrt = 0, wint = 0, t0 = 0, st0 = 0,
     s1 = 1, s2 = 1, s = NULL, smu1 = 0, smu2 = 0, sza = 0, szb = 0,
     time_scaled = TRUE, delta = 0.01, maxrt = 15)

rPCRM(n, mu1, mu2, a, b, wx = 1, wrt = 0, wint = 0, t0 = 0, st0 = 0,
     s1 = 1, s2 = 1, s = NULL, smu1 = 0, smu2 = 0, sza = 0, szb = 0,
     time_scaled = TRUE, delta = 0.01, maxrt = 15)

```

**Arguments**

<code>rt</code>	a numeric vector of RTs. For convenience also a <code>data.frame</code> with columns <code>rt</code> and <code>response</code> is possible.
<code>response</code>	numeric vector with values in <code>c(1, 2)</code> , giving the accumulator that hit its boundary first.
<code>mu1</code>	numeric. Drift rate for the first accumulator
<code>mu2</code>	numeric. Drift rate for the second accumulator
<code>a</code>	positive numeric. Distance from starting point to boundary of the first accumulator.
<code>b</code>	positive numeric. Distance from starting point to boundary of the second accumulator.
<code>th1</code>	numeric. Lower bound of interval range for the confidence measure.
<code>th2</code>	numeric. Upper bound of interval range for the confidence measure.
<code>wx</code>	numeric. Weight on losing accumulator for the computation of the confidence measure. (Used only if <code>time_scale=TRUE</code> , 1)
<code>wrt</code>	numeric. Weight on reaction time for the computation of the confidence measure. (Used only if <code>time_scale=TRUE</code> , Default 0)

wint	numeric. Weight on the interaction of losing accumulator and reaction time for the computation of the confidence measure. (Used only if time_scale=TRUE, Default 0)
t0	numeric. Lower bound of non-decision time component in observable response times. Range: $t_0 \geq 0$ . Default: 0.
st0	numeric. Range of a uniform distribution for non-decision time. Range: $st_0 \geq 0$ . Default: 0.
s1	numeric. Diffusion constant of the first accumulator. Usually fixed to 1 for most purposes as it scales other parameters (see Details). Range: $s_1 > 0$ , Default: 1.
s2	numeric. Diffusion constant of the second accumulator. Usually fixed to 1 for most purposes as it scales other parameters (see Details). Range: $s_2 > 0$ , Default: 1.
s	numeric. Alternative way to specify diffusion constants, if both are assumed to be equal. If both ( $s_1$ , $s_2$ and $s$ ) are given, only $s_1$ and $s_2$ will be used.
time_scaled	logical. Whether the confidence measure should be time-dependent. See Details.
precision	numerical scalar value. Precision of calculation. Determines the step size of integration w.r.t. $t_0$ . Represents the number of decimals precisely computed on average. Default is 6.
step_width	numeric. Alternative way to define the precision of integration w.r.t. $t_0$ by directly providing the step size for the integration.
smu1	numeric. Between-trial variability in the drift rate of the first accumulator.
smu2	numeric. Between-trial variability in the drift rate of the second accumulator.
sza	numeric. Between-trial variability in starting point of the first accumulator.
szb	numeric. Between-trial variability in starting point of the second accumulator.
n	integer. The number of samples generated.
delta	numeric. Discretization step size for simulations in the stochastic process
maxrt	numeric. Maximum decision time returned. If the simulation of the stochastic process exceeds a decision time of $maxrt$ , the response will be set to 0 and the $maxrt$ will be returned as $rt$ .

## Details

The parameters are formulated, s.t. both accumulators start at 0 and trigger a decision at their positive boundary  $a$  and  $b$  respectively. That means, both parameters have to be positive. Internally the computations adapt the parametrization of Moreno-Bote (2010).

`time_scaled` determines whether the confidence measure is computed in accordance to the Balance of Evidence hypothesis (if `time_scaled=FALSE`), i.e. if response is 1 at time  $T$  and  $X_2$  is the second accumulator, then

$$conf = b - X_2(T)$$

. Otherwise, if `time_scaled=TRUE` (default), confidence is computed as linear combination of Balance of Evidence, decision time, and an interaction term, i.e.

$$conf = wx(b - X_2(T)) + wrt \frac{1}{\sqrt{T}} + wint \frac{b - X_2(T)}{\sqrt{T}}.$$

Usually the weights ( $w_x$ ,  $w_{rt}$ ,  $w_{int}$ ) should sum to 1, as the confidence thresholds ( $th1$  and  $th2$ ) may be scaled according to their sum. If this is not the case, they will be scaled accordingly internally! Usually, this formula results in lower confidence when the reaction time is longer but the state of the second accumulator is held constant. It is based on the optimal decision confidence in Moreno-Bote (2010).

For convenience, the likelihood function allows that the first argument is a `data.frame` containing the information of the first and second argument in the columns (i.e., `rt` and `response`). Other columns (as well as passing response separately as argument) will be ignored.

The simulations are done by simulating normal variables in discretized steps until one process reaches the boundary. If no boundary is met within the maximum time, response is set to 0.

### Value

`dIRM` and `dPCRM` return the numerical value of the probability density in a numerical vector of the same length as `rt`.

`rIRM` and `dPCRM` return a `data.frame` with four columns and `n` rows. Column names are `rt` (response time), `response` (1 or 2, indicating which accumulator hit its boundary first), `x1` (the final state of the loosing accumulator), and `conf` (the value of the confidence measure; not discretized!).

The race parameters (as well as `response`, `th1`, and `th2`) are recycled to the length of the result (either `rt` or `n`). In other words, the functions are completely vectorized for all parameters and even the response.

### Note

Similarly to the drift diffusion models (like `ddiffusion` and `ddynaViTE`), `s1` and `s2` are scaling factors (`s1` scales:  $\mu_1$  and  $a$ , `s2` scales:  $\mu_2$  and  $b$ , and depending on response: if `response=2`, `s1` scales  $th1$ ,  $th2$ , and  $w_{rt}$ ), otherwise `s2` is the scaling factor. It is sometimes assumed (Moreno-Bote, 2010), that both noise terms are equal, then they should definitely be fixed for fitting.

### Author(s)

Sebastian Hellmann

### References

- Hellmann, S., Zehetleitner, M., & Rausch, M. (2023). Simultaneous modeling of choice, confidence and response time in visual perception. *Psychological Review* 2023 Mar 13. doi: 10.1037/rev0000411. Epub ahead of print. PMID: 36913292.
- Moreno-Bote, R. (2010). Decision confidence and uncertainty in diffusion models with partially correlated neuronal integrators. *Neural Computation*, 22(7), 1786–1811. doi:10.1162/neco.2010.12-08-930

### Examples

```
# Plot rt distribution ignoring confidence
curve(dPCRM(x, 1, mu1=0.5, mu2=-0.5, a=1, b=1, th1=-Inf, th2=Inf, t0=0.1), xlim=c(0,2.5))
curve(dPCRM(x, 2, mu1=0.5, mu2=-0.5, a=1, b=1, th1=-Inf, th2=Inf, t0=0.1), col="red", add=TRUE)
curve(dIRM(x, 1, mu1=0.5, mu2=-0.5, a=1, b=1, th1=-Inf, th2=Inf, t0=0.1), lty=2, add=TRUE)
```

```

curve(dIRM(x, 2, mu1=0.5, mu2=-0.5, a=1, b=1, th1=-Inf, th2=Inf, t0=0.1),
      col="red", lty=2, add=TRUE)
# t0 indicates minimal response time possible
abline(v=0.1)
## Following example may be equivalently used for the IRM model functions.
# Generate a random sample
df1 <- rPCRM(5000, mu1=0.2, mu2=-0.2, a=1, b=1, t0=0.1,
            wx = 1) # Balance of Evidence
# Same RT and response distribution but different confidence distribution
df2 <- rPCRM(5000, mu1=0.2, mu2=-0.2, a=1, b=1, t0=0.1,
            wint = 0.2, wrt=0.8)
head(df1)

# Compute density with rt and response as separate arguments
dPCRM(seq(0, 2, by =0.4), response= 2, mu1=0.2, mu2=-0.2, a=1, b=1, th1=0.5,
      th2=2, wx = 0.3, wint=0.4, wrt=0.1, t0=0.1)
# Compute density with rt and response in data.frame argument
df1 <- subset(df1, response !=0) # drop trials where no accumulation hit its boundary
dPCRM(df1[1:5,], mu1=0.2, mu2=-0.2, a=1, b=1, th1=0, th2=Inf, t0=0.1)
# s1 and s2 scale other decision relevant parameters
s <- 2 # common (equal) standard deviation
dPCRM(df1[1:5,], mu1=0.2*s, mu2=-0.2*s, a=1*s, b=1*s, th1=0, th2=Inf, t0=0.1, s1=s, s2=s)
s1 <- 2 # different standard deviations
s2 <- 1.5
dPCRM(df1[1:5,], mu1=0.2*s1, mu2=-0.2*s2, a=1*s1, b=1*s2, th1=0, th2=Inf, t0=0.1, s1=s1, s2=s2)

# s1 and s2 scale also confidence parameters
df1[1:5,]$response <- 2 # set response to 2
# for confidence it is important to scale confidence parameters with
# the right variation parameter (the one of the loosing accumulator)
dPCRM(df1[1:5,], mu1=0.2, mu2=-0.2, a=1, b=1,
      th1=0.5, th2=2, wx = 0.3, wint=0.4, wrt=0.1, t0=0.1)
dPCRM(df1[1:5,], mu1=0.2*s1, mu2=-0.2*s2, a=1*s1, b=1*s2,
      th1=0.5, th2=2, wx = 0.3/s1, wint = 0.4/s1, wrt = 0.1, t0=0.1, s1=s1, s2=s2)
dPCRM(df1[1:5,], mu1=0.2*s1, mu2=-0.2*s2, a=1*s1, b=1*s2,
      th1=0.5*s1, th2=2*s1, wx = 0.3, wint = 0.4, wrt = 0.1*s1, t0=0.1, s1=s1, s2=s2)

two_samples <- rbind(cbind(df1, ws="BoE"),
                    cbind(df2, ws="RT"))
# drop not finished decision processes
two_samples <- two_samples[two_samples$response!=0,]
# no difference in RT distributions
boxplot(rt~ws+response, data=two_samples)
# but different confidence distributions
boxplot(conf~ws+response, data=two_samples)
if (requireNamespace("ggplot2", quietly = TRUE)) {
  require(ggplot2)
  ggplot(two_samples, aes(x=rt, y=conf, fill = after_stat(density)))+
    stat_density_2d(geom="raster", contour=FALSE, na.rm=TRUE)+
    xlim(c(0.2, 1.3))+ ylim(c(0, 2.5))+
    facet_grid(cols=vars(ws), rows=vars(response), labeller = "label_both")
}

```

```
# Restricting to specific confidence region
df1 <- df1[df1$conf >0 & df1$conf <1,]
dPCRm(df1[1:5,], th1=0, th2=1, mu1=0.2, mu2=-0.2, a=1, b=1, t0=0.1, wx = 1 )
```

---

rLCA	<i>Simulation of confidence ratings and RTs in leaky competing accumulator model</i>
------	--

---

## Description

Simulates the decision responses, reaction times and state of the loosing accumulator together with a confidence measure in the leaky competing accumulator model. Optionally, there is a post-decisional accumulation period, where the processes continues.

## Usage

```
rLCA(n, mu1, mu2, th1, th2, k = 0, beta = 0, SPV = 0, tau = 0,
     wx = 1, wrt = 0, wint = 0, t0 = 0, st0 = 0, pi = 0, sig = 1,
     time_scaled = TRUE, simult_conf = FALSE, delta = 0.01, maxrt = 15)
```

## Arguments

n	integer. number of samples.
mu1	mean momentary evidence for alternative 1
mu2	mean momentary evidence for alternative 2
th1	decision threshold for alternative 1
th2	decision threshold for alternative 2
k	leakage (default: 0)
beta	inhibition (default: 0)
SPV	variation in starting points (default: 0)
tau	fixed post decisional accumulation period (default: 0)
wx	weight on balance of evidence in confidence measure (default: 1)
wrt	weight on RT in confidence measure (default: 0)
wint	weight on interaction of evidence and RT in confidence measure (default: 0)
t0	minimal non-decision time (default: 0)
st0	range of uniform distribution of non-decision time (default: 0)
pi	factor for input dependent noise of infinitesimal variance of processes (default: 0)
sig	input independent component of infinitesimal variance of processes (default: 1)
time_scaled	logical. Whether a time_scaled transformation for the confidence measure should be used.

<code>simult_conf</code>	logical. Whether in the experiment confidence was reported simultaneously with the decision. If that is the case decision and confidence judgment are assumed to have happened subsequent before the response. Therefore <code>tau</code> is included in the response time. If the decision was reported before the confidence report, <code>simult_conf</code> should be FALSE.
<code>delta</code>	numerical. Size of steps for the discretized simulation (see details).
<code>maxrt</code>	numerical. Maximum reaction time to be simulated (see details). Default: 15.

### Details

The simulation is done by simulating discretized steps until one process reaches the boundary with an update rule:

$$\delta X_i(t) = \max(0, X_i(t) + \delta_t((k-1)X_i(t) - \beta X_{j=i}(t) + \mu_i + \varepsilon_i(t)),$$

with  $\varepsilon_i(t) \sim N(0, (\pi\mu_i)^2 + \sigma^2)$ . If no boundary is met within the maximum time, response is set to 0. After the decision, the accumulation continues for a time period (`tau`), until the final state is used for the computation of confidence.

### Value

Returns a `data.frame` with three columns and `n` rows. Column names are `rt` (response time), `response` (1 or 2, indicating which accumulator hit its boundary first), and `conf` (the value of the confidence measure; not discretized!).

### Author(s)

Sebastian Hellmann.

### Examples

```
# minimal arguments
simus<- rLCA(n=20, mu1=1, mu2=-0.5, th1=1, th2=0.8)
head(simus)

# specifying all relevant parameters
simus <- rLCA(n=1000, mu1 = 2.5, mu2=1, th1=1.5, th2=1.6,
             k=0.1, beta=0.1, SPV=0.2, tau=0.1,
             wx=0.8, wrt=0.2, wint=0, t0=0.2, st0=0.1,
             pi=0.2, sig=1)
if (requireNamespace("ggplot2", quietly = TRUE)) {
  if (requireNamespace("MASS", quietly = TRUE)) {
    require(MASS)
    require(ggplot2)
    ggplot(simus, aes(x=rt, y=conf))+
      geom_bin2d()+
      facet_wrap(~response)
  }
}
boxplot(conf~response, data=simus)
```

---

simulateMTLNR	<i>Simulation of confidence ratings and RTs in the Multiple-threshold Log-normal Race Model</i>
---------------	---

---

## Description

Simulates the decision responses and reaction times together with a discrete confidence judgment in the MTLNR (Reynolds et al., 2020), given specific parameter constellations. See [dMTLNR](#) for more information about parameters. Also computes the Gamma rank correlation between the confidence ratings and condition (task difficulty), reaction times and accuracy in the simulated output. Basically, this function is a wrapper for [rMTLNR](#) for application in confidence experiments with manipulation of specific parameters.

## Usage

```
simulateMTLNR(paramDf, n = 10000, gamma = FALSE, agg_simus = FALSE,
  stimulus = c(1, 2), seed = NULL)
```

## Arguments

paramDf	a list or data frame with one row. Column names should match the following names (see <a href="#">dMTLNR</a> ): For different stimulus quality/mean drift rates, names should be v1, v2, v3,... (corresponding to the mean parameter for the accumulation rate for the stimulus-corresponding accumulator, therefore mu_v1 and mu_v2 are not used in this context but replaced by the parameter v); mu_d1 and mu_d2 correspond to the mean parameters for boundary distance of the two accumulators; s1 and s2 correspond to the variance parameters of the first and second boundary hitting time; rho corresponds to the correlation of boundary hitting times. Note that s_v1,s_v2,rho_v,s_d1,s_d2, and rho_d are not used in this context, although the accumulation rate-related parameters can be used to replace the above-mentioned variance parameters. Additionally, the confidence thresholds should be given by names with thetaUpper1, thetaUpper2,..., thetaLower1,... or, for symmetric thresholds only by theta1, theta2,... (see Details for the correspondence to the data)
n	integer. The number of samples (per condition and stimulus direction) generated. Total number of samples is n*nConditions*length(stimulus).
gamma	logical. If TRUE, the gamma correlation between confidence ratings, rt and accuracy is computed.
agg_simus	logical. Simulation is done on a trial basis with RTs outcome. If TRUE, the simulations will be aggregated over RTs to return only the distribution of response and confidence ratings. Default: FALSE.
stimulus	numeric vector. Either 1, 2 or c(1,2) (default). Together with condition represents the experimental situation. In a binary decision task the presented stimulus belongs to one of two categories. In the default setting trials with both categories presented are simulated but one can choose to simulate only trials with the stimulus coming from one category (1 for the category that is associated with

positive drift in the decision process where 1 responses are considered correct and 2 correspondingly for negative drifts and 2 correct decisions).

seed                      numerical. Seeding for non-random data generation.

## Details

Simulation is done by simulating normally distributed logarithms of boundary crossing times for both accumulators based on the MTLNR model. The smaller time determines decision time and response (i.e. the winning accumulator). The confidence variable is computed based on the log-ratio of the loosing boundary crossing time over the winning boundary crossing time.

The confidence values are then binned according to the given thresholds. The output of the fitting function `fitRTConf` with the respective model fits the argument `paramDf` for simulation. The Gamma coefficients are computed separately for correct/incorrect responses for the correlation of confidence ratings with condition and rt and separately for conditions for the correlation of accuracy and confidence. The resulting data frames in the output thus have two columns. One for the grouping variable and one for the Gamma coefficient.

## Value

Depending on `gamma` and `agg_simus`.

If `gamma` is FALSE, returns a `data.frame` with columns: `condition`, `stimulus`, `response`, `correct`, `rt`, `conf` (the continuous confidence measure) and `rating` (the discrete confidence rating), and `dec` and `vis` (only if `process_results=TRUE`) for the final states of accumulators in the simulation or (if `agg_simus=TRUE`): `condition`, `stimulus`, `response`, `correct`, `rating` and `p` (for the probability of a response and rating, given the condition and stimulus).

If `gamma` is TRUE, returns a list with elements: `simus` (the simulated data frame) and `gamma`, which is again a list with elements `condition`, `rt` and `correct`, each a tibble with two columns (see details for more information).

## Note

Different parameters for different conditions are only allowed for drift rate, `v`. All other parameters are used for all conditions.

## Author(s)

Sebastian Hellmann.

## References

Reynolds, A., Kvam, P. D., Osth, A. F., & Heathcote, A. (2020). Correlated racing evidence accumulator models. *Journal of Mathematical Psychology*, 96, 102331. doi: 10.1016/j.jmp.2020.102331

## Examples

```
# 1. Define some parameter set in a data.frame
paramDf <- data.frame(v1=0.5, v2=1.0, t0=0.1, st0=0,
                      mu_d1=1, mu_d2=1,
                      s1=0.5, s2=0.5, rho=0.2,
```

```

        theta1=0.8, theta2=1.5)

# 2. Simulate trials for both stimulus categories and all conditions (2)
simus <- simulateMTLNR(paramDf)
head(simus)

library(ggplot2)
simus <- simus[simus$response != 0, ]
simus$rating <- factor(simus$rating, labels = c("unsure", "medium", "sure"))
ggplot(simus, aes(x = rt, group = interaction(correct, rating),
  color = as.factor(correct), linetype = rating)) +
  geom_density(linewidth = 1.2, na.rm=TRUE) + xlim(c(0, 5)) +
  facet_grid(rows = vars(condition), labeller = "label_both")

# automatically aggregate simulation distribution
# to get only accuracy x confidence rating distribution for
# all conditions
agg_simus <- simulateMTLNR(paramDf, agg_simus = TRUE)
head(agg_simus)

agg_simus$rating <- factor(agg_simus$rating, labels = c("unsure", "medium", "sure"))
library(ggplot2)
ggplot(agg_simus, aes(x = rating, group = correct, fill = as.factor(correct), y = p)) +
  geom_bar(stat = "identity", position = "dodge") +
  facet_grid(cols = vars(condition), labeller = "label_both")

# Compute Gamma correlation coefficients between
# confidence and other behavioral measures
# output will be a list
simu_list <- simulateMTLNR(paramDf, n = 400, gamma = TRUE)
simu_list

# Example with asymmetric confidence thresholds
paramDf_asym <- data.frame(v1=0.5, v2=1.0, t0=0.1, st0=0,
  mu_d1=1, mu_d2=1,
  s1=0.5, s2=0.5, rho=0.2,
  thetaLower1=0.5, thetaLower2=1.2,
  thetaUpper1=0.7, thetaUpper2=1.8)

simus_asym <- simulateMTLNR(paramDf_asym, n = 1000)
head(simus_asym)

# Example with multiple conditions
paramDf_multi <- data.frame(v1=0.3, v2=0.6, v3=1.2, t0=0.1, st0=0,
  mu_d1=1, mu_d2=1,
  s1=0.5, s2=0.5, rho=0.2,
  theta1=0.8, theta2=1.5)

simus_multi <- simulateMTLNR(paramDf_multi, n = 1000)
table(simus_multi$condition, simus_multi$correct)

```

simulateRM

*Simulation of confidence ratings and RTs in race confidence models***Description**

Simulates the decision responses, reaction times and state of the loosing accumulator together with a discrete confidence judgment in the independent and partially anti-correlated race model (IRM and PCRM) (Hellmann et al., 2023), given specific parameter constellations. See [RaceModels](#) for more information about parameters. Also computes the Gamma rank correlation between the confidence ratings and condition (task difficulty), reaction times and accuracy in the simulated output. Basically, this function is a wrapper for [rIRM](#) and [rPCRM](#) for application in confidence experiments with manipulation of specific parameters. `rRM_Kiani` simulates a different version of race models, presented in Kiani et al. (2014), but without a confidence measure.

**Usage**

```
simulateRM(paramDf, n = 10000, model = "IRM", time_scaled = FALSE,
  gamma = FALSE, agg_simus = FALSE, stimulus = c(1, 2), delta = 0.01,
  maxrt = 15, seed = NULL)
```

```
rRM_Kiani(paramDf, n = 10000, time_scaled = FALSE, gamma = FALSE,
  agg_simus = FALSE, stimulus = c(1, 2), delta = 0.01, maxrt = 15,
  seed = NULL)
```

**Arguments**

<code>paramDf</code>	a list or data frame with one row. Column names should match the names of <a href="#">RaceModels</a> parameter names (only <code>mu1</code> and <code>mu2</code> are not used in this context but replaced by the parameter <code>v</code> ). For different stimulus quality/mean drift rates, names should be <code>v1</code> , <code>v2</code> , <code>v3</code> ,.... Different <code>s</code> parameters are possible with <code>s1</code> , <code>s2</code> , <code>s3</code> ,... with equally many steps as for drift rates. Additionally, the confidence thresholds should be given by names with <code>thetaUpper1</code> , <code>thetaUpper2</code> ,..., <code>thetaLower1</code> ,... or, for symmetric thresholds only by <code>theta1</code> , <code>theta2</code> ,....
<code>n</code>	integer. The number of samples (per condition and stimulus direction) generated. Total number of samples is <code>n*nConditions*length(stimulus)</code> .
<code>model</code>	character scalar. One of "IRM" or "PCRM". ("IRMt" and "PCRMt" will also be accepted. In that case, <code>time_scaled</code> is set to TRUE.)
<code>time_scaled</code>	logical. Whether a <code>time_scaled</code> transformation for the confidence measure should be used.
<code>gamma</code>	logical. If TRUE, the gamma correlation between confidence ratings, <code>rt</code> and accuracy is computed.
<code>agg_simus</code>	logical. Simulation is done on a trial basis with RTs outcome. If TRUE, the simulations will be aggregated over RTs to return only the distribution of response and confidence ratings. Default: FALSE.

stimulus	numeric vector. Either 1, 2 or c(1, 2) (default). Together with condition represents the experimental situation. In a binary decision task the presented stimulus belongs to one of two categories. In the default setting trials with both categories presented are simulated but one can choose to simulate only trials with the stimulus coming from one category (each associated with positive drift in one of two accumulators).
delta	numerical. Size of steps for the discretized simulation (see details).
maxrt	numerical. Maximum reaction time to be simulated (see details). Default: 15.
seed	numerical. Seeding for non-random data generation. (Also possible outside of the function.)

### Details

The simulation is done by simulating normal variables in discretized steps until one process reaches the boundary. If no boundary is met within the maximum time, response is set to 0. The output of the fitting function `fitRTConf` with the respective model fits the argument `paramDf` for simulation. The Gamma coefficients are computed separately for correct/incorrect responses for the correlation of confidence ratings with condition and rt and separately for conditions for the correlation of accuracy and confidence. The resulting data frames in the output thus have two columns. One for the grouping variable and one for the Gamma coefficient.

### Value

Depending on `gamma` and `agg_simus`.

If `gamma` is FALSE, returns a `data.frame` with columns: `condition`, `stimulus`, `response`, `correct`, `rt`, `conf` (the continuous confidence measure) and `rating` (the discrete confidence rating) or (if `agg_simus=TRUE`): `condition`, `stimulus`, `response`, `correct`, `rating` and `p` (for the probability of a response and rating, given the condition and stimulus).

If `gamma` is TRUE, returns a list with elements: `simus` (the simulated data frame) and `gamma`, which is again a list with elements `condition`, `rt` and `correct`, each a tibble with two columns (see details for more information).

### Note

Different parameters for different conditions are only allowed for drift rate,  $v$ , and process variability,  $s$ . All other parameters are used for all conditions.

### Author(s)

Sebastian Hellmann.

### References

- Hellmann, S., Zehetleitner, M., & Rausch, M. (2023). Simultaneous modeling of choice, confidence and response time in visual perception. *Psychological Review* 2023 Mar 13. doi: 10.1037/rev0000411. Epub ahead of print. PMID: 36913292.
- Kiani, R., Corthell, L., & Shadlen, M.N. (2014) Choice certainty is informed by both evidence and decision time. *Neuron*, 84(6), 1329-1342. doi:10.1016/j.neuron.2014.12.015

## Examples

```
# Examples for "PCRM" model (equivalent applicable for "IRM" model)
# 1. Define some parameter set in a data.frame
paramDf <- data.frame(a=2,b=2, v1=0.5, v2=1, t0=0.1,st0=0,
                      wx=0.6, wint=0.2, wrt=0.2,
                      theta1=4)

# 2. Simulate trials for both stimulus categories and all conditions (2)
simus <- simulateRM(paramDf, n=30,model="PCRM", time_scaled=TRUE)
head(simus)
# equivalent:
simus <- simulateRM(paramDf, model="PCRMt")

library(ggplot2)
simus <- simus[simus$response!=0,]
simus$rating <- factor(simus$rating, labels=c("unsure", "sure"))
ggplot(simus, aes(x=rt, group=interaction(correct, rating),
                 color=as.factor(correct), linetype=rating))+
  geom_density(linewidth=1.2, na.rm=TRUE)+
  facet_grid(rows=vars(condition), labeller = "label_both")

# automatically aggregate simulation distribution
# to get only accuracy x confidence rating distribution for
# all conditions
agg_simus <- simulateRM(paramDf, n = 20, model="PCRMt", agg_simus = TRUE)
head(agg_simus)

agg_simus$rating <- factor(agg_simus$rating, labels=c("unsure", "sure"))
library(ggplot2)
ggplot(agg_simus, aes(x=rating, group=correct, fill=as.factor(correct), y=p))+
  geom_bar(stat="identity", position="dodge")+
  facet_grid(cols=vars(condition), labeller = "label_both")

# Compute Gamma correlation coefficients between
# confidence and other behavioral measures
# output will be a list
simu_list <- simulateRM(paramDf, model="IRMt", gamma=TRUE, n=200)
simu_list
```

## Description

Simulates the decision responses, reaction times and confidence measure together with a discrete confidence judgment for the sequential sampling confidence model specified by the argument `model`, given specific parameter constellations. This function is a wrapper that calls the respective functions for diffusion based models (dynaViTE and 2DSD: [simulateWEV](#)) and race models (IRM, PCRM, IRMt, and PCRMt: [simulateRM](#); and MTLNR: [simulateMTLNR](#)). It also computes the Gamma rank correlation between the confidence ratings and condition (task difficulty), reaction times and accuracy in the simulated output.

## Usage

```
simulateRTConf(paramDf, n = 10000, model = NULL, gamma = FALSE,
  agg_simus = FALSE, simult_conf = FALSE, stimulus = c(1, 2),
  delta = 0.01, maxrt = 15, seed = NULL)
```

## Arguments

<code>paramDf</code>	a list or dataframe with one row with the required parameters.
<code>n</code>	integer. The number of samples (per condition and stimulus direction) generated. Total number of samples is $n \times nConditions \times \text{length}(stimulus)$ .
<code>model</code>	character scalar. One of "dynaViTE", "dynWEV", "2DSD", "2DSDT", "IRM", "PCRM", "IRMt", "PCRMt", or "MTLNR". Could also be passed as a column in the <code>paramDf</code> argument.
<code>gamma</code>	logical. If TRUE, the gamma correlation between confidence ratings, rt and accuracy is computed.
<code>agg_simus</code>	logical. Simulation is done on a trial basis with RTs outcome. If TRUE, the simulations will be aggregated over RTs to return only the distribution of response and confidence ratings. Default: FALSE.
<code>simult_conf</code>	logical. Whether in the experiment confidence was reported simultaneously with the decision. If that is the case decision and confidence judgment are assumed to have happened subsequent before the response. Therefore tau is included in the response time. If the decision was reported before the confidence report, <code>simult_conf</code> should be FALSE.
<code>stimulus</code>	numeric vector. Either 1, 2 or <code>c(1, 2)</code> (default). Together with condition represents the experimental situation. In a 2AFC task the presented stimulus belongs to one of two categories. In the default setting trials with both categories presented are simulated but one can choose to simulate only trials with the stimulus coming from one category.
<code>delta</code>	numerical. Size of steps for the discretized simulation.
<code>maxrt</code>	numerical. Maximum reaction time to be simulated. Default: 15.
<code>seed</code>	numerical. Seeding for non-random data generation. (Also possible outside of the function.)

## Details

The output of the fitting function `fitRTConf` with the respective model fits the argument `paramDf` for simulation. The function calls the respective simulation function for diffusion based models, i.e. `dynaViTE` and `2DSD` (`simulateWEV`) or race models, i.e. `IRM(t)` and `PCRM(t)`, (`simulateRM`). See there for more information.

**Simulation Method:** The simulation is done by simulating normal variables in discretized steps until the processes reach the boundary. If no boundary is met within the maximum time, response is set to 0. The MTLNR model is simulated without discretization, but simply by simulating the necessary log-normally distributed random variables.

**Gamma correlations:** The Gamma coefficients are computed separately for correct/incorrect responses for the correlation of confidence ratings with condition and rt and separately for conditions for the correlation of accuracy and confidence. The resulting data frames in the output thus have two columns. One for the grouping variable and one for the Gamma coefficient.

## Value

Depending on `gamma` and `agg_simus`.

If `gamma` is `FALSE`, returns a `data.frame` with columns: `condition`, `stimulus`, `response`, `correct`, `rt`, `conf` (the continuous confidence measure) and `rating` (the discrete confidence rating) or (if `agg_simus=TRUE`): `condition`, `stimulus`, `response`, `correct`, `rating` and `p` (for the probability of a response and rating, given the condition and stimulus).

If `gamma` is `TRUE`, returns a list with elements: `simus` (the simulated data frame) and `gamma`, which is again a list with elements `condition`, `rt` and `correct`, each a tibble with two columns (see details for more information).

## Author(s)

Sebastian Hellmann.

## Examples

```
# The function is particularly useful, when having a collection
# of parameter sets for different models (e.g. output by fitRTConfModels for
# more than one model).
library(dplyr)
# 1. Generate only one parameter set but for two different models
paramDf1 <- data.frame(model="dynWEV", a=1.5, v1=0.2, v2=1, t0=0.1, z=0.52,
  sz=0.3, sv=0.4, st0=0, tau=3, w=0.5,
  theta1=1, svis=0.5, sigvis=0.8)
paramDf2 <- data.frame(model="PCRMt", a=2, b=2, v1=0.5, v2=1, t0=0.1, st0=0,
  wx=0.6, wint=0.2, wrt=0.2, theta1=4)
paramDf <- full_join(paramDf1, paramDf2)
paramDf # each model parameters sets hat its relevant parameters
# Split paramDf by model (maybe also other columns) and simulate data
simus <- paramDf %>% group_by(model) %>%
  reframe(simulateRTConf(cbind(cur_group(), pick(everything()))), n=200, simult_conf = TRUE))
head(simus)
```

---

simulateWEV	<i>Simulation of confidence ratings and RTs in dynWEV and 2DSD confidence models</i>
-------------	--

---

## Description

Simulates the decision responses and reaction times together with a discrete confidence judgment in the dynaViTE model, the 2DSD model (Pleskac & Busemeyer, 2010) and the dynWEV model (Hellmann et al., 2023), given specific parameter constellations. See [ddynaViTE](#) and [d2DSD](#) for more information about parameters. Also computes the Gamma rank correlation between the confidence ratings and condition (task difficulty), reaction times and accuracy in the simulated output. Basically, this function is a wrapper for [rdynaViTE](#) and [r2DSD](#) for application in confidence experiments with manipulation of specific parameters.

## Usage

```
simulateWEV(paramDf, n = 10000, model = "dynWEV", simult_conf = FALSE,
  gamma = FALSE, agg_simus = FALSE, stimulus = c(-1, 1), delta = 0.01,
  maxrt = 15, seed = NULL, process_results = FALSE)
```

## Arguments

paramDf	a list or dataframe with one row. Column names should match the names of <a href="#">dynaViTE</a> and <a href="#">2DSD</a> model specific parameter names. For different stimulus quality/mean drift rates, names should be v1, v2, v3,... Different sv and/or s parameters are possible with sv1, sv2, sv3... (s1, s2, s3,... respectively) with equally many steps as for drift rates. Additionally, the confidence thresholds should be given by names with thetaUpper1, thetaUpper2,..., thetaLower1,... or, for symmetric thresholds only by theta1, theta2,....
n	integer. The number of samples (per condition and stimulus direction) generated. Total number of samples is $n \times nConditions \times length(stimulus)$ .
model	character scalar. One of "dynaViTE", "dynWEV", or "2DSD".
simult_conf	logical. TRUE, if in the experiment confidence was reported simultaneously with the decision, as then decision and confidence judgment are assumed to have happened subsequent before response and tau is added to the simulated decision time. If FALSE returned response time will only be decision time plus non-judgment time component.
gamma	logical. If TRUE, the gamma correlation between confidence ratings, rt and accuracy is computed.
agg_simus	logical. Simulation is done on a trial basis with RTs outcome. If TRUE, the simulations will be aggregated over RTs to return only the distribution of response and confidence ratings. Default: FALSE.
stimulus	numeric vector. Either 1, -1 or c(-1, 1) (default). Together with condition represents the experimental situation. In a binary decision task the presented stimulus belongs to one of two categories. In the default setting trials with both categories

	presented are simulated but one can choose to simulate only trials with the stimulus coming from one category (1 for the category that is associated with positive drift in the decision process where "upper"/1 responses are considered correct and -1 correspondingly for negative drifts and "lower"/-1 correct decisions).
delta	numeric. Discretization steps for simulations with the stochastic process.
maxrt	numeric. Maximum reaction time returned. If the simulation of the stochastic process exceeds a rt of maxrt, the response will be set to 0 and maxrt will be returned as rt.
seed	numerical. Seeding for non-random data generation.
process_results	logical. Whether the output simulations should contain the final state of the decision (and visibility) process as additional column. Default is FALSE, meaning that no additional columns for the final process states are returned.

### Details

Simulation of response and decision times is done by simulating normal variables in discretized steps until the lower or upper boundary is met (or the maximal rt is reached). Afterwards, a confidence measure is simulated according to the respective model.

The confidence outputs are then binned according to the given thresholds. The output of the fitting function `fitRTConf` with the respective model fits the argument `paramDf` for simulation. The Gamma coefficients are computed separately for correct/incorrect responses for the correlation of confidence ratings with condition and rt and separately for conditions for the correlation of accuracy and confidence. The resulting data frames in the output thus have two columns. One for the grouping variable and one for the Gamma coefficient.

### Value

Depending on `gamma` and `agg_simus`.

If `gamma` is FALSE, returns a `data.frame` with columns: `condition`, `stimulus`, `response`, `correct`, `rt`, `conf` (the continuous confidence measure) and `rating` (the discrete confidence rating), and `dec` and `vis` (only if `process_results=TRUE`) for the final states of accumulators in the simulation or (if `agg_simus=TRUE`): `condition`, `stimulus`, `response`, `correct`, `rating` and `p` (for the probability of a response and rating, given the condition and stimulus).

If `gamma` is TRUE, returns a list with elements: `simus` (the simulated data frame) and `gamma`, which is again a list with elements `condition`, `rt` and `correct`, each a tibble with two columns (see details for more information).

### Note

Different parameters for different conditions are only allowed for drift rate, `v`, drift rate variability, `sv` and diffusion constant `s`. All other parameters are used for all conditions.

### Author(s)

Sebastian Hellmann.

## References

Hellmann, S., Zehetleitner, M., & Rausch, M. (2023). Simultaneous modeling of choice, confidence and response time in visual perception. *Psychological Review* 2023 Mar 13. doi: 10.1037/rev0000411. Epub ahead of print. PMID: 36913292.

## Examples

```
# Examples for "dynWEV" model (equivalent applicable
# for "2DSD" model (with less parameters))
# 1. Define some parameter set in a data.frame
paramDf <- data.frame(a=2.5,v1=0.1, v2=1, t0=0.1,z=0.55,
                      sz=0.3,sv=0.8, st0=0, tau=3, w=0.1,
                      theta1=0.8, svis=0.5, sigvis=0.8)

# 2. Simulate trials for both stimulus categories and all conditions (2)
simus <- simulateWEV(paramDf, model="dynWEV")
head(simus)

library(ggplot2)
simus <- simus[simus$response!=0,]
simus$rating <- factor(simus$rating, labels=c("unsure", "sure"))
ggplot(simus, aes(x=rt, group=interaction(correct, rating),
                 color=as.factor(correct), linetype=rating))+
  geom_density(linewidth=1.2, na.rm=TRUE)+xlim(c(0,5))+
  facet_grid(rows=vars(condition), labeller = "label_both")

# automatically aggregate simulation distribution
# to get only accuracy x confidence rating distribution for
# all conditions
agg_simus <- simulateWEV(paramDf, model="dynWEV", agg_simus = TRUE)
head(agg_simus)

agg_simus$rating <- factor(agg_simus$rating, labels=c("unsure", "sure"))
library(ggplot2)
ggplot(agg_simus, aes(x=rating, group=correct, fill=as.factor(correct), y=p))+
  geom_bar(stat="identity", position="dodge")+
  facet_grid(cols=vars(condition), labeller = "label_both")

# Compute Gamma correlation coefficients between
# confidence and other behavioral measures
# output will be a list
simu_list <- simulateWEV(paramDf,n = 400, model="dynWEV", gamma=TRUE)
simu_list
```

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