

# Package ‘SimDesign’

August 17, 2024

**Title** Structure for Organizing Monte Carlo Simulation Designs

**Version** 2.17.1

**Description** Provides tools to safely and efficiently organize and execute Monte Carlo simulation experiments in R. The package controls the structure and back-end of Monte Carlo simulation experiments by utilizing a generate-analyse-summarise workflow. The workflow safeguards against common simulation coding issues, such as automatically re-simulating non-convergent results, prevents inadvertently overwriting simulation files, catches error and warning messages during execution, implicitly supports parallel processing with high-quality random number generation, and provides tools for managing high-performance computing (HPC) array jobs submitted to schedulers such as SLURM. For a pedagogical introduction to the package see Sigal and Chalmers (2016) <[doi:10.1080/10691898.2016.1246953](https://doi.org/10.1080/10691898.2016.1246953)>. For a more in-depth overview of the package and its design philosophy see Chalmers and Adkins (2020) <[doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)>.

**VignetteBuilder** knitr

**Depends** R (>= 3.5.0)

**Imports** methods, testthat, parallel, parallelly, snow, dplyr, sessioninfo, beepr, pbapply (>= 1.3-0), RPushbullet, future, future.apply, progressr, R.utils, stats

**Suggests** knitr, ggplot2, tidyr, purrr, shiny, copula, extraDistr, renv, cli, job, future.batchtools, FrF2, rmarkdown

**License** GPL (>= 2)

**ByteCompile** yes

**LazyData** true

**URL** <http://philchalmers.github.io/SimDesign/>,  
<https://github.com/philchalmers/SimDesign/wiki>

**RoxygenNote** 7.3.2

**NeedsCompilation** no

**Author** Phil Chalmers [aut, cre] (<<https://orcid.org/0000-0001-5332-2810>>),  
Matthew Sigal [ctb],  
Ogreden Oguzhan [ctb],  
Mikko Ronkko [ctb]

**Maintainer** Phil Chalmers <rphilip.chalmers@gmail.com>

**Repository** CRAN

**Date/Publication** 2024-08-17 05:00:02 UTC

## Contents

|                        |    |
|------------------------|----|
| addMissing             | 3  |
| Analyse                | 5  |
| AnalyseIf              | 7  |
| Attach                 | 9  |
| BF_sim                 | 11 |
| BF_sim_alternative     | 12 |
| bias                   | 13 |
| bootPredict            | 15 |
| Bradley1978            | 18 |
| CC                     | 20 |
| clusterSetRNGSubStream | 21 |
| colVars                | 21 |
| createDesign           | 22 |
| ECR                    | 24 |
| EDR                    | 27 |
| expandDesign           | 28 |
| Generate               | 29 |
| GenerateIf             | 31 |
| genSeeds               | 33 |
| getArrayID             | 35 |
| IRMSE                  | 36 |
| MAE                    | 38 |
| manageMessages         | 39 |
| manageWarnings         | 42 |
| MSRSE                  | 47 |
| nc                     | 48 |
| PBA                    | 50 |
| quiet                  | 53 |
| RAB                    | 54 |
| rbind.SimDesign        | 55 |
| RD                     | 57 |
| RE                     | 58 |
| rejectionSampling      | 60 |
| reSummarise            | 63 |
| rHeadrick              | 66 |
| rint                   | 68 |
| rinvWishart            | 69 |
| rmgh                   | 70 |
| RMSE                   | 71 |
| rmvnorm                | 73 |
| rmvt                   | 74 |

|                              |     |
|------------------------------|-----|
| RobbinsMonro . . . . .       | 76  |
| RSE . . . . .                | 78  |
| rtruncate . . . . .          | 79  |
| runArraySimulation . . . . . | 81  |
| runSimulation . . . . .      | 87  |
| rValeMaurelli . . . . .      | 104 |
| Serlin2000 . . . . .         | 106 |
| SFA . . . . .                | 107 |
| SimAnova . . . . .           | 111 |
| SimCheck . . . . .           | 113 |
| SimClean . . . . .           | 114 |
| SimCollect . . . . .         | 115 |
| SimDesign . . . . .          | 118 |
| SimExtract . . . . .         | 119 |
| SimFunctions . . . . .       | 121 |
| SimResults . . . . .         | 123 |
| SimShiny . . . . .           | 125 |
| SimSolve . . . . .           | 127 |
| Summarise . . . . .          | 136 |
| timeFormater . . . . .       | 137 |

**Index****139**


---

|            |   |
|------------|---|
| addMissing | <i>Add missing values to a vector given a MCAR, MAR, or MNAR scheme</i> |
|------------|---|

---

**Description**

Given an input vector, replace elements of this vector with missing values according to some scheme. Default method replaces input values with a MCAR scheme (where on average 10% of the values will be replaced with NAs). MAR and MNAR are supported by replacing the default FUN argument.

**Usage**

```
addMissing(y, fun = function(y, rate = 0.1, ...) rep(rate, length(y)), ...)
```

**Arguments**

|     |   |
|-----|---|
| y   | an input vector that should contain missing data in the form of NA's  |
| fun | a user defined function indicating the missing data mechanism for each element in y. Function must return a vector of probability values with the length equal to the length of y. Each value in the returned vector indicates the probability that the respective element in y will be replaced with NA. Function must contain the argument y, representing the input vector, however any number of additional arguments can be included |
| ... | additional arguments to be passed to FUN  |

## Details

Given an input vector  $y$ , and other relevant variables inside ( $X$ ) and outside ( $Z$ ) the data-set, the three types of missingness are:

**MCAR** Missing completely at random (MCAR). This is realized by randomly sampling the values of the input vector ( $y$ ) irrespective of the possible values in  $X$  and  $Z$ . Therefore missing values are randomly sampled and do not depend on any data characteristics and are truly random

**MAR** Missing at random (MAR). This is realized when values in the dataset ( $X$ ) predict the missing data mechanism in  $y$ ; conceptually this is equivalent to  $P(y = NA|X)$ . This requires the user to define a custom missing data function

**MNAR** Missing not at random (MNAR). This is similar to MAR except that the missing mechanism comes from the value of  $y$  itself or from variables outside the working dataset; conceptually this is equivalent to  $P(y = NA|X, Z, y)$ . This requires the user to define a custom missing data function

## Value

the input vector  $y$  with the sampled NA values (according to the FUN scheme)

## Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

## References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

## Examples

```
## Not run:

set.seed(1)
y <- rnorm(1000)

## 10% missing rate with default FUN
head(ymiss <- addMissing(y), 10)

## 50% missing with default FUN
head(ymiss <- addMissing(y, rate = .5), 10)

## missing values only when female and low
X <- data.frame(group = sample(c('male', 'female'), 1000, replace=TRUE),
                level = sample(c('high', 'low'), 1000, replace=TRUE))
head(X)
```

```

fun <- function(y, X, ...){
  p <- rep(0, length(y))
  p[X$group == 'female' & X$level == 'low'] <- .2
  p
}

ymiss <- addMissing(y, X, fun=fun)
tail(cbind(ymiss, X), 10)

## missingness as a function of elements in X (i.e., a type of MAR)
fun <- function(y, X){
  # missingness with a logistic regression approach
  df <- data.frame(y, X)
  mm <- model.matrix(y ~ group + level, df)
  cfs <- c(-5, 2, 3) #intercept, group, and level coeffs
  z <- cfs %*% t(mm)
  plogis(z)
}

ymiss <- addMissing(y, X, fun=fun)
tail(cbind(ymiss, X), 10)

## missing values when y elements are large (i.e., a type of MNAR)
fun <- function(y) ifelse(abs(y) > 1, .4, 0)
ymiss <- addMissing(y, fun=fun)
tail(cbind(y, ymiss), 10)

## End(Not run)

```

---

Analyse

---

*Compute estimates and statistics*


---

## Description

Compute all relevant test statistics, parameter estimates, detection rates, and so on. This is the computational heavy lifting portion of the Monte Carlo simulation. Users may define a single Analysis function to perform all the analyses in the same function environment, or may define a list of named functions to [runSimulation](#) to allow for a more modularized approach to performing the analyses in independent blocks (but that share the same generated data). Note that if a suitable [Generate](#) function was not supplied then this function can be used to generate and analyse the Monte Carlo data (though in general this setup is not recommended for larger simulations).

## Usage

```
Analyse(condition, dat, fixed_objects)
```

**Arguments**

|                            |  |
|----------------------------|--|
| <code>condition</code>     | a single row from the design input (as a <code>data.frame</code> ), indicating the simulation conditions   |
| <code>dat</code>           | the <code>dat</code> object returned from the <code>Generate</code> function (usually a <code>data.frame</code> , <code>matrix</code> , <code>vector</code> , or <code>list</code> ) |
| <code>fixed_objects</code> | object passed down from <code>runSimulation</code>   |

**Details**

In some cases, it may be easier to change the output to a named `list` containing different parameter configurations (e.g., when determining RMSE values for a large set of population parameters).

The use of `try` functions is generally not required in this function because `Analyse` is internally wrapped in a `try` call. Therefore, if a function stops early then this will cause the function to halt internally, the message which triggered the `stop` will be recorded, and `Generate` will be called again to obtain a different dataset. That said, it may be useful for users to throw their own `stop` commands if the data should be re-drawn for other reasons (e.g., an estimated model terminated correctly but the maximum number of iterations were reached).

**Value**

returns a named numeric vector or `data.frame` with the values of interest (e.g., p-values, effects sizes, etc), or a `list` containing values of interest (e.g., separate matrix and vector of parameter estimates corresponding to elements in parameters). If a `data.frame` is returned with more than 1 row then these objects will be wrapped into suitable `list` objects

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

**See Also**

`stop`, `AnalyseIf`, `manageWarnings`

**Examples**

```
## Not run:

analyse <- function(condition, dat, fixed_objects) {

  # require packages/define functions if needed, or better yet index with the :: operator
  require(stats)
  mygreatfunction <- function(x) print('Do some stuff')

  #wrap computational statistics in try() statements to control estimation problems
  welch <- t.test(DV ~ group, dat)
```

```

ind <- stats::t.test(DV ~ group, dat, var.equal=TRUE)

# In this function the p values for the t-tests are returned,
# and make sure to name each element, for future reference
ret <- c(welch = welch$p.value,
         independent = ind$p.value)

return(ret)
}

# A more modularized example approach

analysis_welch <- function(condition, dat, fixed_objects) {
  welch <- t.test(DV ~ group, dat)
  ret <- c(p=welch$p.value)
  ret
}

analysis_ind <- function(condition, dat, fixed_objects) {
  ind <- t.test(DV ~ group, dat, var.equal=TRUE)
  ret <- c(p=ind$p.value)
  ret
}

# pass functions as a named list
# runSimulation(..., analyse=list(welch=analyse_welch, independent=analysis_ind))

## End(Not run)

```

---

|           |  |
|-----------|--|
| AnalyseIf | <i>Perform a test that indicates whether a given Analyse() function should be executed</i> |
|-----------|--|

---

## Description

This function is designed to prevent specific analysis function executions when the design conditions are not met. Primarily useful when the analyse argument to [runSimulation](#) was input as a named list object, however some of the analysis functions are not interesting/compatible with the generated data and should therefore be skipped.

## Usage

```
AnalyseIf(x, condition = NULL)
```

## Arguments

|   |  |
|---|--|
| x | logical statement to evaluate. If the statement evaluates to TRUE then the remainder of the defined function will be evaluated |
|---|--|

condition (optional) the current design condition. This does not need to be supplied if the expression in `x` evaluates to valid logical (e.g., use `Attach(condition)` prior to using `AnalyseIf`, or use `with(condition, AnalyseIf(someLogicalTest))`)

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

### See Also

[Analyse](#), [runSimulation](#)

### Examples

```
## Not run:

Design <- createDesign(N=c(10,20,30), var.equal = c(TRUE, FALSE))

Generate <- function(condition, fixed_objects) {
  Attach(condition)
  dat <- data.frame(DV = rnorm(N*2), IV = gl(2, N, labels=c('G1', 'G2')))
  dat
}

# always run this analysis for each row in Design
Analyse1 <- function(condition, dat, fixed_objects) {
  mod <- t.test(DV ~ IV, data=dat)
  mod$p.value
}

# Only perform analysis when variances are equal and N = 20 or 30
Analyse2 <- function(condition, dat, fixed_objects) {
  AnalyseIf(var.equal && N %in% c(20, 30), condition)
  mod <- t.test(DV ~ IV, data=dat, var.equal=TRUE)
  mod$p.value
}

Summarise <- function(condition, results, fixed_objects) {
  ret <- EDR(results, alpha=.05)
  ret
}

#-----
```



```

# append names 'Welch' and 'independent' to associated output
res <- runSimulation(design=Design, replications=100, generate=Generate,
                    analyse=list(Welch=Analyse1, independent=Analyse2),
                    summarise=Summarise)

res

# leave results unnamed
res <- runSimulation(design=Design, replications=100, generate=Generate,
                    analyse=list(Analyse1, Analyse2),
                    summarise=Summarise)

## End(Not run)

```

---

 Attach

*Attach objects for easier reference*


---

## Description

The behaviour of this function is very similar to `attach`, however it is environment specific, and therefore only remains defined in a given function rather than in the Global Environment. Hence, this function is much safer to use than the `attach`, which incidentally should never be used in your code. This is useful primarily as a convenience function when you prefer to call the variable names in condition directly rather than indexing with `condition$sample_size` or `with(condition, sample_size)`, for example.

## Usage

```

Attach(
  ...,
  omit = NULL,
  check = TRUE,
  attach_listone = TRUE,
  RStudio_flags = FALSE
)

```

## Arguments

|                   |   |
|-------------------|---|
| <code>...</code>  | a comma separated list of <code>data.frame</code> , <code>tibble</code> , <code>list</code> , or <code>matrix</code> objects containing (column) elements that should be placed in the current working environment  |
| <code>omit</code> | an optional character vector containing the names of objects that should not be attached to the current environment. For instance, if the objects named 'a' and 'b' should not be attached then use <code>omit = c('a', 'b')</code> . When <code>NULL</code> (default) all objects are attached |

|                |  |
|----------------|--|
| check          | logical; check to see if the function will accidentally replace previously defined variables with the same names as in condition? Default is TRUE, which will avoid this error   |
| attach_listone | logical; if the element to be assign is a list of length one then assign the first element of this list with the associated name. This generally avoids adding an often unnecessary list 1 index, such as <code>name &lt;- list[[1L]]</code>           |
| RStudio_flags  | logical; print R script output comments that disable flagged missing variables in RStudio? Requires the form <code>Attach(Design, RStudio_flags=TRUE)</code> or in an interactive debugging session <code>Attach(condition, RStudio_flags=TRUE)</code> |

### Details

Note that if you are using RStudio with the *"Warn if variable used has no definition in scope"* diagnostic flag then using `Attach()` will raise suspensions. To suppress such issues, you can either disable such flags (the atomic solution) or evaluate the following output in the R console and place the output in your working simulation file.

```
Attach(Design, RStudio_flags = TRUE)
```

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

### See Also

[runSimulation](#), [Generate](#)

### Examples

```
Design <- createDesign(N1=c(10,20),
                     N2=c(10,20),
                     sd=c(1,2))

Design

# does not use Attach()
Generate <- function(condition, fixed_objects ) {
  # condition = single row of Design input (e.g., condition <- Design[1,])
  N1 <- condition$N1
  N2 <- condition$N2
  sd <- condition$sd

  group1 <- rnorm(N1)
```

```

group2 <- rnorm(N2, sd=sd)
dat <- data.frame(group = c(rep('g1', N1), rep('g2', N2)),
                  DV = c(group1, group2))
dat
}

# similar to above, but using the Attach() function instead of indexing
Generate <- function(condition, fixed_objects ) {
  Attach(condition) # N1, N2, and sd are now 'attached' and visible

  group1 <- rnorm(N1)
  group2 <- rnorm(N2, sd=sd)
  dat <- data.frame(group = c(rep('g1', N1), rep('g2', N2)),
                    DV = c(group1, group2))
  dat
}

#####
# NOTE: if you're using RStudio with code diagnostics on then evaluate + add the
# following output to your source file to manually support the flagged variables

Attach(Design, RStudio_flags=TRUE)

# Below is the same example, however with false positive missing variables suppressed
# when # !diagnostics ... is added added to the source file(s)

# !diagnostics suppress=N1,N2,sd
Generate <- function(condition, fixed_objects ) {
  Attach(condition) # N1, N2, and sd are now 'attached' and visible

  group1 <- rnorm(N1)
  group2 <- rnorm(N2, sd=sd)
  dat <- data.frame(group = c(rep('g1', N1), rep('g2', N2)),
                    DV = c(group1, group2))
  dat
}

```

---

BF\_sim

*Example simulation from Brown and Forsythe (1974)*


---

## Description

Example results from the Brown and Forsythe (1974) article on robust estimators for variance ratio tests. Statistical tests are organized by columns and the unique design conditions are organized by rows. See [BF\\_sim\\_alternative](#) for an alternative form of the same simulation. Code for this simulation is available of the wiki (<https://github.com/philchalmers/SimDesign/wiki>).

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Brown, M. B. and Forsythe, A. B. (1974). Robust tests for the equality of variances. *Journal of the American Statistical Association*, 69(346), 364–367.

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

**Examples**

```
## Not run:  
data(BF_sim)  
head(BF_sim)  
  
#Type I errors  
subset(BF_sim, var_ratio == 1)  
  
## End(Not run)
```

---

BF\_sim\_alternative      (Alternative) Example simulation from Brown and Forsythe (1974)

---

**Description**

Example results from the Brown and Forsythe (1974) article on robust estimators for variance ratio tests. Statistical tests and distributions are organized by columns and the unique design conditions are organized by rows. See [BF\\_sim](#) for an alternative form of the same simulation where distributions are also included in the rows. Code for this simulation is available on the wiki (<https://github.com/philchalmers/SimDesign/wiki>).

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Brown, M. B. and Forsythe, A. B. (1974). Robust tests for the equality of variances. *Journal of the American Statistical Association*, 69(346), 364–367.

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

**Examples**

```
## Not run:
data(BF_sim_alternative)
head(BF_sim_alternative)

#' #Type I errors
subset(BF_sim_alternative, var_ratio == 1)

## End(Not run)
```

---

bias

---

*Compute (relative/standardized) bias summary statistic*


---

**Description**

Computes the (relative) bias of a sample estimate from the parameter value. Accepts estimate and parameter values, as well as estimate values which are in deviation form. If relative bias is requested the estimate and parameter inputs are both required.

**Usage**

```
bias(
  estimate,
  parameter = NULL,
  type = "bias",
  abs = FALSE,
  percent = FALSE,
  unname = FALSE
)
```

**Arguments**

|           |   |
|-----------|---|
| estimate  | a numeric vector, matrix/data.frame, or list of parameter estimates. If a vector, the length is equal to the number of replications. If a matrix/data.frame, the number of rows must equal the number of replications. list objects will be looped over using the same rules after above after first translating the information into one-dimensional vectors and re-creating the structure upon return |
| parameter | a numeric scalar/vector indicating the fixed parameters. If a single value is supplied and estimate is a matrix/data.frame then the value will be recycled for each column; otherwise, each element will be associated with each respective column in the estimate input. If NULL then it will be assumed that the estimate input is in a deviation form (therefore mean(estimate)) will be returned)   |
| type      | type of bias statistic to return. Default ('bias') computes the standard bias (average difference between sample and population), 'relative' computes the relative bias statistic (i.e., divide the bias by the value in parameter; note that multiplying this by 100 gives the "percent bias" measure, or if Type I error rates  |

( $\alpha$ ) are supplied will result in the "percentage error"), 'abs\_relative' computes the relative bias but the absolute values of the parameters are used in the denominator rather than the (potentially) signed input values, and 'standardized' computes the standardized bias estimate (standard bias divided by the standard deviation of the sample estimates)

|         |   |
|---------|---|
| abs     | logical; find the absolute bias between the parameters and estimates? This effectively just applies the <a href="#">abs</a> transformation to the returned result. Default is FALSE |
| percent | logical; change returned result to percentage by multiplying by 100? Default is FALSE   |
| unnname | logical; apply <a href="#">unnname</a> to the results to remove any variable names?   |

### Value

returns a numeric vector indicating the overall (relative/standardized) bias in the estimates

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. [doi:10.1080/10691898.2016.1246953](https://doi.org/10.1080/10691898.2016.1246953)

### See Also

[RMSE](#)

### Examples

```
pop <- 2
samp <- rnorm(100, 2, sd = 0.5)
bias(samp, pop)
bias(samp, pop, type = 'relative')
bias(samp, pop, type = 'standardized')

dev <- samp - pop
bias(dev)

# equivalent here
bias(mean(samp), pop)

# matrix input
mat <- cbind(M1=rnorm(100, 2, sd = 0.5), M2 = rnorm(100, 2, sd = 1))
bias(mat, parameter = 2)
bias(mat, parameter = 2, type = 'relative')
```

```

bias(mat, parameter = 2, type = 'standardized')

# different parameter associated with each column
mat <- cbind(M1=rnorm(1000, 2, sd = 0.25), M2 = rnorm(1000, 3, sd = .25))
bias(mat, parameter = c(2,3))

# same, but with data.frame
df <- data.frame(M1=rnorm(100, 2, sd = 0.5), M2 = rnorm(100, 2, sd = 1))
bias(df, parameter = c(2,2))

# parameters of the same size
parameters <- 1:10
estimates <- parameters + rnorm(10)
bias(estimates, parameters)

# relative difference dividing by the magnitude of parameters
bias(estimates, parameters, type = 'abs_relative')

# relative bias as a percentage
bias(estimates, parameters, type = 'abs_relative', percent = TRUE)

# percentage error (PE) statistic given alpha (Type I error) and EDR() result
# edr <- EDR(results, alpha = .05)
edr <- c(.04, .05, .06, .08)
bias(matrix(edr, 1L), .05, type = 'relative', percent = TRUE)

```

---

bootPredict

---

*Compute prediction estimates for the replication size using bootstrap  
MSE estimates*


---

## Description

This function computes bootstrap mean-square error estimates to approximate the sampling behavior of the meta-statistics in SimDesign's summarise functions. A single design condition is supplied, and a simulation with  $\max(Rstar)$  replications is performed whereby the generate-analyse results are collected. After obtaining these replication values, the replications are further drawn from (with replacement) using the differing sizes in Rstar to approximate the bootstrap MSE behavior given different replication sizes. Finally, given these bootstrap estimates linear regression models are fitted using the predictor term  $one\_sqrtR = 1 / \sqrt{Rstar}$  to allow extrapolation to replication sizes not observed in Rstar. For more information about the method and subsequent bootstrap MSE plots, refer to Koehler, Brown, and Haneuse (2009).

## Usage

```

bootPredict(
  condition,
  generate,

```

```

  analyse,
  summarise,
  fixed_objects = NULL,
  ...,
  Rstar = seq(100, 500, by = 100),
  boot_draws = 1000
)

boot_predict(...)

```

### Arguments

|               |  |
|---------------|--|
| condition     | a data.frame consisting of one row from the original design input object used within <code>runSimulation</code>  |
| generate      | see <code>runSimulation</code>   |
| analyse       | see <code>runSimulation</code>   |
| summarise     | see <code>runSimulation</code>   |
| fixed_objects | see <code>runSimulation</code>   |
| ...           | additional arguments to be passed to <code>runSimulation</code>  |
| Rstar         | a vector containing the size of the bootstrap subsets to obtain. Default investigates the vector [100, 200, 300, 400, 500] to compute the respective MSE terms |
| boot_draws    | number of bootstrap replications to draw. Default is 1000  |

### Value

returns a list of linear model objects (via `lm`) for each meta-statistics returned by the `summarise()` function

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

- Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248
- Koehler, E., Brown, E., & Haneuse, S. J.-P. A. (2009). On the Assessment of Monte Carlo Error in Simulation-Based Statistical Analyses. *The American Statistician*, 63, 155-162.
- Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953



**Examples**

```

set.seed(4321)
Design <- createDesign(sigma = c(1, 2))

#-----

Generate <- function(condition, fixed_objects) {
  dat <- rnorm(100, 0, condition$sigma)
  dat
}

Analyse <- function(condition, dat, fixed_objects) {
  CIs <- t.test(dat)$conf.int
  names(CIs) <- c('lower', 'upper')
  ret <- c(mean = mean(dat), CIs)
  ret
}

Summarise <- function(condition, results, fixed_objects) {
  ret <- c(mu_bias = bias(results[, "mean"], 0),
          mu_coverage = ECR(results[, c("lower", "upper")], parameter = 0))
  ret
}

## Not run:
# boot_predict supports only one condition at a time
out <- bootPredict(condition=Design[1L, , drop=FALSE],
  generate=Generate, analyse=Analyse, summarise=Summarise)
out # list of fitted linear model(s)

# extract first meta-statistic
mu_bias <- out$mu_bias

dat <- model.frame(mu_bias)
print(dat)

# original R metric plot
R <- 1 / dat$one_sqrtR^2
plot(R, dat$MSE, type = 'b', ylab = 'MSE', main = "Replications by MSE")

plot(MSE ~ one_sqrtR, dat, main = "Bootstrap prediction plot", xlim = c(0, max(one_sqrtR)),
  ylim = c(0, max(MSE)), ylab = 'MSE', xlab = expression(1/sqrt(R)))
beta <- coef(mu_bias)
abline(a = 0, b = beta, lty = 2, col='red')

# what is the replication value when x-axis = .02? What's its associated expected MSE?
1 / .02^2 # number of replications
predict(mu_bias, data.frame(one_sqrtR = .02)) # y-axis value

# approximately how many replications to obtain MSE = .001?
(beta / .001)^2

```

```
## End(Not run)
```

---

Bradley1978

*Bradley's (1978) empirical robustness interval*


---

## Description

Robustness interval criteria for empirical detection rate estimates and empirical coverage estimates defined by Bradley (1978). See [EDR](#) and [ECR](#) to obtain such estimates.

## Usage

```
Bradley1978(
  rate,
  alpha = 0.05,
  type = "liberal",
  CI = FALSE,
  out.logical = FALSE,
  out.labels = c("conservative", "robust", "liberal"),
  unname = FALSE
)
```

## Arguments

|             |  |
|-------------|--|
| rate        | (optional) numeric vector containing the empirical detection rate(s) or empirical confidence interval estimates. If supplied a character vector with elements defined in <code>out.labels</code> or a logical vector will be returned indicating whether the detection rate estimate is considered 'robust'.<br>When the input is an empirical coverage rate the argument <code>CI</code> must be set to <code>TRUE</code> .<br>If this input is missing, the interval criteria will be printed to the console |
| alpha       | Type I error rate to evaluated (default is .05)  |
| type        | character vector indicating the type of interval classification to use. Default is 'liberal', however can be 'stringent' to use Bradley's more stringent robustness criteria   |
| CI          | logical; should this robust interval be constructed on empirical detection rates ( <code>FALSE</code> ) or empirical coverage rates ( <code>TRUE</code> )?   |
| out.logical | logical; should the output vector be <code>TRUE/FALSE</code> indicating whether the supplied empirical detection rate/CI should be considered "robust"? Default is <code>FALSE</code> , in which case the <code>out.labels</code> elements are used instead  |
| out.labels  | character vector of length three indicating the classification labels according to the desired robustness interval   |
| unname      | logical; apply <code>unname</code> to the results to remove any variable names?  |

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Bradley, J. V. (1978). Robustness? *British Journal of Mathematical and Statistical Psychology*, 31, 144-152.

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

**See Also**

[EDR](#), [ECR](#), [Serlin2000](#)

**Examples**

```
# interval criteria used for empirical detection rates
Bradley1978()
Bradley1978(type = 'stringent')
Bradley1978(alpha = .01, type = 'stringent')

# intervals applied to empirical detection rate estimates
edr <- c(test1 = .05, test2 = .027, test3 = .051, test4 = .076, test5 = .024)

Bradley1978(edr)
Bradley1978(edr, out.logical=TRUE) # is robust?

#####
# interval criteria used for coverage estimates

Bradley1978(CI = TRUE)
Bradley1978(CI = TRUE, type = 'stringent')
Bradley1978(CI = TRUE, alpha = .01, type = 'stringent')

# intervals applied to empirical coverage rate estimates
ecr <- c(test1 = .950, test2 = .973, test3 = .949, test4 = .924, test5 = .976)

Bradley1978(ecr, CI=TRUE)
Bradley1978(ecr, CI=TRUE, out.logical=TRUE) # is robust?
```

---

**CC** *Compute congruence coefficient*

---

**Description**

Computes the congruence coefficient, also known as an "unadjusted" correlation or Tucker's congruence coefficient.

**Usage**

```
CC(x, y = NULL, unname = FALSE)
```

**Arguments**

**x** a vector or data.frame/matrix containing the variables to use. If a vector then the input y is required, otherwise the congruence coefficient is computed for all bivariate combinations

**y** (optional) the second vector input to use if x is a vector

**unname** logical; apply [unname](#) to the results to remove any variable names?

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. [doi:10.1080/10691898.2016.1246953](https://doi.org/10.1080/10691898.2016.1246953)

**See Also**

[cor](#)

**Examples**

```
vec1 <- runif(1000)
vec2 <- runif(1000)

CC(vec1, vec2)
# compare to cor()
cor(vec1, vec2)

# column input
df <- data.frame(vec1, vec2, vec3 = runif(1000))
CC(df)
```

```
cor(df)
```

---

```
clusterSetRNGSubStream
```

*Set RNG sub-stream for Pierre L'Ecuyer's RngStreams*

---

### Description

Sets the sub-stream RNG state within for Pierre L'Ecuyer's (1999) algorithm. Should be used within distributed array jobs after suitable L'Ecuyer's (1999) have been distributed to each array, and each array is further defined to use multi-core processing. See [clusterSetRNGStream](#) for further information.

### Usage

```
clusterSetRNGSubStream(cl, seed)
```

### Arguments

|      |   |
|------|---|
| cl   | A cluster from the <code>parallel</code> or <code>snow</code> package, or (if NULL) the registered cluster  |
| seed | An integer vector of length 7 as given by <code>.Random.seed</code> when the L'Ecuyer-CMR RNG is in use. See <a href="#">RNG</a> for the valid values |

### Value

invisible NULL

---

```
colVars
```

*Form Column Standard Deviation and Variances*

---

### Description

Form column standard deviation and variances for numeric arrays (or data frames).

### Usage

```
colVars(x, na.rm = FALSE, unname = FALSE)
```

```
colSDs(x, na.rm = FALSE, unname = FALSE)
```

### Arguments

|        |  |
|--------|--|
| x      | an array of two dimensions containing numeric, complex, integer or logical values, or a numeric data frame |
| na.rm  | logical; remove missing values in each respective column?  |
| unname | logical; apply <a href="#">unname</a> to the results to remove any variable names?                         |

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**See Also**

[colMeans](#)

**Examples**

```
results <- matrix(rnorm(100), ncol=4)
colnames(results) <- paste0('stat', 1:4)

colVars(results)
colSDs(results)

results[1,1] <- NA
colSDs(results)
colSDs(results, na.rm=TRUE)
colSDs(results, na.rm=TRUE, unname=TRUE)
```

---

createDesign

*Create the simulation design object*

---

**Description**

Create a partially or fully-crossed data object reflecting the unique simulation design conditions. Each row of the returned object represents a unique simulation condition, and each column represents the named factor variables under study.

**Usage**

```
createDesign(
  ...,
  subset,
  fractional = NULL,
  tibble = TRUE,
  stringsAsFactors = FALSE
)

## S3 method for class 'Design'
print(x, list2char = TRUE, pillar.sigfig = 5, ...)
```

**Arguments**

... comma separated list of named input objects representing the simulation factors to completely cross. Note that these arguments are passed to [expand.grid](#) to perform the complete crossings

|                  |  |
|------------------|--|
| subset           | (optional) a logical vector indicating elements or rows to keep to create a partially crossed simulation design  |
| fractional       | a fractional design matrix returned from the FrF2 package. Note that the order of the factor names/labels are associated with the respective ... inputs  |
| tibble           | logical; return a tibble object instead of a data.frame? Default is TRUE   |
| stringsAsFactors | logical; should character variable inputs be coerced to factors when building a data.frame? Default is FALSE   |
| x                | object returned by <code>createDesign</code>   |
| list2char        | logical; for tibble object re-evaluate list elements as character vectors for better printing of the levels? Note that this does not change the original classes of the object, just how they are printed. Default is TRUE |
| pillar.sigfig    | number of significant digits to print. Default is 5  |

**Value**

a tibble or data.frame containing the simulation experiment conditions to be evaluated in `runSimulation`

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

**See Also**

`expandDesign`

**Examples**

```
## Not run:

# modified example from runSimulation()

Design <- createDesign(N = c(10, 20),
                      SD = c(1, 2))

Design

# remove N=10, SD=2 row from initial definition
Design <- createDesign(N = c(10, 20),
                      SD = c(1, 2),
                      subset = !(N == 10 & SD == 2))

Design
```

```

# example with list inputs
Design <- createDesign(N = c(10, 20),
                      SD = c(1, 2),
                      combo = list(c(0,0), c(0,0,1)))
Design # notice levels printed (not typical for tibble)
print(Design, list2char = FALSE) # standard tibble output

Design <- createDesign(N = c(10, 20),
                      SD = c(1, 2),
                      combo = list(c(0,0), c(0,0,1)),
                      combo2 = list(c(5,10,5), c(6,7)))
Design
print(Design, list2char = FALSE) # standard tibble output

#####

## fractional factorial example

library(FrF2)
# help(FrF2)

# 7 factors in 32 runs
fr <- FrF2(32,7)
dim(fr)
fr[1:6,]

# Create working simulation design given -1/1 combinations
fDesign <- createDesign(sample_size=c(100,200),
                       mean_diff=c(.25, 1, 2),
                       variance_ratio=c(1,4, 8),
                       equal_size=c(TRUE, FALSE),
                       dists=c('norm', 'skew'),
                       same_dists=c(TRUE, FALSE),
                       symmetric=c(TRUE, FALSE),
                       # remove same-normal combo
                       subset = !(symmetric & dists == 'norm'),
                       fractional=fr)

fDesign

## End(Not run)

```

### Description

Computes the detection rate for determining empirical coverage rates given a set of estimated confidence intervals. Note that using `1 - ECR(CIs, parameter)` will provide the empirical detection



rate. Also supports computing the average width of the CIs, which may be useful when comparing the efficiency of CI estimators.

### Usage

```
ECR(
  CIs,
  parameter,
  tails = FALSE,
  CI_width = FALSE,
  complement = FALSE,
  names = NULL,
  unname = FALSE
)
```

### Arguments

|            |  |
|------------|--|
| CIs        | a numeric vector or matrix of confidence interval values for a given parameter value, where the first element/column indicates the lower confidence interval and the second element/column the upper confidence interval. If a vector of length 2 is passed instead then the returned value will be either a 1 or 0 to indicate whether the parameter value was or was not within the interval, respectively. Otherwise, the input must be a matrix with an even number of columns |
| parameter  | a numeric scalar indicating the fixed parameter value. Alternative, a numeric vector object with length equal to the number of rows as CIs (use to compare sets of parameters at once)   |
| tails      | logical; when TRUE returns a vector of length 2 to indicate the proportion of times the parameter was lower or higher than the supplied interval, respectively. This is mainly only useful when the coverage region is not expected to be symmetric, and therefore is generally not required. Note that $1 - \text{sum}(\text{ECR}(\text{CIs}, \text{parameter}, \text{tails}=\text{TRUE})) == \text{ECR}(\text{CIs}, \text{parameter})$   |
| CI_width   | logical; rather than returning the overall coverage rate, return the average width of the CIs instead? Useful when comparing the efficiency of different CI estimators   |
| complement | logical; rather than computing the proportion of population parameters within the CI, return the proportion outside the advertised CI ( $1 - \text{ECR} = \alpha$ ). In the case where only one value is provided, which normally would return a 0 if outside the CI or 1 if inside, the values will be switched (useful when using, for example, CI tests of for the significance of parameters)  |
| names      | an optional character vector used to name the returned object. Generally useful when more than one CI estimate is investigated at once   |
| unname     | logical; apply <code>unname</code> to the results to remove any variable names?  |

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

## References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

## See Also

[EDR](#)

## Examples

```

CIs <- matrix(NA, 100, 2)
for(i in 1:100){
  dat <- rnorm(100)
  CIs[i,] <- t.test(dat)$conf.int
}

ECR(CIs, 0)
ECR(CIs, 0, tails = TRUE)
ECR(CIs, 0, complement = TRUE) # proportion outside interval

# single vector input
CI <- c(-1, 1)
ECR(CI, 0)
ECR(CI, 0, complement = TRUE)
ECR(CI, 2)
ECR(CI, 2, complement = TRUE)
ECR(CI, 2, tails = TRUE)

# parameters of the same size as CI
parameters <- 1:10
CIs <- cbind(parameters - runif(10), parameters + runif(10))
parameters <- parameters + rnorm(10)
ECR(CIs, parameters)

# average width of CIs
ECR(CIs, parameters, CI_width=TRUE)

# ECR() for multiple CI estimates in the same object
parameter <- 10
CIs <- data.frame(lowerCI_1=parameter - runif(10),
                  upperCI_1=parameter + runif(10),
                  lowerCI_2=parameter - 2*runif(10),
                  upperCI_2=parameter + 2*runif(10))

head(CIs)
ECR(CIs, parameter)
ECR(CIs, parameter, tails=TRUE)
ECR(CIs, parameter, CI_width=TRUE)

```

```
# often a good idea to provide names for the output
ECR(CIs, parameter, names = c('this', 'that'))
ECR(CIs, parameter, CI_width=TRUE, names = c('this', 'that'))
ECR(CIs, parameter, tails=TRUE, names = c('this', 'that'))
```

---

EDR *Compute the empirical detection/rejection rate for Type I errors and Power*

---

## Description

Computes the detection/rejection rate for determining empirical Type I error and power rates using information from p-values.

## Usage

```
EDR(p, alpha = 0.05, unname = FALSE)
```

## Arguments

|        |  |
|--------|--|
| p      | a numeric vector or matrix/data.frame of p-values from the desired statistical estimator. If a matrix, each statistic must be organized by column, where the number of rows is equal to the number of replications |
| alpha  | the detection threshold (typical values are .10, .05, and .01). Default is .05   |
| unname | logical; apply <a href="#">unname</a> to the results to remove any variable names?   |

## Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

## References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. [doi:10.1080/10691898.2016.1246953](https://doi.org/10.1080/10691898.2016.1246953)

## See Also

[ECR](#), [Bradley1978](#)

**Examples**

```

rates <- numeric(100)
for(i in 1:100){
  dat <- rnorm(100)
  rates[i] <- t.test(dat)$p.value
}

EDR(rates)
EDR(rates, alpha = .01)

# multiple rates at once
rates <- cbind(runif(1000), runif(1000))
EDR(rates)

```

---

expandDesign

*Create the simulation design object*


---

**Description**

Repeat each design row the specified number of times. This is primarily used for cluster computing where jobs are distributed with batches of replications and later aggregated into a complete simulation object (see [runArraySimulation](#) and [SimCollect](#)).

**Usage**

```
expandDesign(Design, repeat_conditions)
```

**Arguments**

|                   |   |
|-------------------|---|
| Design            | object created by <a href="#">createDesign</a> which should have its rows repeated for optimal HPC schedulers   |
| repeat_conditions | integer vector used to repeat each design row the specified number of times. Can either be a single integer, which repeats each row this many times, or an integer vector equal to the number of total rows in the created object.<br><br>This argument is useful when distributing independent row conditions to cluster computing environments, particularly with different replication information. For example, if 1000 replications in total are the target but the condition is repeated over 4 rows then only 250 replications per row would be required across the repeated conditions. See <a href="#">SimCollect</a> for combining the simulation objects once complete |

**Value**

a tibble or data.frame containing the simulation experiment conditions to be evaluated in [runSimulation](#)

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

**See Also**

[createDesign](#), [SimCollect](#), [runArraySimulation](#)

**Examples**

```
## Not run:

# repeat each row 4 times (for cluster computing)
Design <- createDesign(N = c(10, 20),
                      SD.equal = c(TRUE, FALSE))
Design4 <- expandDesign(Design, 4)
Design4

# repeat first two rows 2x and the rest 4 times (for cluster computing
# where first two conditions are faster to execute)
Design <- createDesign(SD.equal = c(TRUE, FALSE),
                      N = c(10, 100, 1000))
Design24 <- expandDesign(Design, c(2,2,rep(4, 4)))
Design24

## End(Not run)
```

---

Generate

*Generate data*

---

**Description**

Generate data from a single row in the design input (see [runSimulation](#)). R contains numerous approaches to generate data, some of which are contained in the base package, as well as in SimDesign (e.g., [rmgh](#), [rValeMaurelli](#), [rHeadrick](#)). However the majority can be found in external packages. See CRAN's list of possible distributions here: <https://CRAN.R-project.org/view=Distributions>. Note that this function technically can be omitted if the data generation is provided in the [Analyse](#) step, though in general this is not recommended.

**Usage**

```
Generate(condition, fixed_objects)
```

**Arguments**

`condition` a single row from the design input (as a `data.frame`), indicating the simulation conditions

`fixed_objects` object passed down from `runSimulation`

**Details**

The use of `try` functions is generally not required in this function because `Generate` is internally wrapped in a `try` call. Therefore, if a function stops early then this will cause the function to halt internally, the message which triggered the `stop` will be recorded, and `Generate` will be called again to obtain a different dataset. That said, it may be useful for users to throw their own `stop` commands if the data should be re-drawn for other reasons (e.g., an estimated model terminated correctly but the maximum number of iterations were reached).

**Value**

returns a single object containing the data to be analyzed (usually a vector, `matrix`, or `data.frame`), or `list`

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

**See Also**

`addMissing`, `Attach`, `rmgh`, `rValeMaurelli`, `rHeadrick`

**Examples**

```
## Not run:

generate <- function(condition, fixed_objects) {
  N1 <- condition$sample_sizes_group1
  N2 <- condition$sample_sizes_group2
  sd <- condition$standard_deviations

  group1 <- rnorm(N1)
  group2 <- rnorm(N2, sd=sd)
  dat <- data.frame(group = c(rep('g1', N1), rep('g2', N2)),
                    DV = c(group1, group2))
  # just a silly example of a simulated parameter
  pars <- list(random_number = rnorm(1))
}
```

```

    list(dat=dat, parameters=pars)
  }

# similar to above, but using the Attach() function instead of indexing
generate <- function(condition, fixed_objects) {
  Attach(condition)
  N1 <- sample_sizes_group1
  N2 <- sample_sizes_group2
  sd <- standard_deviations

  group1 <- rnorm(N1)
  group2 <- rnorm(N2, sd=sd)
  dat <- data.frame(group = c(rep('g1', N1), rep('g2', N2)),
                    DV = c(group1, group2))
  dat
}

generate2 <- function(condition, fixed_objects) {
  mu <- sample(c(-1,0,1), 1)
  dat <- rnorm(100, mu)
  dat      #return simple vector (discard mu information)
}

generate3 <- function(condition, fixed_objects) {
  mu <- sample(c(-1,0,1), 1)
  dat <- data.frame(DV = rnorm(100, mu))
  dat
}

## End(Not run)

```

---

|            |   |
|------------|---|
| GenerateIf | <i>Perform a test that indicates whether a given Generate() function should be executed</i> |
|------------|---|

---

### Description

This function is designed to prevent specific generate function executions when the design conditions are not met. Primarily useful when the generate argument to `runSimulation` was input as a named list object, however should only be applied for some specific design condition (otherwise, the data generation moves to the next function in the list).

### Usage

```
GenerateIf(x, condition = NULL)
```

**Arguments**

|           |  |
|-----------|--|
| x         | logical statement to evaluate. If the statement evaluates to TRUE then the remainder of the defined function will be evaluated   |
| condition | (optional) the current design condition. This does not need to be supplied if the expression in x evaluates to valid logical (e.g., use <code>Attach(condition)</code> prior to using <code>AnalyseIf</code> , or use <code>with(condition, AnalyseIf(someLogicalTest))</code> ) |

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

**See Also**

[Analyse](#), [runSimulation](#)

**Examples**

```
## Not run:

# SimFunctions(nGenerate = 2)

Design <- createDesign(N=c(10,20,30), var.equal = c(TRUE, FALSE))

Generate.G1 <- function(condition, fixed_objects) {
  GenerateIf(condition$var.equal == FALSE) # only run when unequal vars
  Attach(condition)
  dat <- data.frame(DV = c(rnorm(N), rnorm(N, sd=2)),
                   IV = gl(2, N, labels=c('G1', 'G2')))
  dat
}

Generate.G2 <- function(condition, fixed_objects) {
  Attach(condition)
  dat <- data.frame(DV = rnorm(N*2), IV = gl(2, N, labels=c('G1', 'G2')))
  dat
}

# always run this analysis for each row in Design
Analyse <- function(condition, dat, fixed_objects) {
  mod <- t.test(DV ~ IV, data=dat)
  mod$p.value
}
```



```

Summarise <- function(condition, results, fixed_objects) {
  ret <- EDR(results, alpha=.05)
  ret
}

#-----

# append names 'Welch' and 'independent' to associated output
res <- runSimulation(design=Design, replications=1000,
                    generate=list(G1=Generate.G1, G2=Generate.G2),
                    analyse=Analyse,
                    summarise=Summarise)

res

## End(Not run)

```

---

genSeeds

*Generate random seeds*


---

### Description

Generate seeds to be passed to `runSimulation`'s seed input. Values are sampled from 1 to 2147483647, or are generated using L'Ecuyer-CMRG's (2002) method (returning either a list if `arrayID` is omitted, or the specific row value from this list if `arrayID` is included).

### Usage

```
genSeeds(design = 1L, iseed = NULL, arrayID = NULL, old.seeds = NULL)
```

```
gen_seeds(...)
```

### Arguments

|                      |  |
|----------------------|--|
| <code>design</code>  | design matrix that requires a unique seed per condition, or a number indicating the number of seeds to generate. Default generates one number  |
| <code>iseed</code>   | the initial set.seed number used to generate a sequence of independent seeds according to the L'Ecuyer-CMRG (2002) method. This is recommended whenever quality random number generation is required across similar (if not identical) simulation jobs (e.g., see <a href="#">runArraySimulation</a> ). If <code>arrayID</code> is not specified then this will return a list of the associated seed for the full design |
| <code>arrayID</code> | (optional) single integer input corresponding to the specific row in the design object when using the <code>iseed</code> input. This is used in functions such as <a href="#">runArraySimulation</a> to pull out the specific seed rather than manage a complete list, and is therefore more memory efficient  |

old.seeds        (optional) vector or matrix of last seeds used in previous simulations to avoid repeating the same seed on a subsequent run. Note that this approach should be used sparingly as seeds set more frequently are more likely to correlate, and therefore provide less optimal random number behaviour (e.g., if performing a simulation on two runs to achieve  $5000 * 2 = 10,000$  replications this is likely reasonable, but for simulations with  $100 * 2 = 200$  replications this is more likely to be sub-optimal). Length must be equal to the number of rows in design

...                does nothing

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### Examples

```
# generate 1 seed (default)
genSeeds()

# generate 5 unique seeds
genSeeds(5)

# generate from nrow(design)
design <- createDesign(factorA=c(1,2,3),
                      factorB=letters[1:3])
seeds <- genSeeds(design)
seeds

# construct new seeds that are independent from original (use this sparingly)
newseeds <- genSeeds(design, old.seeds=seeds)
newseeds

# can be done in batches too
newseeds2 <- genSeeds(design, old.seeds=cbind(seeds, newseeds))
cbind(seeds, newseeds, newseeds2) # all unique

#####
# generate seeds for runArraySimulation()
(iseed <- genSeeds()) # initial seed
seed_list <- genSeeds(design, iseed=iseed)
seed_list

# expand number of unique seeds given iseed (e.g., in case more replications
# are required at a later date)
seed_list_tmp <- genSeeds(nrow(design)*2, iseed=iseed)
str(seed_list_tmp) # first 9 seeds identical to seed_list

# more usefully for HPC, extract only the seed associated with an arrayID
arraySeed.15 <- genSeeds(nrow(design)*2, iseed=iseed, arrayID=15)
arraySeed.15
```

---

|            |   |
|------------|---|
| getArrayID | <i>Get job array ID (e.g., from SLURM or other HPC array distributions)</i> |
|------------|---|

---

### Description

Get the array ID from an HPC array distribution job (e.g., from SLURM or from optional command line arguments). The array ID is used to index the rows in the design object in [runArraySimulation](#). For instance, a SLURM array with 10 independent jobs might have the following shell instructions.

### Usage

```
getArrayID(type = "slurm", trailingOnly = TRUE, ID.shift = 0L)
```

### Arguments

|              |  |
|--------------|--|
| type         | an integer indicating the element from the result of <a href="#">commandArgs</a> to extract, or a character specifying the the type of. Default is 'slurm'   |
| trailingOnly | logical value passed to <a href="#">commandArgs</a> . Only used when type is an integer  |
| ID.shift     | single integer value used to shift the array ID by a constant. Useful when there are array range limitation that must be specified in the shell files (e.g., array can only be 10000 but there are more rows in the design object). For example, if the array ID should be 10000 through 12000, but the cluster computer environment does not allow these indices, then including the arrange range as 1-2000 in the shell file with <code>shift=9999</code> would add this constant to the detected arrayID, thereby indexing the remaining row elements in the design object |

### Details

```
#!/bin/bash -l
```

```
#SBATCH -time=00:01:00
```

```
#SBATCH -array=1-10
```

which names the associated jobs with the numbers 1 through 10. `getArrayID()` then extracts this information per array, which is used as the `runArraySimulation(design, ..., arrayID=getArrayID())` to pass specific rows for the design object.

### See Also

[runArraySimulation](#)

### Examples

```
## Not run:
```

```
# get slurm array ID  
arrayID <- getArrayID()
```

```

# get ID based on first optional argument in shell specification
arrayID <- getArrayID(type = 1)

# pass to
# runArraySimulation(design, . . . . , arrayID = arrayID)

# increase detected arrayID by constant 9999 (for array
  specification limitations)
arrayID <- getArrayID(ID.shift=9999)

## End(Not run)

```

---

 IRMSE

---

*Compute the integrated root mean-square error*


---

### Description

Computes the average/cumulative deviation given two continuous functions and an optional function representing the probability density function. Only one-dimensional integration is supported.

### Usage

```

IRMSE(
  estimate,
  parameter,
  fn,
  density = function(theta, ...) 1,
  lower = -Inf,
  upper = Inf,
  ...
)

```

### Arguments

|           |  |
|-----------|--|
| estimate  | a vector of parameter estimates  |
| parameter | a vector of population parameters  |
| fn        | a continuous function where the first argument is to be integrated and the second argument is a vector of parameters or parameter estimates. This function represents a implied continuous function which uses the sample estimates or population parameters   |
| density   | (optional) a density function used to marginalize (i.e., average), where the first argument is to be integrated, and must be of the form <code>density(theta, ...)</code> or <code>density(theta, param1, param2)</code> , where <code>param1</code> is a placeholder name for the hyper-parameters associated with the probability density function. If omitted then the cumulative different between the respective functions will be computed instead |

lower            lower bound to begin numerical integration from  
 upper            upper bound to finish numerical integration to  
 ...                additional parameters to pass to `fnest`, `fnparam`, `density`, and `integrate`,

### Details

The integrated root mean-square error (IRMSE) is of the form

$$IRMSE(\theta) = \sqrt{\int [f(\theta, \hat{\psi}) - f(\theta, \psi)]^2 g(\theta, \dots)}$$

where  $g(\theta, \dots)$  is the density function used to marginalize the continuous sample ( $f(\theta, \hat{\psi})$ ) and population ( $f(\theta, \psi)$ ) functions.

### Value

returns a single numeric term indicating the average/cumulative deviation given the supplied continuous functions

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

### See Also

[RMSE](#)

### Examples

```
# logistic regression function with one slope and intercept
fn <- function(theta, param) 1 / (1 + exp(-(param[1] + param[2] * theta)))

# sample and population sets
est <- c(-0.4951, 1.1253)
pop <- c(-0.5, 1)

theta <- seq(-10,10,length.out=1000)
plot(theta, fn(theta, pop), type = 'l', col='red', ylim = c(0,1))
lines(theta, fn(theta, est), col='blue', lty=2)

# cumulative result (i.e., standard integral)
IRMSE(est, pop, fn)
```

```
# integrated RMSE result by marginalizing over a N(0,1) distribution
den <- function(theta, mean, sd) dnorm(theta, mean=mean, sd=sd)

IRMSE(est, pop, fn, den, mean=0, sd=1)

# this specification is equivalent to the above
den2 <- function(theta, ...) dnorm(theta, ...)

IRMSE(est, pop, fn, den2, mean=0, sd=1)
```

---

MAE

---

*Compute the mean absolute error*


---

### Description

Computes the average absolute deviation of a sample estimate from the parameter value. Accepts estimate and parameter values, as well as estimate values which are in deviation form.

### Usage

```
MAE(estimate, parameter = NULL, type = "MAE", percent = FALSE, unname = FALSE)
```

### Arguments

|           |  |
|-----------|--|
| estimate  | a numeric vector, matrix/data.frame, or list of parameter estimates. If a vector, the length is equal to the number of replications. If a matrix/data.frame the number of rows must equal the number of replications. list objects will be looped over using the same rules after above after first translating the information into one-dimensional vectors and re-creating the structure upon return                     |
| parameter | a numeric scalar/vector or matrix indicating the fixed parameter values. If a single value is supplied and estimate is a matrix/data.frame then the value will be recycled for each column; otherwise, each element will be associated with each respective column in the estimate input. If NULL, then it will be assumed that the estimate input is in a deviation form (therefore mean(abs(estimate)) will be returned) |
| type      | type of deviation to compute. Can be 'MAE' (default) for the mean absolute error, 'NMSE' for the normalized MAE (MAE / (max(estimate) - min(estimate))), or 'SMSE' for the standardized MAE (MAE / sd(estimate))   |
| percent   | logical; change returned result to percentage by multiplying by 100? Default is FALSE  |
| unname    | logical; apply <code>unname</code> to the results to remove any variable names?  |

### Value

returns a numeric vector indicating the overall mean absolute error in the estimates

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

**See Also**

RMSE

**Examples**

```
pop <- 1
samp <- rnorm(100, 1, sd = 0.5)
MAE(samp, pop)

dev <- samp - pop
MAE(dev)
MAE(samp, pop, type = 'NMAE')
MAE(samp, pop, type = 'SMAE')

# matrix input
mat <- cbind(M1=rnorm(100, 2, sd = 0.5), M2 = rnorm(100, 2, sd = 1))
MAE(mat, parameter = 2)

# same, but with data.frame
df <- data.frame(M1=rnorm(100, 2, sd = 0.5), M2 = rnorm(100, 2, sd = 1))
MAE(df, parameter = c(2,2))

# parameters of the same size
parameters <- 1:10
estimates <- parameters + rnorm(10)
MAE(estimates, parameters)
```

---

manageMessages

*Increase the intensity or suppress the output of an observed message*

---

**Description**

Function provides more nuanced management of known message outputs messages that appear in function calls outside the front-end users control (e.g., functions written in third-party packages). Specifically, this function provides a less nuclear approach than `quiet` and friends, which suppresses all cat and messages raised, and instead allows for specific messages to be raised either

to warnings or, even more extremely, to errors. Note that for messages that are not suppressed the order with which the output and message calls appear in the original function is not retained.

### Usage

```
manageMessages(  
  expr,  
  allow = NULL,  
  message2warning = NULL,  
  message2error = NULL,  
  ...  
)
```

### Arguments

|                 |  |
|-----------------|--|
| expr            | expression to be evaluated (e.g., <code>ret &lt;- myfun(args)</code> ). Function should either be used as a wrapper, such as <code>manageMessages(ret &lt;- myfun(args), ...)</code> or <code>ret &lt;- manageMessages(myfun(args), ...)</code> , or more readably as a pipe, <code>ret &lt;- myfun(args)  &gt; manageMessages(...)</code>   |
| allow           | (optional) a character vector indicating messages that should still appear, while all other messages should remain suppressed. Each supplied message is matched using a <a href="#">grepl</a> expression, so partial matching is supported (though more specific messages are less likely to throw false positives). If <code>NULL</code> , all messages will be suppressed unless they appear in <code>message2error</code> or <code>message2warning</code> |
| message2warning | (optional) Input can be a character vector containing messages that should probably be considered warning messages for the current application instead. Each supplied character vector element is matched using a <a href="#">grepl</a> expression, so partial matching is supported (though more specific messages are less likely to throw false positives).   |
| message2error   | (optional) Input can be a character vector containing known-to-be-severe messages that should be converted to errors for the current application. See <code>message2warning</code> for details.  |
| ...             | additional arguments passed to <a href="#">grepl</a>   |

### Value

returns the original result of `eval(expr)`, with warning messages either left the same, increased to errors, or suppressed (depending on the input specifications)

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)



**See Also**

[manageWarnings](#), [quiet](#)

**Examples**

```
## Not run:

myfun <- function(x, warn=FALSE){
  message('This function is rather chatty')
  cat("It even prints in different output forms!\n")
  message('And even at different ')
  cat(" many times!\n")
  cat("Too many messages can be annoying \n")
  if(warn)
    warning('It may even throw warnings ')
  x
}

out <- myfun(1)
out

# tell the function to shhhh
out <- quiet(myfun(1))
out

# same default behaviour as quiet(), but potential for nuance
out2 <- manageMessages(myfun(1))
identical(out, out2)

# allow some messages to still get printed
out2 <- manageMessages(myfun(1), allow = "many times!")
out2 <- manageMessages(myfun(1), allow = "This function is rather chatty")

# note: . matches single character (regex)
out2 <- manageMessages(myfun(1), allow = c("many times.",
                                           "This function is rather chatty"))

# convert specific message to warning
out3 <- manageMessages(myfun(1), message2warning = "many times!")
identical(out, out3)

# other warnings also get through
out3 <- manageMessages(myfun(1, warn=TRUE), message2warning = "times!")
identical(out, out3)

# convert message to error
manageMessages(myfun(1), message2error = "m... times!")

# multiple message intensity changes
manageMessages(myfun(1),
  message2warning = "It even prints in different output forms",
  message2error = "many times!")
```

```
manageMessages(myfun(1),
  allow = c("This function is rather chatty",
    "Too many messages can be annoying"),
  message2warning = "It even prints in different output forms",
  message2error = "many times!")
```

```
## End(Not run)
```

---

manageWarnings

*Manage specific warning messages*

---

## Description

Function provides more nuanced management of known warning messages that appear in function calls outside the front-end users control (e.g., functions written in third-party packages). Specifically, this function provides a less nuclear approach than [suppressWarnings](#), which suppresses all warning messages rather than those which are known to be innocuous to the current application, or when globally setting `options(warn=2)`, which has the opposite effect of treating all warnings messages as errors in the function executions. To avoid these two extreme behaviors, character vectors can instead be supplied to this function to either leave the raised warnings as-is (default behaviour), raise only specific warning messages to errors, or specify specific warning messages that can be generally be ignored (and therefore suppressed) while allowing new or yet to be discovered warnings to still be raised.

## Usage

```
manageWarnings(expr, warning2error = FALSE, suppress = NULL, ...)
```

## Arguments

|               |   |
|---------------|---|
| expr          | expression to be evaluated (e.g., <code>ret &lt;- myfun(args)</code> ). Function should either be used as a wrapper, such as <code>manageWarnings(ret &lt;- myfun(args), ...)</code> or <code>ret &lt;- manageWarnings(myfun(args), ...)</code> , or more readably as a pipe, <code>ret &lt;- myfun(args)  &gt; manageWarnings(...)</code>  |
| warning2error | logical or character vector to control the conversion of warnings to errors. Setting this input to <code>TRUE</code> will treat all observed warning messages as errors (same behavior as <code>options(warn=2)</code> , though defined on a per expression basis rather than globally), while setting to <code>FALSE</code> (default) will leave all warning messages as-is, retaining the default behavior<br><br>Alternatively, and more useful for specificity reasons, input can be a character vector containing known-to-be-severe warning messages that should be converted to errors. Each supplied character vector element is matched using a <a href="#">grepl</a> expression, so partial matching is supported (though more specific messages are less likely to throw false positives). |

|          |   |
|----------|---|
| suppress | a character vector indicating warning messages that are known to be innocuous a priori and can therefore be suppressed. Each supplied warning message is matched using a <code>grepl</code> expression, so partial matching is supported (though more specific messages are less likely to throw false positives). If NULL, no warning message will be suppressed |
| ...      | additional arguments passed to <code>grepl</code>   |

### Details

In general, global/nuclear behaviour of warning messages should be avoided as they are generally bad practice. On one extreme, when suppressing all warning messages using `suppressWarnings`, potentially important warning messages will become muffled, which can be problematic if the code developer has not become aware of these (now muffled) warnings. Moreover, this can become a long-term sustainability issue when third-party functions that the developer's code depends upon throw new warnings in the future as the code developer will be less likely to become aware of these newly implemented warnings.

On the other extreme, where all warning messages are turned into errors using `options(warn=2)`, innocuous warning messages can and will be (unwantingly) raised to an error. This negatively affects the logical workflow of the developer's functions, where more error messages must now be manually managed (e.g., via `tryCatch`), including the known to be innocuous warning messages as these will now be considered as errors.

To avoid these extremes, front-end users should first make note of the warning messages that have been raised in their prior executions, and organized these messages into vectors of ignorable warnings (least severe), known/unknown warnings that should remain as warnings (even if not known by the code developer yet), and explicit warnings that ought to be considered errors for the current application (most severe). Once collected, these can be passed to the respective `warning2error` argument to increase the intensity of a specific warning raised, or to the `suppress` argument to suppress only the messages that have been deemed ignorable a priori (and therefore allowing all other warning messages to be raised).

### Value

returns the original result of `eval(expr)`, with warning messages either left the same, increased to errors, or suppressed (depending on the input specifications)

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)

### See Also

[manageMessages](#), [quiet](#)

**Examples**

```

## Not run:

fun <- function(warn1=FALSE, warn2=FALSE, warn3=FALSE,
               warn_trailing = FALSE, error=FALSE){
  if(warn1) warning('Message one')
  if(warn2) warning('Message two')
  if(warn3) warning('Message three')
  if(warn_trailing) warning(sprintf('Message with lots of random trailings: %s',
                                   paste0(sample(letters, sample(1:20, 1)), collapse=',')))
  if(error) stop('terminate function call')
  return('Returned from fun()')
}

# normal run (no warnings or errors)
out <- fun()
out

# these are all the same
manageWarnings(out <- fun())
out <- manageWarnings(fun())
out <- fun() |> manageWarnings()

# errors treated normally
fun(error=TRUE)
fun(error=TRUE) |> manageWarnings()

# all warnings/returns treated normally by default
ret1 <- fun(warn1=TRUE)
ret2 <- fun(warn1=TRUE) |> manageWarnings()
identical(ret1, ret2)

# all warnings converted to errors (similar to options(warn=2), but local)
fun(warn1=TRUE) |> manageWarnings(warning2error=TRUE)
fun(warn2=TRUE) |> manageWarnings(warning2error=TRUE)

# Specific warnings treated as errors (others stay as warnings)
# Here, treat first warning message as error but not the second or third
ret <- fun(warn1=TRUE) # warning
ret <- fun(warn1=TRUE) |> manageWarnings("Message one") # now error
ret <- fun(warn2=TRUE) |> manageWarnings("Message one") # still a warning

# multiple warnings raised but not converted as they do not match criteria
fun(warn2=TRUE, warn3=TRUE)
fun(warn2=TRUE, warn3=TRUE) |> manageWarnings("Message one")

# Explicitly convert multiple warning messages, allowing others through.
# This is generally the best use of the function's specificity
fun(warn1=TRUE, warn2=TRUE)
fun(warn1=TRUE) |> # error given either message
  manageWarnings(c("Message one", "Message two"))
fun(warn2=TRUE) |>

```

```

        manageWarnings(c("Message one", "Message two"))

# last warning gets through (left as valid warning)
ret <- fun(warn3=TRUE) |>
        manageWarnings(c("Message one", "Message two"))
ret

# suppress warnings that have only partial matching
fun(warn_trailing=TRUE)
fun(warn_trailing=TRUE)
fun(warn_trailing=TRUE)

# partial match, therefore suppressed
fun(warn_trailing=TRUE) |>
  manageWarnings(suppress="Message with lots of random trailings: ")

# multiple suppress strings
fun(warn_trailing=TRUE) |>
  manageWarnings(suppress=c("Message with lots of random trailings: ",
                            "Suppress this too"))

# could also use .* to catch all remaining characters (finer regex control)
fun(warn_trailing=TRUE) |>
  manageWarnings(suppress="Message with lots of random trailings: .*")

#####
# Combine with quiet() and suppress argument to suppress innocuous messages

fun <- function(warn1=FALSE, warn2=FALSE, warn3=FALSE, error=FALSE){
  message('This function is rather chatty')
  cat("It even prints in different output forms!\n")
  if(warn1) warning('Message one')
  if(warn2) warning('Message two')
  if(warn3) warning('Message three')
  if(error) stop('terminate function call')
  return('Returned from fun()')
}

# normal run (no warnings or errors, but messages)
out <- fun()
out <- quiet(fun()) # using "indoor voice"

# suppress all print messages and warnings (not recommended)
fun(warn2=TRUE) |> quiet()
fun(warn2=TRUE) |> quiet() |> suppressWarnings()

# convert all warning to errors, and keep suppressing messages via quiet()
fun(warn2=TRUE) |> quiet() |> manageWarnings(warning2error=TRUE)

# define tolerable warning messages (only warn1 deemed ignorable)
ret <- fun(warn1=TRUE) |> quiet() |>
  manageWarnings(suppress = 'Message one')
```

```

# all other warnings raised to an error except ignorable ones
fun(warn1=TRUE, warn2=TRUE) |> quiet() |>
  manageWarnings(warning2error=TRUE, suppress = 'Message one')

# only warn2 raised to an error explicitly (warn3 remains as warning)
ret <- fun(warn1=TRUE, warn3=TRUE) |> quiet() |>
  manageWarnings(warning2error = 'Message two',
    suppress = 'Message one')

fun(warn1=TRUE, warn2 = TRUE, warn3=TRUE) |> quiet() |>
  manageWarnings(warning2error = 'Message two',
    suppress = 'Message one')

#####
# Practical example, converting warning into error for model that
# failed to converged normally

library(lavaan)

## The industrialization and Political Democracy Example
## Bollen (1989), page 332
model <- '
  # latent variable definitions
  ind60 =~ x1 + x2 + x3
  dem60 =~ y1 + a*y2 + b*y3 + c*y4
  dem65 =~ y5 + a*y6 + b*y7 + c*y8

  # regressions
  dem60 ~ ind60
  dem65 ~ ind60 + dem60

  # residual correlations
  y1 ~~ y5
  y2 ~~ y4 + y6
  y3 ~~ y7
  y4 ~~ y8
  y6 ~~ y8
  ,

# throws a warning
fit <- sem(model, data = PoliticalDemocracy, control=list(iter.max=60))

# for a simulation study, often better to treat this as an error
fit <- sem(model, data = PoliticalDemocracy, control=list(iter.max=60)) |>
  manageWarnings(warning2error = "the optimizer warns that a solution has NOT been found!")

## End(Not run)

```

---

|       |  |
|-------|--|
| MSRSE | <i>Compute the relative performance behavior of collections of standard errors</i> |
|-------|--|

---

### Description

The mean-square relative standard error (MSRSE) compares standard error estimates to the standard deviation of the respective parameter estimates. Values close to 1 indicate that the behavior of the standard errors closely matched the sampling variability of the parameter estimates.

### Usage

```
MSRSE(SE, SD, percent = FALSE, unname = FALSE)
```

### Arguments

|         |   |
|---------|---|
| SE      | a numeric scalar/vector indicating the average standard errors across the replications, or a matrix of collected standard error estimates themselves to be used to compute the average standard errors. Each column/element in this input corresponds to the column/element in SD |
| SD      | a numeric scalar/vector indicating the standard deviation across the replications, or a matrix of collected parameter estimates themselves to be used to compute the standard deviations. Each column/element in this input corresponds to the column/element in SE               |
| percent | logical; change returned result to percentage by multiplying by 100? Default is FALSE   |
| unname  | logical; apply <code>unname</code> to the results to remove any variable names?   |

### Details

Mean-square relative standard error (MSRSE) is expressed as

$$MSRSE = \frac{E(SE(\psi)^2)}{SD(\psi)^2} = \frac{1/R * \sum_{r=1}^R SE(\psi_r)^2}{SD(\psi)^2}$$

where  $SE(\psi_r)$  represents the estimate of the standard error at the  $r$ th simulation replication, and  $SD(\psi)$  represents the standard deviation estimate of the parameters across all  $R$  replications. Note that  $SD(\psi)^2$  is used, which corresponds to the variance of  $\psi$ .

### Value

returns a vector of ratios indicating the relative performance of the standard error estimates to the observed parameter standard deviation. Values less than 1 indicate that the standard errors were larger than the standard deviation of the parameters (hence, the SEs are interpreted as more conservative), while values greater than 1 were smaller than the standard deviation of the parameters (i.e., more liberal SEs)

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

**Examples**

```
Generate <- function(condition, fixed_objects) {
  X <- rep(0:1, each = 50)
  y <- 10 + 5 * X + rnorm(100, 0, .2)
  data.frame(y, X)
}

Analyse <- function(condition, dat, fixed_objects) {
  mod <- lm(y ~ X, dat)
  so <- summary(mod)
  ret <- c(SE = so$coefficients[,"Std. Error"],
          est = so$coefficients[,"Estimate"])
  ret
}

Summarise <- function(condition, results, fixed_objects) {
  MSRSE(SE = results[,1:2], SD = results[,3:4])
}

results <- runSimulation(replications=500, generate=Generate,
                        analyse=Analyse, summarise=Summarise)
results
```

---

 nc

*Auto-named Concatenation of Vector or List*

---

**Description**

This is a wrapper to the function `c`, however names the respective elements according to their input object name. For this reason, nesting `nc()` calls is not recommended (joining independent `nc()` calls via `c()` is however reasonable).

**Usage**

```
nc(..., use.names = FALSE, error.on.duplicate = TRUE)
```



## Arguments

... objects to be concatenated

use.names logical indicating if names should be preserved (unlike `c`, default is FALSE)

error.on.duplicate logical; if the same object name appears in the returning object should an error be thrown? Default is TRUE

## References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

## Examples

```
A <- 1
B <- 2
C <- 3

names(C) <- 'LetterC'

# compare the following
c(A, B, C) # unnamed

nc(A, B, C) # named
nc(this=A, B, C) # respects override named (same as c() )
nc(this=A, B, C, use.names = TRUE) # preserve original name

## Not run:
# throws errors if names not unique
nc(this=A, this=B, C)
nc(LetterC=A, B, C, use.names=TRUE)

## End(Not run)

# poor input choice names
nc(t.test(c(1:2))$p.value, t.test(c(3:4))$p.value)

# better to explicitly provide name
nc(T1 = t.test(c(1:2))$p.value,
    T2 = t.test(c(3:4))$p.value)

# vector of unnamed inputs
A <- c(5,4,3,2,1)
B <- c(100, 200)

nc(A, B, C) # A's and B's numbered uniquely
c(A, B, C) # compare
```

```

nc(beta=A, B, C) # replacement of object name

# retain names attributes (but append object name, when appropriate)
names(A) <- letters[1:5]
nc(A, B, C)
nc(beta=A, B, C)
nc(A, B, C, use.names=TRUE)

# mix and match if some named elements work while others do not
c( nc(A, B, use.names=TRUE), nc(C))

## Not run:
# error, 'b' appears twice
names(B) <- c('b', 'b2')
nc(A, B, C, use.names=TRUE)

## End(Not run)

# List input
A <- list(1)
B <- list(2:3)
C <- list('C')

names(C) <- 'LetterC'

# compare the following
c(A, B, C) # unnamed

nc(A, B, C) # named
nc(this=A, B, C) # respects override named (same as c() and list() )
nc(this=A, B, C, use.names = TRUE) # preserve original name

```

## Description

The function PBA searches a specified interval for a root (i.e., zero) of the function  $f(x)$  with respect to its first argument. However, this function differs from deterministic cousins such as [uniroot](#) in that  $f$  may contain stochastic error components, and instead provides a Bayesian interval where the root is likely to lie. Note that it is assumed that  $E[f(x)]$  is non-decreasing in  $x$  and that the root is between the search interval (evaluated approximately when `check.interval=TRUE`). See Waeber, Frazier, and Henderson (2013) for details.

## Usage

```

PBA(
  f,

```

```

    interval,
    ...,
    p = 0.6,
    integer = FALSE,
    tol = if (integer) 0.01 else 1e-04,
    maxiter = 300L,
    miniter = 100L,
    wait.time = NULL,
    f.prior = NULL,
    resolution = 10000L,
    check.interval = TRUE,
    check.interval.only = FALSE,
    verbose = TRUE
)

## S3 method for class 'PBA'
print(x, ...)

## S3 method for class 'PBA'
plot(x, type = "posterior", main = "Probabilistic Bisection Posterior", ...)

```

### Arguments

|                             |  |
|-----------------------------|--|
| <code>f</code>              | noisy function for which the root is sought  |
| <code>interval</code>       | a vector containing the end-points of the interval to be searched for the root   |
| <code>...</code>            | additional named arguments to be passed to <code>f</code>  |
| <code>p</code>              | assumed constant for probability of correct responses (must be > 0.5)  |
| <code>integer</code>        | logical; should the values of the root be considered integer or numeric? The former uses a discreet grid to track the updates, while the latter currently creates a grid with resolution points  |
| <code>tol</code>            | tolerance criteria for convergence based on average of the $f(x)$ evaluations  |
| <code>maxiter</code>        | the maximum number of iterations (default 300)   |
| <code>miniter</code>        | minimum number of iterations (default 100)   |
| <code>wait.time</code>      | (optional) instead of terminating after specific estimate criteria are satisfied (e.g., <code>tol</code> ), terminate after a specific wait time. Input is specified either as a numeric vector in seconds or as a character vector to be formatted by <code>timeFormater</code> . Note that users should increase the number of <code>maxiter</code> as well so that termination can occur if either the maximum iterations are satisfied or the specified wait time has elapsed (whichever occurs first) |
| <code>f.prior</code>        | density function indicating the likely location of the prior (e.g., if root is within [0,1] then <code>dunif</code> works, otherwise custom functions will be required)  |
| <code>resolution</code>     | constant indicating the number of equally spaced grid points to track when <code>integer = FALSE</code> .  |
| <code>check.interval</code> | logical; should an initial check be made to determine whether $f(\text{interval}[1L])$ and $f(\text{interval}[2L])$ have opposite signs? Default is TRUE   |

|                                  |  |
|----------------------------------|--|
| <code>check.interval.only</code> | logical; return only TRUE or FALSE to test whether there is a likely root given interval? Setting this to TRUE can be useful when you are unsure about the root location interval and may want to use a higher replication input from <a href="#">SimSolve</a> |
| <code>verbose</code>             | logical; should the iterations and estimate be printed to the console?   |
| <code>x</code>                   | an object of class PBA   |
| <code>type</code>                | type of plot to draw for PBA object. Can be either 'posterior' or 'history' to plot the PBA posterior distribution or the mediation iteration history  |
| <code>main</code>                | plot title   |

## References

- Horstein, M. (1963). Sequential transmission using noiseless feedback. *IEEE Trans. Inform. Theory*, 9(3):136-143.
- Waeber, R., Frazier, P. I. & Henderson, S. G. (2013). Bisection Search with Noisy Responses. *SIAM Journal on Control and Optimization, Society for Industrial & Applied Mathematics (SIAM)*, 51, 2261-2279.

## See Also

[uniroot](#), [RobbinsMonro](#)

## Examples

```
# find x that solves f(x) - b = 0 for the following
f.root <- function(x, b = .6) 1 / (1 + exp(-x)) - b
f.root(.3)

xs <- seq(-3,3, length.out=1000)
plot(xs, f.root(xs), type = 'l', ylab = "f(x)", xlab='x', las=1)
abline(h=0, col='red')

retuni <- uniroot(f.root, c(0,1))
retuni
abline(v=retuni$root, col='blue', lty=2)

# PBA without noisy root
retpba <- PBA(f.root, c(0,1))
retpba
retpba$root
plot(retpba)
plot(retpba, type = 'history')

# Same problem, however root function is now noisy. Hence, need to solve
# fhat(x) - b + e = 0, where E(e) = 0
f.root_noisy <- function(x) 1 / (1 + exp(-x)) - .6 + rnorm(1, sd=.02)
sapply(rep(.3, 10), f.root_noisy)

# uniroot "converges" unreliably
```

```

set.seed(123)
uniroot(f.root_noisy, c(0,1))$root
uniroot(f.root_noisy, c(0,1))$root
uniroot(f.root_noisy, c(0,1))$root

# probabilistic bisection provides better convergence
retpba.noise <- PBA(f.root_noisy, c(0,1))
retpba.noise
plot(retpba.noise)
plot(retpba.noise, type = 'history')

## Not run:
# ignore termination criteria and instead run for 30 seconds or 30000 iterations
retpba.noise_30sec <- PBA(f.root_noisy, c(0,1), wait.time = "0:30", maxiter=30000)
retpba.noise_30sec

## End(Not run)

```

---

quiet

*Suppress verbose function messages*


---

## Description

This function is used to suppress information printed from external functions that make internal use of `link{message}` and `cat`, which provide information in interactive R sessions. For simulations, the session is not interactive, and therefore this type of output should be suppressed. For similar behaviour for suppressing warning messages, see `link{manageWarnings}`.

## Usage

```
quiet(..., cat = TRUE, keep = FALSE, attr.name = "quiet.messages")
```

## Arguments

|                        |  |
|------------------------|--|
| <code>...</code>       | the functional expression to be evaluated  |
| <code>cat</code>       | logical; also capture calls from <code>cat</code> ? If FALSE only <code>message</code> will be suppressed  |
| <code>keep</code>      | logical; return a character vector of the messages/concatenate and print strings as an attribute to the resulting object from <code>expr(...)</code> ? |
| <code>attr.name</code> | attribute name to use when <code>keep = TRUE</code>  |

## References

- Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248
- Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

**See Also**[manageWarnings](#)**Examples**

```

myfun <- function(x, warn=FALSE){
  message('This function is rather chatty')
  cat("It even prints in different output forms!\n")
  message('And even at different....')
  cat("...times!\n")
  if(warn)
    warning('It may even throw warnings!')
  x
}

out <- myfun(1)
out

# tell the function to shhhh
out <- quiet(myfun(1))
out

# which messages are suppressed? Extract stored attribute
out <- quiet(myfun(1), keep = TRUE)
attr(out, 'quiet.messages')

# Warning messages still get through (see manageWarnings(suppress))
# for better alternative than using suppressWarnings())
out2 <- myfun(2, warn=TRUE) |> quiet() # warning gets through
out2

# suppress warning message explicitly, allowing others to be raised if present
myfun(2, warn=TRUE) |> quiet() |>
  manageWarnings(suppress='It may even throw warnings!')

```

---

RAB

---

*Compute the relative absolute bias of multiple estimators*


---

**Description**

Computes the relative absolute bias given the bias estimates for multiple estimators.

**Usage**

```
RAB(x, percent = FALSE, unname = FALSE)
```

**Arguments**

|         |   |
|---------|---|
| x       | a numeric vector of bias estimates (see <a href="#">bias</a> ), where the first element will be used as the reference |
| percent | logical; change returned result to percentage by multiplying by 100? Default is FALSE                                 |
| unname  | logical; apply <a href="#">unname</a> to the results to remove any variable names?                                    |

**Value**

returns a vector of absolute bias ratios indicating the relative bias effects compared to the first estimator. Values less than 1 indicate better bias estimates than the first estimator, while values greater than 1 indicate worse bias than the first estimator

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. [doi:10.1080/10691898.2016.1246953](https://doi.org/10.1080/10691898.2016.1246953)

**Examples**

```
pop <- 1
samp1 <- rnorm(5000, 1)
bias1 <- bias(samp1, pop)
samp2 <- rnorm(5000, 1)
bias2 <- bias(samp2, pop)

RAB(c(bias1, bias2))
RAB(c(bias1, bias2), percent = TRUE) # as a percentage
```

---

rbind.SimDesign

*Combine two separate SimDesign objects by row*

---

**Description**

This function combines two Monte Carlo simulations executed by SimDesign's [runSimulation](#) function which, for all intents and purposes, could have been executed in a single run. This situation arises when a simulation has been completed, however the Design object was later modified to include more levels in the defined simulation factors. Rather than re-executing the previously completed simulation combinations, only the new combinations need to be evaluated into a different object and then rbind together to create the complete object combinations.

**Usage**

```
## S3 method for class 'SimDesign'
rbind(...)
```

**Arguments**

```
...          two or more SimDesign objects that should be combined by rows
```

**Value**

same object that is returned by [runSimulation](#)

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. [doi:10.1080/10691898.2016.1246953](https://doi.org/10.1080/10691898.2016.1246953)

**Examples**

```
## Not run:

# modified example from runSimulation()

Design <- createDesign(N = c(10, 20),
                      SD = c(1, 2))

Generate <- function(condition, fixed_objects) {
  dat <- with(condition, rnorm(N, 10, sd=SD))
  dat
}

Analyse <- function(condition, dat, fixed_objects) {
  ret <- mean(dat) # mean of the sample data vector
  ret
}

Summarise <- function(condition, results, fixed_objects) {
  ret <- c(mu=mean(results), SE=sd(results)) # mean and SD summary of the sample means
  ret
}

Final1 <- runSimulation(design=Design, replications=1000,
                      generate=Generate, analyse=Analyse, summarise=Summarise)

Final1
```



```

###
# later decide that N = 30 should have also been investigated. Rather than
# running the following object ...
newDesign <- createDesign(N = c(10, 20, 30),
                          SD = c(1, 2))

# ... only the new subset levels are executed to save time
subDesign <- subset(newDesign, N == 30)
subDesign

Final2 <- runSimulation(design=subDesign, replications=1000,
                       generate=Generate, analyse=Analyse, summarise=Summarise)
Final2

# glue results together by row into one object as though the complete 'Design'
# object were run all at once
Final <- rbind(Final1, Final2)
Final

summary(Final)

## End(Not run)

```

---

RD

---

*Compute the relative difference*


---

### Description

Computes the relative difference statistic of the form  $(\text{est} - \text{pop}) / \text{pop}$ , which is equivalent to the form  $\text{est}/\text{pop} - 1$ . If matrices are supplied then an equivalent matrix variant will be used of the form  $(\text{est} - \text{pop}) * \text{solve}(\text{pop})$ . Values closer to 0 indicate better relative parameter recovery. Note that for single variable inputs this is equivalent to `bias(..., type = 'relative')`.

### Usage

```
RD(est, pop, as.vector = TRUE, unname = FALSE)
```

### Arguments

|                        |  |
|------------------------|--|
| <code>est</code>       | a numeric vector, matrix/data.frame, or list containing the parameter estimates                                      |
| <code>pop</code>       | a numeric vector or matrix containing the true parameter values. Must be of comparable dimension to <code>est</code> |
| <code>as.vector</code> | logical; always wrap the result in a <code>as.vector</code> function before returning?                               |
| <code>unname</code>    | logical; apply <code>unname</code> to the results to remove any variable names?                                      |

**Value**

returns a vector or matrix depending on the inputs and whether `as.vector` was used

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

**Examples**

```
# vector
pop <- seq(1, 100, length.out=9)
est1 <- pop + rnorm(9, 0, .2)
(rds <- RD(est1, pop))
summary(rds)

# matrix
pop <- matrix(c(1:8, 10), 3, 3)
est2 <- pop + rnorm(9, 0, .2)
RD(est2, pop, as.vector = FALSE)
(rds <- RD(est2, pop))
summary(rds)
```

---

RE

*Compute the relative efficiency of multiple estimators*

---

**Description**

Computes the relative efficiency given the RMSE (default) or MSE values for multiple estimators.

**Usage**

```
RE(x, MSE = FALSE, percent = FALSE, unname = FALSE)
```

**Arguments**

|         |   |
|---------|---|
| x       | a numeric vector of root mean square error values (see <a href="#">RMSE</a> ), where the first element will be used as the reference. Otherwise, the object could contain MSE values if the flag <code>MSE = TRUE</code> is also included |
| MSE     | logical; are the input value mean squared errors instead of root mean square errors?  |
| percent | logical; change returned result to percentage by multiplying by 100? Default is FALSE   |
| unname  | logical; apply <a href="#">unname</a> to the results to remove any variable names?  |

**Value**

returns a vector of variance ratios indicating the relative efficiency compared to the first estimator. Values less than 1 indicate better efficiency than the first estimator, while values greater than 1 indicate worse efficiency than the first estimator

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. [doi:10.1080/10691898.2016.1246953](https://doi.org/10.1080/10691898.2016.1246953)

**Examples**

```
pop <- 1
samp1 <- rnorm(100, 1, sd = 0.5)
RMSE1 <- RMSE(samp1, pop)
samp2 <- rnorm(100, 1, sd = 1)
RMSE2 <- RMSE(samp2, pop)

RE(c(RMSE1, RMSE2))
RE(c(RMSE1, RMSE2), percent = TRUE) # as a percentage

# using MSE instead
mse <- c(RMSE1, RMSE2)^2
RE(mse, MSE = TRUE)
```

---

rejectionSampling      *Rejection sampling (i.e., accept-reject method)*

---

### Description

This function supports the rejection sampling (i.e., accept-reject) approach to drawing values from seemingly difficult (probability) density functions by sampling values from more manageable proxy distributions.

### Usage

```
rejectionSampling(
  n,
  df,
  dg,
  rg,
  M,
  method = "optimize",
  interval = NULL,
  logfun = FALSE,
  maxM = 1e+05,
  parstart = rg(1L),
  ESRS_Mstart = 1.0001
)
```

### Arguments

|    |   |
|----|---|
| n  | number of samples to draw   |
| df | the desired (potentially un-normed) density function to draw independent samples from. Must be in the form of a function with a single input corresponding to the values sampled from rg. Function is assumed to be vectorized (if not, see <a href="#">Vectorize</a> )   |
| dg | the proxy (potentially un-normed) density function to draw samples from in lieu of drawing samples from df. The support for this density function should be the same as df (i.e., when $df(x) > 0$ then $dg(x) > 0$ ). Must be in the form of a function with a single input corresponding to the values sampled from rg. Function is assumed to be vectorized (if not, see <a href="#">Vectorize</a> )                 |
| rg | the proxy random number generation function, associated with dg, used to draw proposal samples from. Must be in the form of a function with a single input corresponding to the number of values to draw, while the output can either be a vector or a matrix (if a matrix, each independent observation must be stored in a unique row). Function is assumed to be vectorized (if not, see <a href="#">Vectorize</a> ) |
| M  | the upper-bound of the ratio of probability density functions to help minimize the number of discarded draws and define the corresponding rescaled proposal envelope. When missing, M is computed internally by finding a reasonable maximum of $\log(df(x)) - \log(dg(x))$ , and this value is returned to the console.  |

|                          |  |
|--------------------------|--|
|                          | When both <code>df</code> and <code>dg</code> are true probability density functions (i.e., integrate to 1) the acceptance probability is equal to $1/M$   |
| <code>method</code>      | when <code>M</code> is missing, the optimization of <code>M</code> is done either by finding the mode of the log-density values (" <code>optimize</code> ") or by using the "Empirical Supremum Rejection Sampling" method (" <code>ESRS</code> ")   |
| <code>interval</code>    | when <code>M</code> is missing, for univariate density function draws, the interval to search within via <code>optimize</code> . If not specified, a sample of 5000 values from the <code>rg</code> function definition will be collected, and the min/max will be obtained via this random sample                       |
| <code>logfuns</code>     | logical; have the <code>df</code> and <code>dg</code> function been written so as to return log-densities instead of the original densities? The <code>FALSE</code> default assumes the original densities are returned (use <code>TRUE</code> when higher accuracy is required when generating each density definition) |
| <code>maxM</code>        | logical; if when optimizing <code>M</code> the value is greater than this cut-off then stop; ampler would likelihood be too efficient, or optimization is failing  |
| <code>parstart</code>    | starting value vector for optimization of <code>M</code> in multidimensional distributions   |
| <code>ESRS_Mstart</code> | starting <code>M</code> value for the <code>ESRS</code> algorithm  |

### Details

The accept-reject algorithm is a flexible approach to obtaining i.i.d.'s from a difficult to sample from (probability) density function where either the transformation method fails or inverse transform method is difficult to manage. The algorithm does so by sampling from a more "well-behaved" proxy distribution (with identical support, up to some proportionality constant  $M$  that reshapes the proposal density to envelope the target density), and accepts the draws if they are likely within the target density. Hence, the closer the shape of  $dg(x)$  is to the desired  $df(x)$ , the more likely the draws are to be accepted; otherwise, many iterations of the accept-reject algorithm may be required, which decreases the computational efficiency.

### Value

returns a vector or matrix of draws (corresponding to the output class from `rg`) from the desired `df`

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

- Caffo, B. S., Booth, J. G., and Davison, A. C. (2002). Empirical supremum rejection sampling. *Biometrika*, 89, 745–754.
- Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the `SimDesign` Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248
- Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

**Examples**

```

## Not run:

# Generate  $X \sim \text{beta}(a,b)$ , where  $a$  and  $b$  are  $a = 2.7$  and  $b = 6.3$ ,
# and the support is  $Y \sim \text{Unif}(0,1)$ 
dfn <- function(x) dbeta(x, shape1 = 2.7, shape2 = 6.3)
dgn <- function(x) dunif(x, min = 0, max = 1)
rgn <- function(n) runif(n, min = 0, max = 1)

# when df and dg both integrate to 1, acceptance probability = 1/M
M <- rejectionSampling(df=dfn, dg=dgn, rg=rgn)
M
dat <- rejectionSampling(10000, df=dfn, dg=dgn, rg=rgn, M=M)
hist(dat, 100)
hist(rbeta(10000, 2.7, 6.3), 100) # compare

# obtain empirical estimate of M via ESRS method
M <- rejectionSampling(1000, df=dfn, dg=dgn, rg=rgn, method='ESRS')
M

# generate using better support function (here,  $Y \sim \text{beta}(2,6)$ ),
# and use log setup in initial calls (more numerically accurate)
dfn <- function(x) dbeta(x, shape1 = 2.7, shape2 = 6.3, log = TRUE)
dgn <- function(x) dbeta(x, shape1 = 2, shape2 = 6, log = TRUE)
rgn <- function(n) rbeta(n, shape1 = 2, shape2 = 6)
M <- rejectionSampling(df=dfn, dg=dgn, rg=rgn, logfuns=TRUE) # better M
M

## Alternative estimation of M
## M <- rejectionSampling(10000, df=dfn, dg=dgn, rg=rgn, logfuns=TRUE,
## method='ESRS')
dat <- rejectionSampling(10000, df=dfn, dg=dgn, rg=rgn, M=M, logfuns=TRUE)
hist(dat, 100)

#-----
# sample from wonky (and non-normalized) density function, like below
dfn <- function(x){
  ret <- numeric(length(x))
  ret[x <= .5] <- dnorm(x[x <= .5])
  ret[x > .5] <- dnorm(x[x > .5]) + dchisq(x[x > .5], df = 2)
  ret
}
y <- seq(-5,5, length.out = 1000)
plot(y, dfn(y), type = 'l', main = "Function to sample from")

# choose dg/rg functions that have support within the range  $[-\text{inf}, \text{inf}]$ 
rgn <- function(n) rnorm(n, sd=4)
dgn <- function(x) dnorm(x, sd=4)

## example M height from above graphic
## (M selected using ESRS to help stochastically avoid local mins)
M <- rejectionSampling(10000, df=dfn, dg=dgn, rg=rgn, method='ESRS')

```

```

M
lines(y, dgn(y)*M, lty = 2)
dat <- rejectionSampling(10000, df=dfn, dg=dgn, rg=rgn, M=M)
hist(dat, 100, prob=TRUE)

# true density (normalized)
C <- integrate(dfn, -Inf, Inf)$value
ndfn <- function(x) dfn(x) / C
curve(ndfn, col='red', lwd=2, add=TRUE)

#-----
# multivariate distribution
dfn <- function(x) sum(log(c(dnorm(x[1]) + dchisq(x[1], df = 5),
                           dnorm(x[2], -1, 2))))
rgn <- function(n) c(rnorm(n, sd=3), rnorm(n, sd=3))
dgn <- function(x) sum(log(c(dnorm(x[1], sd=3), dnorm(x[1], sd=3))))

# M <- rejectionSampling(df=dfn, dg=dgn, rg=rgn, logfuns=TRUE)
dat <- rejectionSampling(5000, df=dfn, dg=dgn, rg=rgn, M=4.6, logfuns=TRUE)
hist(dat[,1], 30)
hist(dat[,2], 30)
plot(dat)

## End(Not run)

```

---

reSummarise

*Run a summarise step for results that have been saved to the hard drive*


---

## Description

When `runSimulation()` uses the option `save_results = TRUE` the R replication results from the Generate-Analyse functions are stored to the hard drive. As such, additional summarise components may be required at a later time, whereby the respective `.rds` files must be read back into R to be summarised. This function performs the reading of these files, application of a provided summarise function, and final collection of the respective results.

## Usage

```

reSummarise(
  summarise,
  dir = NULL,
  files = NULL,
  results = NULL,
  Design = NULL,
  fixed_objects = NULL,

```

```

boot_method = "none",
boot_draws = 1000L,
CI = 0.95,
prefix = "results-row"
)

```

### Arguments

|               |  |
|---------------|--|
| summarise     | a summarise function to apply to the read-in files. See <a href="#">runSimulation</a> for details  |
| dir           | directory pointing to the .rds files to be read-in that were saved from <code>runSimulation(..., save_results=TRUE)</code> . If NULL, it is assumed the current working directory contains the .rds files  |
| files         | (optional) names of files to read-in. If NULL all files located within dir will be used  |
| results       | (optional) the results of <a href="#">runSimulation</a> when no summarise function was provided. Can be either a tibble or matrix (indicating that exactly one design condition was evaluated), or a list of matrix/tibble objects indicating that multiple conditions were performed with no summarise evaluation.<br>Alternatively, if <code>store_results = TRUE</code> in the <code>runSimulation()</code> execution then the final SimDesign object may be passed, where the generate-analyse information will be extracted from the object instead |
| Design        | (optional) if results input used, and design condition information important in the summarise step, then the original design object from <a href="#">runSimulation</a> should be included  |
| fixed_objects | (optional) see <a href="#">runSimulation</a> for details   |
| boot_method   | method for performing non-parametric bootstrap confidence intervals for the respective meta-statistics computed by the Summarise function. See <a href="#">runSimulation</a> for details   |
| boot_draws    | number of non-parametric bootstrap draws to sample for the summarise function after the generate-analyse replications are collected. Default is 1000   |
| CI            | bootstrap confidence interval level (default is 95%)   |
| prefix        | character indicating prefix used for stored files  |

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

- Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248
- Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953



**Examples**

```

Design <- createDesign(N = c(10, 20, 30))

Generate <- function(condition, fixed_objects) {
  dat <- with(condition, rnorm(N, 10, 5)) # distributed N(10, 5)
  dat
}

Analyse <- function(condition, dat, fixed_objects) {
  ret <- c(mean=mean(dat), median=median(dat)) # mean/median of sample data
  ret
}

Summarise <- function(condition, results, fixed_objects){
  colMeans(results)
}

## Not run:
# run the simulation
runSimulation(design=Design, replications=50,
              generate=Generate, analyse=Analyse,
              summarise=Summarise, save_results=TRUE,
              save_details = list(save_results_dirname='simresults'))

res <- reSummarise(Summarise, dir = 'simresults/')
res

Summarise2 <- function(condition, results, fixed_objects){
  ret <- c(mean_ests=colMeans(results), SE=colSDs(results))
  ret
}

res2 <- reSummarise(Summarise2, dir = 'simresults/')
res2

SimClean(dir='simresults/')

## End(Not run)

###
# Similar, but with results stored within the final object

res <- runSimulation(design=Design, replications=50, store_results = TRUE,
                    generate=Generate, analyse=Analyse, summarise=Summarise)
res

# same summarise but with bootstrapping
res2 <- reSummarise(Summarise, results = res, boot_method = 'basic')
res2

```

---

rHeadrick

*Generate non-normal data with Headrick's (2002) method*


---

### Description

Generate multivariate non-normal distributions using the fifth-order polynomial method described by Headrick (2002).

### Usage

```
rHeadrick(
  n,
  mean = rep(0, nrow(sigma)),
  sigma = diag(length(mean)),
  skew = rep(0, nrow(sigma)),
  kurt = rep(0, nrow(sigma)),
  gam3 = NaN,
  gam4 = NaN,
  return_coefs = FALSE,
  coefs = NULL,
  control = list(trace = FALSE, max.ntry = 15, obj.tol = 1e-10, n.valid.sol = 1)
)
```

### Arguments

|              |   |
|--------------|---|
| n            | number of samples to draw   |
| mean         | a vector of k elements for the mean of the variables  |
| sigma        | desired k x k covariance matrix between bivariate non-normal variables  |
| skew         | a vector of k elements for the skewness of the variables  |
| kurt         | a vector of k elements for the kurtosis of the variables  |
| gam3         | (optional) explicitly supply the gamma 3 value? Default computes this internally  |
| gam4         | (optional) explicitly supply the gamma 4 value? Default computes this internally  |
| return_coefs | logical; return the estimated coefficients only? See below regarding why this is useful.  |
| coefs        | (optional) supply previously estimated coefficients? This is useful when there must be multiple data sets drawn and will avoid repetitive computations. Must be the object returned after passing return_coefs = TRUE |
| control      | a list of control parameters when locating the polynomial coefficients  |

### Details

This function is primarily a wrapper for the code written by Oscar L. Olvera Astivia (last edited Feb 26, 2015) with some modifications (e.g., better starting values for the Newton optimizer, passing previously saved coefs, etc).



---

rint *Generate integer values within specified range*

---

### Description

Efficiently generate positive and negative integer values with (default) or without replacement. This function is mainly a wrapper to the `sample.int` function (which itself is much more efficient integer sampler than the more general `sample`), however is intended to work with both positive and negative integer ranges since `sample.int` only returns positive integer values that must begin at 1L.

### Usage

```
rint(n, min, max, replace = TRUE, prob = NULL)
```

### Arguments

|         |  |
|---------|--|
| n       | number of samples to draw  |
| min     | lower limit of the distribution. Must be finite  |
| max     | upper limit of the distribution. Must be finite  |
| replace | should sampling be with replacement?   |
| prob    | a vector of probability weights for obtaining the elements of the vector being sampled |

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

### Examples

```
set.seed(1)

# sample 1000 integer values within 20 to 100
x <- rint(1000, min = 20, max = 100)
summary(x)

# sample 1000 integer values within 100 to 10 billion
x <- rint(1000, min = 100, max = 1e8)
summary(x)
```

```
# compare speed to sample()
system.time(x <- rint(1000, min = 100, max = 1e8))
system.time(x2 <- sample(100:1e8, 1000, replace = TRUE))

# sample 1000 integer values within -20 to 20
x <- rint(1000, min = -20, max = 20)
summary(x)
```

---

rinvWishart

*Generate data with the inverse Wishart distribution*

---

### Description

Function generates data in the form of symmetric matrices from the inverse Wishart distribution given a covariance matrix and degrees of freedom.

### Usage

```
rinvWishart(n = 1, df, sigma)
```

### Arguments

|       |   |
|-------|---|
| n     | number of matrix observations to generate. By default n = 1, which returns a single symmetric matrix. If n > 1 then a list of n symmetric matrices are returned instead |
| df    | degrees of freedom  |
| sigma | positive definite covariance matrix   |

### Value

a numeric matrix with columns equal to `ncol(sigma)` when n = 1, or a list of n matrices with the same properties

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

### See Also

[runSimulation](#)

**Examples**

```
# random inverse Wishart matrix given variances [3,6], covariance 2, and df=15
sigma <- matrix(c(3,2,2,6), 2, 2)
x <- rinwWishart(sigma = sigma, df = 15)
x

# list of matrices
x <- rinwWishart(20, sigma = sigma, df = 15)
x
```

---

 rmgh

---

*Generate data with the multivariate g-and-h distribution*


---

**Description**

Generate non-normal distributions using the multivariate g-and-h distribution. Can be used to generate several different classes of univariate and multivariate distributions.

**Usage**

```
rmgh(n, g, h, mean = rep(0, length(g)), sigma = diag(length(mean)))
```

**Arguments**

|       |  |
|-------|--|
| n     | number of samples to draw  |
| g     | the g parameter(s) which control the skew of a distribution in terms of both direction and magnitude                     |
| h     | the h parameter(s) which control the tail weight or elongation of a distribution and is positively related with kurtosis |
| mean  | a vector of k elements for the mean of the variables   |
| sigma | desired k x k covariance matrix between bivariate non-normal variables   |

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. [doi:10.1080/10691898.2016.1246953](https://doi.org/10.1080/10691898.2016.1246953)

**Examples**

```

set.seed(1)

# univariate
norm <- rmgh(10000,1e-5,0)
hist(norm)

skew <- rmgh(10000,1/2,0)
hist(skew)

neg_skew_platykurtic <- rmgh(10000,-1,-1/2)
hist(neg_skew_platykurtic)

# multivariate
sigma <- matrix(c(2,1,1,4), 2)
mean <- c(-1, 1)
twovar <- rmgh(10000, c(-1/2, 1/2), c(0,0),
              mean=mean, sigma=sigma)
hist(twovar[,1])
hist(twovar[,2])
plot(twovar)

```

---

RMSE

---

*Compute the (normalized) root mean square error*


---

**Description**

Computes the average deviation (root mean square error; also known as the root mean square deviation) of a sample estimate from the parameter value. Accepts estimate and parameter values, as well as estimate values which are in deviation form.

**Usage**

```

RMSE(
  estimate,
  parameter = NULL,
  type = "RMSE",
  MSE = FALSE,
  percent = FALSE,
  unname = FALSE
)

```

```

RMSD(
  estimate,
  parameter = NULL,
  type = "RMSE",
  MSE = FALSE,

```

```

percent = FALSE,
unname = FALSE
)

```

### Arguments

|           |  |
|-----------|--|
| estimate  | a numeric vector, <code>matrix/data.frame</code> , or list of parameter estimates. If a vector, the length is equal to the number of replications. If a <code>matrix/data.frame</code> , the number of rows must equal the number of replications. list objects will be looped over using the same rules after above after first translating the information into one-dimensional vectors and re-creating the structure upon return  |
| parameter | a numeric scalar/vector indicating the fixed parameter values. If a single value is supplied and <code>estimate</code> is a <code>matrix/data.frame</code> then the value will be recycled for each column; otherwise, each element will be associated with each respective column in the <code>estimate</code> input. If <code>NULL</code> then it will be assumed that the <code>estimate</code> input is in a deviation form (therefore <code>sqrt(mean(estimate^2))</code> will be returned) |
| type      | type of deviation to compute. Can be 'RMSE' (default) for the root mean square-error, 'NRMSE' for the normalized RMSE ( $RMSE / (\max(\text{estimate}) - \min(\text{estimate}))$ ), 'SRMSE' for the standardized RMSE ( $RMSE / \text{sd}(\text{estimate})$ ), 'CV' for the coefficient of variation, or 'RMSLE' for the root mean-square log-error  |
| MSE       | logical; return the mean square error equivalent of the results instead of the root mean-square error (in other words, the result is squared)? Default is <code>FALSE</code>   |
| percent   | logical; change returned result to percentage by multiplying by 100? Default is <code>FALSE</code>   |
| unname    | logical; apply <code>unname</code> to the results to remove any variable names?  |

### Value

returns a numeric vector indicating the overall average deviation in the estimates

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

- Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248
- Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

### See Also

[bias](#)

MAE



**Examples**

```

pop <- 1
samp <- rnorm(100, 1, sd = 0.5)
RMSE(samp, pop)

dev <- samp - pop
RMSE(dev)

RMSE(samp, pop, type = 'NRMSE')
RMSE(dev, type = 'NRMSE')
RMSE(dev, pop, type = 'SRMSE')
RMSE(samp, pop, type = 'CV')
RMSE(samp, pop, type = 'RMSLE')

# percentage reported
RMSE(samp, pop, type = 'NRMSE')
RMSE(samp, pop, type = 'NRMSE', percent = TRUE)

# matrix input
mat <- cbind(M1=rnorm(100, 2, sd = 0.5), M2 = rnorm(100, 2, sd = 1))
RMSE(mat, parameter = 2)
RMSE(mat, parameter = c(2, 3))

# different parameter associated with each column
mat <- cbind(M1=rnorm(1000, 2, sd = 0.25), M2 = rnorm(1000, 3, sd = .25))
RMSE(mat, parameter = c(2,3))

# same, but with data.frame
df <- data.frame(M1=rnorm(100, 2, sd = 0.5), M2 = rnorm(100, 2, sd = 1))
RMSE(df, parameter = c(2,2))

# parameters of the same size
parameters <- 1:10
estimates <- parameters + rnorm(10)
RMSE(estimates, parameters)

```

---

rmvnorm

*Generate data with the multivariate normal (i.e., Gaussian) distribution*


---

**Description**

Function generates data from the multivariate normal distribution given some mean vector and/or covariance matrix.

**Usage**

```
rmvnorm(n, mean = rep(0, nrow(sigma)), sigma = diag(length(mean)))
```

**Arguments**

n                    number of observations to generate  
mean                mean vector, default is `rep(0, length = ncol(sigma))`  
sigma               positive definite covariance matrix, default is `diag(length(mean))`

**Value**

a numeric matrix with columns equal to `length(mean)`

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. [doi:10.1080/10691898.2016.1246953](https://doi.org/10.1080/10691898.2016.1246953)

**See Also**

[runSimulation](#)

**Examples**

```
# random normal values with mean [5, 10] and variances [3,6], and covariance 2
sigma <- matrix(c(3,2,2,6), 2, 2)
mu <- c(5,10)
x <- rmvnorm(1000, mean = mu, sigma = sigma)
head(x)
summary(x)
plot(x[,1], x[,2])
```

---

rmvt

*Generate data with the multivariate t distribution*

---

**Description**

Function generates data from the multivariate t distribution given a covariance matrix, non-centrality parameter (or mode), and degrees of freedom.

**Usage**

```
rmvt(n, sigma, df, delta = rep(0, nrow(sigma)), Kshirsagar = FALSE)
```

**Arguments**

|            |  |
|------------|--|
| n          | number of observations to generate   |
| sigma      | positive definite covariance matrix  |
| df         | degrees of freedom. $df = 0$ and $df = \text{Inf}$ corresponds to the multivariate normal distribution   |
| delta      | the vector of non-centrality parameters of length n which specifies the either the modes (default) or non-centrality parameters                                      |
| Kshirsagar | logical; triggers whether to generate data with non-centrality parameters or to adjust the simulated data to the mode of the distribution. The default uses the mode |

**Value**

a numeric matrix with columns equal to `ncol(sigma)`

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

**See Also**

[runSimulation](#)

**Examples**

```
# random t values given variances [3,6], covariance 2, and df = 15
sigma <- matrix(c(3,2,2,6), 2, 2)
x <- rmvt(1000, sigma = sigma, df = 15)
head(x)
summary(x)
plot(x[,1], x[,2])
```

RobbinsMonro

*Robbins-Monro (1951) stochastic root-finding algorithm***Description**

Function performs stochastic root solving for the provided  $f(x)$  using the Robbins-Monro (1951) algorithm. Differs from deterministic cousins such as [uniroot](#) in that  $f$  may contain stochastic error components, where the root is obtained through the running average method provided by noise filter (see also [PBA](#)). Assumes that  $E[f(x)]$  is non-decreasing in  $x$ .

**Usage**

```
RobbinsMonro(
  f,
  p,
  ...,
  Polyak_Juditsky = FALSE,
  maxiter = 500L,
  miniter = 100L,
  k = 3L,
  tol = 1e-05,
  verbose = TRUE,
  fn.a = function(iter, a = 1, b = 1/2, c = 0, ...) a/(iter + c)^b
)

## S3 method for class 'RM'
print(x, ...)
```

```
## S3 method for class 'RM'
plot(x, par = 1, main = NULL, Polyak_Juditsky = FALSE, ...)
```

**Arguments**

|                              |   |
|------------------------------|---|
| <code>f</code>               | noisy function for which the root is sought   |
| <code>p</code>               | vector of starting values to be passed as $f(p, \dots)$   |
| <code>...</code>             | additional named arguments to be passed to $f$  |
| <code>Polyak_Juditsky</code> | logical; apply the Polyak and Juditsky (1992) running-average method? Returns the final running average estimate using the Robbins-Monro updates (also applies to <code>plot</code> ). Note that this should only be used when the step-sizes are sufficiently large so that the Robbins-Monro have the ability to stochastically explore around the root (not just approach it from one side, which occurs when using small steps) |
| <code>maxiter</code>         | the maximum number of iterations (default 500)  |
| <code>miniter</code>         | minimum number of iterations (default 100)  |

|         |  |
|---------|--|
| k       | number of consecutive tol criteria required before terminating   |
| tol     | tolerance criteria for convergence on the changes in the updated p elements. Must be achieved on k (default 3) successive occasions  |
| verbose | logical; should the iterations and estimate be printed to the console?   |
| fn.a    | function to create the a coefficient in the Robbins-Monro noise filter. Requires the first argument is the current iteration (i ter), provide one or more arguments, and (optionally) the . . . . Sequence function is of the form recommended by Spall (2000).<br>Note that if a different function is provided it must satisfy the property that $\sum_{i=1}^{\infty} a_i = \infty$ and $\sum_{i=1}^{\infty} a_i^2 < \infty$ |
| x       | an object of class RM  |
| par     | which parameter in the original vector p to include in the plot  |
| main    | plot title   |

## References

- Polyak, B. T. and Juditsky, A. B. (1992). Acceleration of Stochastic Approximation by Averaging. *SIAM Journal on Control and Optimization*, 30(4):838.
- Robbins, H. and Monro, S. (1951). A stochastic approximation method. *Ann.Math.Statistics*, 22:400-407.
- Spall, J.C. (2000). Adaptive stochastic approximation by the simultaneous perturbation method. *IEEE Trans. Autom. Control* 45, 1839-1853.

## See Also

[uniroot](#), [PBA](#)

## Examples

```
# find x that solves f(x) - b = 0 for the following
f.root <- function(x, b = .6) 1 / (1 + exp(-x)) - b
f.root(.3)

xs <- seq(-3,3, length.out=1000)
plot(xs, f.root(xs), type = 'l', ylab = "f(x)", xlab='x')
abline(h=0, col='red')

retuni <- uniroot(f.root, c(0,1))
retuni
abline(v=retuni$root, col='blue', lty=2)

# Robbins-Monro without noisy root, start with p=.9
retrm <- RobbinsMonro(f.root, .9)
retrm
plot(retrm)

# Same problem, however root function is now noisy. Hence, need to solve
# fhat(x) - b + e = 0, where E(e) = 0
```

```
f.root_noisy <- function(x) 1 / (1 + exp(-x)) - .6 + rnorm(1, sd=.02)
sapply(rep(.3, 10), f.root_noisy)

# uniroot "converges" unreliably
set.seed(123)
uniroot(f.root_noisy, c(0,1))$root
uniroot(f.root_noisy, c(0,1))$root
uniroot(f.root_noisy, c(0,1))$root

# Robbins-Monro provides better convergence
retrm.noise <- RobbinsMonro(f.root_noisy, .9)
retrm.noise
plot(retrm.noise)

# different power (b) for fn.a()
retrm.b2 <- RobbinsMonro(f.root_noisy, .9, b = .01)
retrm.b2
plot(retrm.b2)

# use Polyak-Juditsky averaging (b should be closer to 0 to work well)
retrm.PJ <- RobbinsMonro(f.root_noisy, .9, b = .01,
                        Polyak_Juditsky = TRUE)
retrm.PJ # final Polyak_Juditsky estimate
plot(retrm.PJ) # Robbins-Monro history
plot(retrm.PJ, Polyak_Juditsky = TRUE) # Polyak_Juditsky history
```

---

RSE

---

*Compute the relative standard error ratio*


---

## Description

Computes the relative standard error ratio given the set of estimated standard errors (SE) and the deviation across the R simulation replications (SD). The ratio is formed by finding the expectation of the SE terms, and compares this expectation to the general variability of their respective parameter estimates across the R replications (ratio should equal 1). This is used to roughly evaluate whether the SEs being advertised by a given estimation method matches the sampling variability of the respective estimates across samples.

## Usage

```
RSE(SE, ests, unname = FALSE)
```

## Arguments

SE a numeric matrix of SE estimates across the replications (extracted from the results object in the Summarise step). Alternatively, can be a vector containing the mean of the SE estimates across the R simulation replications

`ests` a numeric matrix object containing the parameter estimates under investigation found within the [Summarise](#) function. This input is used to compute the standard deviation/variance estimates for each column to evaluate how well the expected SE matches the standard deviation

`unname` logical; apply [unname](#) to the results to remove any variable names?

**Value**

returns vector of variance ratios, ( $RSV = SE^2/SD^2$ )

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. [doi:10.1080/10691898.2016.1246953](https://doi.org/10.1080/10691898.2016.1246953)

**Examples**

```
R <- 10000
par_ests <- cbind(rnorm(R), rnorm(R, sd=1/10),
                 rnorm(R, sd=1/15))
colnames(par_ests) <- paste0("par", 1:3)
(SDs <- colSDs(par_ests))

SEs <- cbind(1 + rnorm(R, sd=.01),
             1/10 + rnorm(R, sd=.01),
             1/15 + rnorm(R, sd=.01))
(E_SEs <- colMeans(SEs))
RSE(SEs, par_ests)

# equivalent to the form
colMeans(SEs) / SDs
```

---

rtruncate

*Generate a random set of values within a truncated range*


---

**Description**

Function generates data given a supplied random number generating function that are constructed to fall within a particular range. Sampled values outside this range are discarded and re-sampled until the desired criteria has been met.

**Usage**

```
rtruncate(n, rfun, range, ..., redraws = 100L)
```

**Arguments**

|         |  |
|---------|--|
| n       | number of observations to generate. This should be the first argument passed to rfun   |
| rfun    | a function to generate random values. Function can return a numeric/integer vector or matrix, and additional arguments required for this function are passed through the argument ...  |
| range   | a numeric vector of length two, where the first element indicates the lower bound and the second the upper bound. When values are generated outside these two bounds then data are redrawn until the bounded criteria is met. When the output of rfun is a matrix then this input can be specified as a matrix with two rows, where each the first row corresponds to the lower bound and the second row the upper bound for each generated column in the output |
| ...     | additional arguments to be passed to rfun  |
| redraws | the maximum number of redraws to take before terminating the iterative sequence. This is in place as a safety in case the range is too small given the random number generator, causing too many consecutive rejections. Default is 100  |

**Details**

In simulations it is often useful to draw numbers from truncated distributions rather than across the full theoretical range. For instance, sampling parameters within the range [-4,4] from a normal distribution. The `rtruncate` function has been designed to accept any sampling function, where the first argument is the number of values to sample, and will draw values iteratively until the number of values within the specified bound are obtained. In situations where it is unlikely for the bounds to be located (e.g., sampling from a standard normal distribution where all values are within [-10,-6]) then the sampling scheme will throw an error if too many re-sampling executions are required (default will stop if more that 100 calls to `rfun` are required).

**Value**

either a numeric vector or matrix, where all values are within the desired range

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

- Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)
- Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. [doi:10.1080/10691898.2016.1246953](https://doi.org/10.1080/10691898.2016.1246953)



**See Also**[runSimulation](#)**Examples**

```
# n = 1000 truncated normal vector between [-2,3]
vec <- rtruncate(1000, rnorm, c(-2,3))
summary(vec)

# truncated correlated multivariate normal between [-1,4]
mat <- rtruncate(1000, rmvnorm, c(-1,4),
  sigma = matrix(c(2,1,1,1),2))
summary(mat)

# truncated correlated multivariate normal between [-1,4] for the
# first column and [0,3] for the second column
mat <- rtruncate(1000, rmvnorm, cbind(c(-1,4), c(0,3)),
  sigma = matrix(c(2,1,1,1),2))
summary(mat)

# truncated chi-square with df = 4 between [2,6]
vec <- rtruncate(1000, rchisq, c(2,6), df = 4)
summary(vec)
```

---

|                    |   |
|--------------------|---|
| runArraySimulation | <i>Run a Monte Carlo simulation using array job submissions per condition</i> |
|--------------------|---|

---

**Description**

This function has the same purpose as [runSimulation](#), however rather than evaluating each row in a design object (potentially with parallel computing architecture) this function evaluates the simulation per independent row condition. This is mainly useful when distributing the jobs to HPC clusters where a job array number is available (e.g., via SLURM), where the simulation results must be saved to independent files as they complete. Use of [expandDesign](#) is useful for distributing replications to different jobs, while [genSeeds](#) is required to ensure high-quality random number generation across the array submissions. See the associated vignette for a brief tutorial of this setup.

**Usage**

```
runArraySimulation(
  design,
  ...,
  replications,
  iseed,
  filename,
```

```

dirname = NULL,
arrayID = getArrayID(),
array2row = function(arrayID) arrayID,
addArrayInfo = TRUE,
parallel = FALSE,
cl = NULL,
ncores = parallelly::availableCores(omit = 1L),
save_details = list(),
control = list()
)

```

### Arguments

|              |   |
|--------------|---|
| design       | design object containing simulation conditions on a per row basis. This function is design to submit each row as in independent job on a HPC cluster. See <a href="#">runSimulation</a> for further details   |
| ...          | additional arguments to be passed to <a href="#">runSimulation</a>  |
| replications | number of independent replications to perform per condition (i.e., each row in design). See <a href="#">runSimulation</a> for further details   |
| iseed        | initial seed to be passed to <a href="#">genSeeds</a> 's argument of the same name, along with the supplied arrayID   |
| filename     | file name to save simulation files to (does not need to specify extension). However, the array ID will be appended to each filename. For example, if filename = 'mysim' then files stored will be 'mysim-1.rds', 'mysim-2.rds', and so on for each row ID in design   |
| dirname      | directory to save the files associated with filename to. If omitted the files will be stored in the same working directory where the script was submitted   |
| arrayID      | array identifier from the scheduler. Must be a number between 1 and nrow(design). If not specified then <a href="#">getArrayID</a> will be called automatically, which assumes the environmental variables are available according the SLURM scheduler  |
| array2row    | user defined function with the single argument arrayID. Used to convert the detected arrayID into a suitable row index in the design object input. By default each arrayID is associated with its respective row in design.<br>For example, if each arrayID should evaluate 10 rows in the design object then the function <code>function(arrayID){1:10 + 10 * (arrayID-1)}</code> can be passed to array2row |
| addArrayInfo | logical; should the array ID and original design row number be added to the <code>SimResults(...)</code> output?  |
| parallel     | logical; use parallel computations via the a "SOCK" cluster? Only useful when the instruction shell file requires more than 1 core (number of cores detected via ncores). For this application the random seeds further distributed using <a href="#">nextRNGSubStream</a>  |
| cl           | cluster definition. If omitted a "SOCK" cluster will be defined   |
| ncores       | number of cores to use when parallel=TRUE. Note that the default uses 1 minus the number of available cores, therefore this will only be useful when ncores > 2 as defined in the shell instruction file  |

|              |  |
|--------------|--|
| save_details | optional list of extra file saving details. See <a href="#">runSimulation</a>  |
| control      | control list passed to <a href="#">runSimulation</a> . In addition to the original control elements two additional arguments have been added: <code>max_time</code> and <code>max_RAM</code> , both of which as specified as character vectors with one element.<br><br><code>max_time</code> specifies the maximum time allowed for a single simulation condition to execute (default does not set any time limits), and is formatted according to the specification in <a href="#">timeFormater</a> . This is primarily useful when the HPC cluster will time out after some known elapsed time. In general, this input should be set to somewhere around 80-90 before the cluster is terminated can be saved. Default applies no time limit<br><br>Similarly, <code>max_RAM</code> controls the (approximate) maximum size that the simulation storage objects can grow before RAM becomes an issue. This can be specified either in terms of megabytes (MB), gigabytes (GB), or terabytes (TB). For example, <code>max_RAM = "4GB"</code> indicates that if the simulation storage objects are larger than 4GB then the workflow will terminate early, returning only the successful results up to this point). Useful for larger HPC cluster jobs with RAM constraints that could terminate abruptly. As a rule of thumb this should be set to around 90 available. Default applies no memory limit |

## Details

Due to the nature of how the replication are split it is important that the L'Ecuyer-CMRG (2002) method of random seeds is used across all array ID submissions (cf. [runSimulation](#)'s parallel approach, which uses this method to distribute random seeds within each isolated condition rather than between all conditions). As such, this function requires the seeds to be generated using [genSeeds](#) with the `iseed` and `arrayID` inputs to ensure that each job is analyzing a high-quality set of random numbers via L'Ecuyer-CMRG's (2002) method, incremented using [nextRNGStream](#).

Additionally, for timed simulations on HPC clusters it is also recommended to pass a `control = list(max_time)` value to avoid discarding conditions that require more than the specified time in the shell script. The `max_time` value should be less than the maximum time allocated on the HPC cluster (e.g., approximately 90 depends on how long each replication takes). Simulations with missing replication information should submit a new set of jobs at a later time to collect the missing replication information.

## Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

## References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)

## See Also

[runSimulation](#), [expandDesign](#), [genSeeds](#), [SimCheck](#), [SimCollect](#), [getArrayID](#)

**Examples**

```

library(SimDesign)

Design <- createDesign(N = c(10, 20, 30))

Generate <- function(condition, fixed_objects) {
  dat <- with(condition, rnorm(N, 10, 5)) # distributed N(10, 5)
  dat
}

Analyse <- function(condition, dat, fixed_objects) {
  ret <- c(mean=mean(dat), median=median(dat)) # mean/median of sample data
  ret
}

Summarise <- function(condition, results, fixed_objects){
  colMeans(results)
}

## Not run:

# define initial seed (do this only once to keep it constant!)
# iseed <- genSeeds()
iseed <- 554184288

### On cluster submission, the active array ID is obtained via getArrayID(),
### and therefore should be used in real SLURM submissions
arrayID <- getArrayID(type = 'slurm')

# However, the following example arrayID is set to
# the first row only for testing purposes
arrayID <- 1L

# run the simulation (results not caught on job submission, only files saved)
res <- runArraySimulation(design=Design, replications=50,
  generate=Generate, analyse=Analyse,
  summarise=Summarise, arrayID=arrayID,
  iseed=iseed, filename='mysim') # saved as 'mysim-1.rds'

res
SimResults(res) # condition and replication count stored

# same, but evaluated with multiple cores
res <- runArraySimulation(design=Design, replications=50,
  generate=Generate, analyse=Analyse,
  summarise=Summarise, arrayID=arrayID,
  parallel=TRUE, ncores=3,
  iseed=iseed, filename='myparsim')

res
SimResults(res) # condition and replication count stored

dir()
SimClean(c('mysim-1.rds', 'myparsim-1.rds'))

```

```
#####
# Same submission job as above, however split the replications over multiple
# evaluations and combine when complete
Design5 <- expandDesign(Design, 5)
Design5

# iseed <- genSeeds()
iseed <- 554184288

# arrayID <- getArrayID(type = 'slurm')
arrayID <- 14L

# run the simulation (replications reduced per row, but same in total)
runArraySimulation(design=Design5, replications=10,
                  generate=Generate, analyse=Analyse,
                  summarise=Summarise, iseed=iseed,
                  filename='mylongsim', arrayID=arrayID)

res <- readRDS('mylongsim-14.rds')
res
SimResults(res) # condition and replication count stored

SimClean('mylongsim-14.rds')

###
# Emulate the arrayID distribution, storing all results in a 'sim/' folder
# (if 'sim/' does not exist in runArraySimulation() it will be
# created automatically)
dir.create('sim/')

# Emulate distribution to nrow(Design5) = 15 independent job arrays
## (just used for presentation purposes on local computer)
sapply(1:nrow(Design5), \(arrayID)
      runArraySimulation(design=Design5, replications=10,
                        generate=Generate, analyse=Analyse,
                        summarise=Summarise, iseed=iseed, arrayID=arrayID,
                        filename='condition', dirname='sim', # files: "sim/condition-#.rds"
                        control = list(max_time="04:00:00", max_RAM="4GB"))) |> invisible()

# If necessary, conditions above will manually terminate before
# 4 hours and 4GB of RAM are used, returning any
# successfully completed results before the HPC session times
# out (provided .slurm script specified more than 4 hours)

# list saved files
dir('sim/')

# check that all files saved (warnings will be raised if missing files)
SimCheck('sim/') |> isTRUE()

condition14 <- readRDS('sim/condition-14.rds')
```

```

condition14
SimResults(condition14)

# aggregate simulation results into single file
final <- SimCollect('sim/')
final

# clean simulation directory
SimClean(dirs='sim/')

#####
# same as above, however passing different amounts of information depending
# on the array ID
array2row <- function(arrayID){
  switch(arrayID,
    "1"=1:8,
    "2"=9:14,
    "3"=15)
}

# arrayID 1 does row 1 though 8, arrayID 2 does 9 to 14
array2row(1)
array2row(2)
array2row(3) # arrayID 3 does 15 only

# emulate remote array distribution with only 3 arrays
sapply(1:3, \(arrayID)
  runArraySimulation(design=Design5, replications=10,
    generate=Generate, analyse=Analyse,
    summarise=Summarise, iseed=iseed, arrayID=arrayID,
    filename='condition', dirname='sim', array2row=array2row)) |> invisible()

# list saved files
dir('sim/')

# note that all row conditions are still stored separately, though note that
# arrayID is now 2 instead
condition14 <- readRDS('sim/condition-14.rds')
condition14
SimResults(condition14)

# aggregate simulation results into single file
final <- SimCollect('sim/')
final

# clean simulation directory
SimClean(dirs='sim/')

## End(Not run)

```

---

|               |   |
|---------------|---|
| runSimulation | <i>Run a Monte Carlo simulation given conditions and simulation functions</i> |
|---------------|---|

---

### Description

This function runs a Monte Carlo simulation study given a set of predefined simulation functions, design conditions, and number of replications. Results can be saved as temporary files in case of interruptions and may be restored by re-running `runSimulation`, provided that the respective temp file can be found in the working directory. `runSimulation` supports parallel and cluster computing (with the `parallel` and `future` packages; see also `runArraySimulation` for submitting array jobs to HPC clusters), global and local debugging, error handling (including fail-safe stopping when functions fail too often, even across nodes), provides bootstrap estimates of the sampling variability (optional), and automatic tracking of error and warning messages with their associated `.Random.seed` states. For convenience, all functions available in the R work-space are exported across all nodes so that they are more easily accessible (however, other R objects are not, and therefore must be passed to the `fixed_objects` input to become available across nodes).

### Usage

```
runSimulation(  
  design,  
  replications,  
  generate,  
  analyse,  
  summarise,  
  fixed_objects = NULL,  
  packages = NULL,  
  filename = NULL,  
  debug = "none",  
  load_seed = NULL,  
  save = any(replications > 2),  
  store_results = TRUE,  
  save_results = FALSE,  
  parallel = FALSE,  
  ncores = parallelly::availableCores(omit = 1L),  
  cl = NULL,  
  notification = "none",  
  beep = FALSE,  
  sound = 1,  
  CI = 0.95,  
  seed = NULL,  
  boot_method = "none",  
  boot_draws = 1000L,  
  max_errors = 50L,  
  resume = TRUE,  
  save_details = list(),
```

```

    control = list(),
    progress = TRUE,
    verbose = TRUE
  )

  ## S3 method for class 'SimDesign'
  summary(object, ...)

  ## S3 method for class 'SimDesign'
  print(x, list2char = TRUE, ...)

```

## Arguments

|              |   |
|--------------|---|
| design       | a tibble or data.frame object containing the Monte Carlo simulation conditions to be studied, where each row represents a unique condition and each column a factor to be varied. See <a href="#">createDesign</a> for the standard approach to create this simulation design object  |
| replications | number of independent replications to perform per condition (i.e., each row in design). Can be a single number, which will be used for each design condition, or an integer vector with length equal to nrow(design). All inputs must be greater than 0, though setting to less than 3 (for initial testing purpose) will disable the save and control\$stop_on_fatal flags   |
| generate     | user-defined data and parameter generating function (or named list of functions). See <a href="#">Generate</a> for details. Note that this argument may be omitted by the user if they wish to generate the data with the analyse step, but for real-world simulations this is generally not recommended. If multiple generate functions are provided as a list then the list of generate functions are executed in order until the first valid generate function is executed, where the subsequent generation functions are then ignored (see <a href="#">GenerateIf</a> to only apply data generation for specific conditions).   |
| analyse      | user-defined analysis function (or named list of functions) that acts on the data generated from <a href="#">Generate</a> (or, if generate was omitted, can be used to generate and analyses the simulated data). See <a href="#">Analyse</a> for details   |
| summarise    | optional (but strongly recommended) user-defined summary function from <a href="#">Summarise</a> to be used to compute meta-statistical summary information after all the replications have completed within each design condition. Return of this function, in order of increasing complexity, should be: a named numeric vector or data.frame with one row, a matrix or data.frame with more than one row, and, failing these more atomic types, a named list. For summary objects that are not easily appended to the original design object use <a href="#">SimExtract</a> with the option what = 'summarise'.<br>Note that unlike the Generate and Analyse steps, the Summarise portion is not as important to perfectly organize as the results can be summarised later on by using the built-in <a href="#">reSummarise</a> function (provided either store_results = TRUE or save_results = TRUE were included).<br>Omitting this function will return a tibble with the Design and associated results information for all nrow(Design) * replications evaluations if the results from each Analyse() call was a one-dimensional vector. For more general |



objects returned by `Analyse()` (such as lists), a list containing the results returned from `Analyse`. This is generally only recommended for didactic purposes because the results will leave out a large amount of information (e.g., try-errors, warning messages, saving files, etc), can witness memory related issues if the `Analyse` function returns larger objects, and generally is not as flexible internally. However, it may be useful when replications are expensive and ANOVA-based decompositions involving the within-condition replication information are of interest, though of course this can be circumvented by using `store_results = TRUE` or `save_results = TRUE` with or without a supplied summarise definition.

- `fixed_objects` (optional) an object (usually a named list) containing additional user-defined objects that should remain fixed across conditions. This is useful when including large vectors/matrices of population parameters, fixed data information that should be used across all conditions and replications (e.g., including a common design matrix for linear regression models), or simply control constant global elements (e.g., a constant for sample size)
- `packages` a character vector of external packages to be used during the simulation (e.g., `c('MASS', 'extraDistr', 'simsem')`). Use this input when running code in parallel to use non-standard functions from additional packages, otherwise the functions must be made available by using explicit `library` or `require` calls within the provided simulation functions. Alternatively, functions can be called explicitly without attaching the package with the `::` operator (e.g., `extraDistr::rgumbel()`)
- `filename` (optional) the name of the `.rds` file to save the final simulation results to. If the extension `.rds` is not included in the file name (e.g. "mysimulation" versus "mysimulation.rds") then the `.rds` extension will be automatically added to the file name to ensure the file extension is correct.
- Note that if the same file name already exists in the working directory at the time of saving then a new file will be generated instead and a warning will be thrown. This helps to avoid accidentally overwriting existing files. Default is `NULL`, indicating no file will be saved by default
- `debug` a string indicating where to initiate a `browser()` call for editing and debugging, and pairs particularly well with the `load_seed` argument for precise debugging. General options are 'none' (default; no debugging), 'error', which starts the debugger when any error in the code is detected in one of three generate-analyse-summarise functions, and 'all', which debugs all the user defined functions regardless of whether an error was thrown or not. Specific options include: 'generate' to debug the data simulation function, 'analyse' to debug the computational function, and 'summarise' to debug the aggregation function.
- If the `Analyse` argument is supplied as a named list of functions then it is also possible to debug the specific function of interest by passing the name of the respective function in the list. For instance, if `analyse = list(A1=Analyse.A1, A2=Analyse.A2)` then passing `debug = 'A1'` will debug only the first function in this list, and all remaining analysis functions will be ignored.
- For debugging specific rows in the `Design` input (e.g., when a number of initial rows successfully complete but the *k*th row fails) the row number can be appended to the standard debug input using a '-' separator. For instance, de-

bugging whenever an error is raised in the second row of Design can be declared via `debug = 'error-2'`.

Finally, users may place `browser` calls within the respective functions for debugging at specific lines, which is useful when debugging based on conditional evaluations (e.g., `if(this == 'problem') browser()`). Note that parallel computation flags will automatically be disabled when a `browser()` is detected or when a debugging argument other than 'none' is supplied

|               |   |
|---------------|---|
| load_seed     | used to replicate an exact simulation state, which is primarily useful for debugging purposes. Input can be a character object indicating which file to load from when the <code>.Random.seeds</code> have been saved (after a call with <code>save_seeds = TRUE</code> ), or an integer vector indicating the actual <code>.Random.seed</code> values (e.g., extracted after using <code>store_seeds</code> ). E.g., <code>load_seed = 'design-row-2/seed-1'</code> will load the first seed in the second row of the design input, or explicitly passing the elements from <code>.Random.seed</code> (see <code>SimExtract</code> to extract the seeds associated explicitly with errors during the simulation, where each column represents a unique seed). If the input is a character vector then it is important NOT to modify the design input object, otherwise the path may not point to the correct saved location, while if the input is an integer vector (or single column tbl object) then it WILL be important to modify the design input in order to load this exact seed for the corresponding design row. Default is NULL                                     |
| save          | logical; save the temporary simulation state to the hard-drive? This is useful for simulations which require an extended amount of time, though for shorter simulations can be disabled to slightly improve computational efficiency. When TRUE, which is the default when evaluating <code>replications &gt; 2</code> , a temp file will be created in the working directory which allows the simulation state to be saved and recovered (in case of power outages, crashes, etc). As well, triggering this flag will save any fatal <code>.Random.seed</code> states when conditions unexpectedly crash (where each seed is stored row-wise in an external <code>.rds</code> file), which provides a much easier mechanism to debug issues (see <code>load_seed</code> for details). Upon completion, this temp file will be removed.<br><br>To recover your simulation at the last known location (having patched the issues in the previous execution code) simply re-run the code you used to initially define the simulation and the external file will automatically be detected and read-in. Default is TRUE when <code>replications &gt; 10</code> and FALSE otherwise |
| store_results | logical; store the complete tables of simulation results in the returned object? This is TRUE default, though if RAM anticipated to be an issue see <code>save_results</code> instead. Note that if the Design object is omitted from the call to <code>runSimulation()</code> , or the number of rows in Design is exactly 1, then this argument is automatically set to TRUE as RAM storage is no longer an issue.<br><br>To extract these results pass the returned object to either <code>SimResults</code> or <code>SimExtract</code> with <code>what = 'results'</code> , which will return a named list of all the simulation results for each condition if <code>nrow(Design) &gt; 1</code> ; otherwise, if <code>nrow(Design) == 1</code> or Design was missing the results object will be stored as-is  |
| save_results  | logical; save the results returned from <code>Analyse</code> to external <code>.rds</code> files located in the defined <code>save_results_dirname</code> directory/folder? Use this if you would like to keep track of the individual parameters returned from the analysis function. Each saved object will contain a list of three elements containing the con-  |

dition (row from design), results (as a list or matrix), and try-errors. See [SimResults](#) for an example of how to read these .rds files back into R after the simulation is complete. Default is FALSE.

WARNING: saving results to your hard-drive can fill up space very quickly for larger simulations. Be sure to test this option using a smaller number of replications before the full Monte Carlo simulation is performed. See also [reSummarise](#) for applying summarise functions from saved simulation results

|              |   |
|--------------|---|
| parallel     | <p>logical; use parallel processing from the <a href="#">parallel</a> package over each unique condition? This distributes the independent replications across the defined nodes, and is repeated for each row condition in the design input.</p> <p>Alternatively, if the <a href="#">future</a> package approach is desired then passing <code>parallel = 'future'</code> to <code>runSimulation()</code> will use the defined <a href="#">plan</a> for execution. This allows for greater flexibility when specifying the general computing plan (e.g., <code>plan(multisession)</code>) for parallel computing on the same machine, <code>plan(future.batchtools::batchtools_slurm)</code> for common MPI/Slurm schedulers, etc). However, it is the responsibility of the user to use <code>plan(sequential)</code> to reset the computing plan when the jobs are completed</p>  |
| ncores       | number of cores to be used in parallel execution (ignored if using the <a href="#">future</a> package approach). Default uses all available minus 1   |
| c1           | <p>cluster object defined by <a href="#">makeCluster</a> used to run code in parallel (ignored if using the <a href="#">future</a> package approach). If NULL and <code>parallel = TRUE</code>, a local cluster object will be defined which selects the maximum number cores available and will be stopped when the simulation is complete. Note that supplying a <code>c1</code> object will automatically set the <code>parallel</code> argument to TRUE. Define and supply this cluster object yourself whenever you have multiple nodes and can link them together manually</p> <p>If the <a href="#">future</a> package has been attached prior to executing <code>runSimulation()</code> then the associated <code>plan()</code> will be followed instead</p>  |
| notification | <p>an optional character vector input that can be used to send Pushbullet notifications from a configured computer. This reports information such as the total execution time, the condition completed, and error/warning messages recorded. This arguments assumes that users have already A) registered for a Pushbullet account, B) installed the application on their mobile device and computer, and C) created an associated JSON file of the form <code>~/rpushbullet.json</code> using <code>RPushbullet::pbSetup()</code>.</p> <p>To utilize the <code>RPushbullet</code> in <code>SimDesign</code> first call <code>library(RPushbullet)</code> before running <code>runSimulation()</code> to read-in the default JSON file. Next, pass one of the following supported options: 'none' (default; send no notification), 'condition' to send a notification after each condition has completed, or 'complete' to send a notification only when the simulation has finished.</p> |
| beep         | logical; call the <code>beepR</code> package when the simulation is completed?  |
| sound        | sound argument passed to <code>beepR::beep()</code>   |
| CI           | bootstrap confidence interval level (default is 95%)  |
| seed         | a vector or list of integers to be used for reproducibility. The length of the vector must be equal the number of rows in design. If the input is a vector  |

then `set.seed` or `clusterSetRNGStream` for each condition will be called, respectively. If a list is provided then these numbers must have been generated from `gen_seeds` with the argument `CMRG.seed` used to specify the initial. The list approach ensures random number generation independence across conditions and replications, while the vector input ensures independence within the replications per conditions but not necessarily across conditions. Default randomly generates seeds within the range 1 to 2147483647 for each condition via `link{gen_seeds}`

|                           |  |
|---------------------------|--|
| <code>boot_method</code>  | method for performing non-parametric bootstrap confidence intervals for the respective meta-statistics computed by the <code>Summarise</code> function. Can be 'basic' for the empirical bootstrap CI, 'percentile' for percentile CIs, 'norm' for normal approximations CIs, or 'studentized' for Studentized CIs (should only be used for simulations with lower replications due to its computational intensity). Alternatively, CIs can be constructed using the argument 'CLT', which computes the intervals according to the large-sample standard error approximation $SD(results)/\sqrt{R}$ . Default is 'none', which performs no CI computations   |
| <code>boot_draws</code>   | number of non-parametric bootstrap draws to sample for the <code>summarise</code> function after the generate-analyse replications are collected. Default is 1000  |
| <code>max_errors</code>   | the simulation will terminate when more than this number of consecutive errors are thrown in any given condition, causing the simulation to continue to the next unique design condition. This is included to avoid getting stuck in infinite re-draws, and to indicate that something fatally problematic is going wrong in the generate-analyse phases. Default is 50  |
| <code>resume</code>       | logical; if a temporary <code>SimDesign</code> file is detected should the simulation resume from this location? Keeping this TRUE is generally recommended, however this should be disabled if using <code>runSimulation</code> within <code>runSimulation</code> to avoid reading improper save states. Alternatively, if an integer is supplied then the simulation will continue at the associated row location in <code>design</code> (e.g., <code>resume=10</code> ). This is useful to overwrite a previously evaluate element in the temporary files that was detected to contain fatal errors that require re-evaluation without discarding the originally valid rows in the simulation   |
| <code>save_details</code> | <p>a list pertaining to information regarding how and where files should be saved when the <code>save</code> or <code>save_results</code> flags are triggered.</p> <p><code>safe</code> logical; trigger whether safe-saving should be performed. When TRUE files will never be overwritten accidentally, and where appropriate the program will either stop or generate new files with unique names. Default is TRUE</p> <p><code>compname</code> name of the computer running the simulation. Normally this doesn't need to be modified, but in the event that a manual node breaks down while running a simulation the results from the temp files may be resumed on another computer by changing the name of the node to match the broken computer. Default is the result of evaluating <code>uname(Sys.info()['nodename'])</code></p> <p><code>out_rootdir</code> root directory to save all files to. Default uses the current working directory</p> <p><code>save_results_dirname</code> a string indicating the name of the folder to save result objects to when <code>save_results = TRUE</code>. If a directory/folder does not exist</p> |

in the current working directory then a unique one will be created automatically. Default is 'SimDesign-results\_' with the associated compname appended if no filename is defined, otherwise the filename is used to replace 'SimDesign' in the string

`save_results_filename` a string indicating the name file to store, where the Design row ID will be appended to ensure uniqueness across rows. Specifying this input will disable any checking for the uniqueness of the file folder, thereby allowing independent runSimulation calls to write to the same `save_results_dirname`. Useful when the files should all be stored in the same working directory, however the rows of Design are evaluated in isolation (e.g., for HPC structures that allow asynchronous file storage). **WARNING:** the uniqueness of the file names are not checked using this approach, therefore please ensure that each generated name will be unique a priori, such as naming the file based on the supplied row condition information

`save_seeds_dirname` a string indicating the name of the folder to save `.Random.seed` objects to when `save_seeds = TRUE`. If a directory/folder does not exist in the current working directory then one will be created automatically. Default is 'SimDesign-seeds\_' with the associated compname appended if no filename is defined, otherwise the filename is used to replace 'SimDesign' in the string

`tmpfilename` string indicating the temporary file name to save provisional information to. If not specified the following will be used: `paste0('SIMDESIGN-TEMPFILE_', compname, '.rds')`

control

a list for extra information flags for controlling less commonly used features. These include

`stop_on_fatal` logical (default is FALSE); should the simulation be terminated immediately when the maximum number of consecutive errors (`max_errors`) is reached? If FALSE, the simulation will continue as though errors did not occur, however a column `FATAL_TERMINATION` will be included in the resulting object indicating the final error message observed, and NA placeholders will be placed in all other row-elements. Default is FALSE, though is automatically set to TRUE when `replications < 3` for the purpose of debugging

`warnings_as_errors` logical (default is FALSE); treat warning messages as error messages during the simulation? Default is FALSE, therefore warnings are only collected and not used to restart the data generation step, and the seeds associated with the warning message conditions are not stored within the final simulation object.

Note that this argument is generally intended for debugging/early planning stages when designing a simulation experiment. If specific warnings are known to be problematic and should be treated as errors then please use [manageWarnings](#) instead

`save_seeds` logical; save the `.Random.seed` states prior to performing each replication into plain text files located in the defined `save_seeds_dirname` directory/folder? Use this if you would like to keep track of every simulation state within each replication and design condition. This can be used to

completely replicate any cell in the simulation if need be. As well, see the `load_seed` input to load a given `.Random.seed` to exactly replicate the generated data and analysis state (mostly useful for debugging). When `TRUE`, temporary files will also be saved to the working directory (in the same way as when `save = TRUE`). Default is `FALSE`

Note, however, that this option is not typically necessary or recommended since the `.Random.seed` states for simulation replications that throw errors during the execution are automatically stored within the final simulation object, and can be extracted and investigated using `SimExtract`. Hence, this option is only of interest when *all* of the replications must be reproducible (which occurs very rarely), otherwise the defaults to `runSimulation` should be sufficient

`store_Random_seeds` logical; store the complete `.Random.seed` states for each simulation replicate? Default is `FALSE` as this can take up a great deal of unnecessary RAM (see related `save_seeds`), however this may be useful when used with `runArraySimulation`. To extract use `SimExtract(..., what = 'stored_Random_seeds')`

`store_warning_seeds` logical (default is `FALSE`); in addition to storing the `.Random.seed` states whenever error messages are raised, also store the `.Random.seed` states when warnings are raised? This is disabled by default since warnings are generally less problematic than errors, and because many more warnings messages may be raised throughout the simulation (potentially causing RAM related issues when constructing the final simulation object as any given simulation replicate could generate numerous warnings, and storing the seeds states could add up quickly).

Set this to `TRUE` when replicating warning messages is important, however be aware that too many warnings messages raised during the simulation implementation could cause RAM related issues.

`include_replication_index` or `include_reps` logical (default is `FALSE`); should a `REPLICATION` element be added to the condition object when performing the simulation to track which specific replication experiment is being evaluated? This is useful when, for instance, attempting to run external software programs (e.g., `Mplus`) that require saving temporary data sets to the hard-drive (see the Wiki for examples)

`try_all_analyse` logical; when `analyse` is a list, should every generated data set be analyzed by each function definition in the `analyse` list? Default is `TRUE`.

Note that this `TRUE` default can be computationally demanding when some analysis functions require more computational resources than others, and the data should be discarded early as an invalid candidate (e.g., estimating a model via maximum-likelihood in on `analyse` component, while estimating a model using MCMC estimation on another). Hence, the main benefit of using `FALSE` instead is that the data set may be rejected earlier, where easier/faster to estimate `analyse` definitions should be placed earlier in the list as the functions are evaluated in sequence (e.g., `Analyse = list(MLE=MLE_definition, MCMC=MCMC_definition)`)

`allow_na` logical (default is `FALSE`); should NAs be allowed in the `analyse` step as a valid result from the simulation analysis?

|           |  |
|-----------|--|
|           | allow_nan logical (default is FALSE); should NaNs be allowed in the analyse step as a valid result from the simulation analysis?   |
|           | type default type of cluster to create for the cl object if not supplied. For Windows OS this defaults to "PSOCK", otherwise "SOCK" is selected (suitable for Linux and Mac OSX). This is ignored if the user specifies their own cl object  |
|           | print_RAM logical (default is TRUE); print the amount of RAM used throughout the simulation? Set to FALSE if unnecessary or if the call to gc is unnecessarily time consuming  |
|           | max_time Similar to runArraySimulation, specifies the (approximate) maximum time that the simulation is allowed to be executed. However, unlike the implementation in runArraySimulation is evaluated on a per condition basis, where max_time is only evaluated after every row in the design object has been completed (hence, is notably more approximate as it has the potential to overshoot by a wider margin). Default sets no time limit. See timeFormater for the input specifications. |
|           | max_RAM Similar to runArraySimulation, specifies the (approximate) maximum RAM that the simulation is allowed to occupy. However, unlike the implementation in runArraySimulation is evaluated on a per condition basis, where max_RAM is only evaluated after every row in the design object has been completed (hence, is notably more approximate as it has the potential to overshoot by a wider margin). Default sets no RAM limit. See runArraySimulation for the input specifications.    |
| progress  | logical; display a progress bar (using the pbapply package) for each simulation condition? This is useful when simulations conditions take a long time to run (see also the notifications argument). Default is TRUE   |
| verbose   | logical; print messages to the R console? Default is TRUE  |
| object    | SimDesign object returned from runSimulation   |
| ...       | additional arguments   |
| x         | SimDesign object returned from runSimulation   |
| list2char | logical; for tibble object re-evaluate list elements as character vectors for better printing of the levels? Note that this does not change the original classes of the object, just how they are printed. Default is TRUE   |

## Details

For an in-depth tutorial of the package please refer to Chalmers and Adkins (2020; doi:10.20982/tqmp.16.4.p248). For an earlier didactic presentation of the package refer to Sigal and Chalmers (2016; doi:10.1080/10691898.2016.1246953). Finally, see the associated wiki on Github (<https://github.com/philchalmers/SimDesign/wiki>) for tutorial material, examples, and applications of SimDesign to real-world simulation experiments, as well as the various vignette files associated with the package.

The strategy for organizing the Monte Carlo simulation work-flow is to

- 1) Define a suitable Design object (a tibble or data.frame) containing fixed conditional information about the Monte Carlo simulations. Each row of this design object pertains to a unique

set of simulation to study, while each column the simulation factor under investigation (e.g., sample size, distribution types, etc). This is often expedited by using the `createDesign` function, and if necessary the argument `subset` can be used to remove redundant or non-applicable rows

- 2) Define the three step functions to generate the data (`Generate`; see also <https://CRAN.R-project.org/view=Distributions> for a list of distributions in R), analyse the generated data by computing the respective parameter estimates, detection rates, etc (`Analyse`), and finally summarise the results across the total number of replications (`Summarise`).
- 3) Pass the design object and three defined R functions to `runSimulation`, and declare the number of replications to perform with the `replications` input. This function will return a suitable tibble object with the complete simulation results and execution details
- 4) Analyze the output from `runSimulation`, possibly using ANOVA techniques (`SimAnova`) and generating suitable plots and tables

Expressing the above more succinctly, the functions to be called have the following form, with the exact functional arguments listed:

```
Design <- createDesign(...)
Generate <- function(condition, fixed_objects) {...}
Analyse <- function(condition, dat, fixed_objects) {...}
Summarise <- function(condition, results, fixed_objects) {...}
res <- runSimulation(design=Design, replications, generate=Generate, analyse=Analyse, summarise=Summarise)
```

The `condition` object above represents a single row from the design object, indicating a unique Monte Carlo simulation condition. The `condition` object also contains two additional elements to help track the simulation's state: an ID variable, indicating the respective row number in the design object, and a `REPLICATION` element indicating the replication iteration number (an integer value between 1 and `replication`). This setup allows users to easily locate the  $r$ th replication (e.g., `REPLICATION == 500`) within the  $j$ th row in the simulation design (e.g., `ID == 2`). The `REPLICATION` input is also useful when temporarily saving files to the hard-drive when calling external command line utilities (see examples on the wiki).

For a template-based version of the work-flow, which is often useful when initially defining a simulation, use the `SimFunctions` function. This function will write a template simulation to one/two files so that modifying the required functions and objects can begin immediately. This means that users can focus on their Monte Carlo simulation details right away rather than worrying about the repetitive administrative code-work required to organize the simulation's execution flow.

Finally, examples, presentation files, and tutorials can be found on the package wiki located at <https://github.com/philchalmers/SimDesign/wiki>.

## Value

a tibble from the `dplyr` package (also of class `'SimDesign'`) with the original design conditions in the left-most columns, simulation results in the middle columns, and additional information in the right-most columns (see below).

The right-most column information for each condition are: `REPLICATIONS` to indicate the number of Monte Carlo replications, `SIM_TIME` to indicate how long (in seconds) it took to complete all



the Monte Carlo replications for each respective design condition, RAM\_USED amount of RAM that was in use at the time of completing each simulation condition, COMPLETED to indicate the date in which the given simulation condition completed, SEED for the integer values in the seed argument (if applicable; not relevant when L'Ecuyer-CMRG method used), and, if applicable, ERRORS and WARNINGS which contain counts for the number of error or warning messages that were caught (if no errors/warnings were observed these columns will be omitted). Note that to extract the specific error and warnings messages see [SimExtract](#). Finally, if boot\_method was a valid input other than 'none' then the final right-most columns will contain the labels BOOT\_ followed by the name of the associated meta-statistic defined in summarise() and bootstrapped confidence interval location for the meta-statistics.

### **Saving data, results, seeds, and the simulation state**

To conserve RAM, temporary objects (such as data generated across conditions and replications) are discarded; however, these can be saved to the hard-disk by passing the appropriate flags. For longer simulations it is recommended to use the save\_results flag to write the analysis results to the hard-drive.

The use of the store\_seeds or the save\_seeds options can be evoked to save R's .Random.seed state to allow for complete reproducibility of each replication within each condition. These individual .Random.seed terms can then be read in with the load\_seed input to reproduce the exact simulation state at any given replication. Most often though, saving the complete list of seeds is unnecessary as problematic seeds are automatically stored in the final simulation object to allow for easier replicability of potentially problematic errors (which incidentally can be extracted using SimExtract(res, 'error\_seeds') and passed to the load\_seed argument). Finally, providing a vector of seeds is also possible to ensure that each simulation condition is macro reproducible under the single/multi-core method selected.

Finally, when the Monte Carlo simulation is complete it is recommended to write the results to a hard-drive for safe keeping, particularly with the filename argument provided (for reasons that are more obvious in the parallel computation descriptions below). Using the filename argument supplied is safer than using, for instance, saveRDS directly because files will never accidentally be overwritten, and instead a new file name will be created when a conflict arises; this type of implementation safety is prevalent in many locations in the package to help avoid unrecoverable (yet surprisingly common) mistakes during the process of designing and executing Monte Carlo simulations.

### **Resuming temporary results**

In the event of a computer crash, power outage, etc, if save = TRUE was used (the default) then the original code used to execute runSimulation() need only be re-run to resume the simulation. The saved temp file will be read into the function automatically, and the simulation will continue one the condition where it left off before the simulation state was terminated. If users wish to remove this temporary simulation state entirely so as to start anew then simply pass SimClean(temp = TRUE) in the R console to remove any previously saved temporary objects.

### **A note on parallel computing**

When running simulations in parallel (either with parallel = TRUE or when using the [future](#) approach with a plan() other than sequential) R objects defined in the global environment will generally *not* be visible across nodes. Hence, you may see errors such as Error: object 'something'

not found if you try to use an object that is defined in the work space but is not passed to `runSimulation`. To avoid this type of error, simply pass additional objects to the `fixed_objects` input (usually it's convenient to supply a named list of these objects). Fortunately, however, *custom functions defined in the global environment are exported across nodes automatically*. This makes it convenient when writing code because custom functions will always be available across nodes if they are visible in the R work space. As well, note the `packages` input to declare packages which must be loaded via `library()` in order to make specific non-standard R functions available across nodes.

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

### See Also

[SimFunctions](#), [createDesign](#), [Generate](#), [Analyse](#), [Summarise](#), [SimExtract](#), [reSummarise](#), [SimClean](#), [SimAnova](#), [SimResults](#), [SimCollect](#), [Attach](#), [AnalyseIf](#), [SimShiny](#), [manageWarnings](#), [runArraySimulation](#)

### Examples

```
#-----
# Example 1: Sampling distribution of mean

# This example demonstrate some of the simpler uses of SimDesign,
# particularly for classroom settings. The only factor varied in this simulation
# is sample size.

# skeleton functions to be saved and edited
SimFunctions()

#### Step 1 --- Define your conditions under study and create design data.frame

Design <- createDesign(N = c(10, 20, 30))

#~~~~~
#### Step 2 --- Define generate, analyse, and summarise functions

# help(Generate)
Generate <- function(condition, fixed_objects) {
  dat <- with(condition, rnorm(N, 10, 5)) # distributed N(10, 5)
  dat
}
```



```

res

# b) approach (store_results = TRUE by default)
res <- runSimulation(design=Design, replications=100,
                    generate=Generate, analyse=Analyse, summarise=Summarise)
res
SimResults(res)

# c) approach
Final <- runSimulation(design=Design, replications=100, save_results=TRUE,
                      generate=Generate, analyse=Analyse, summarise=Summarise)

# read-in all conditions (can be memory heavy)
res <- SimResults(Final)
res
head(res[[1]]$results)

# just first condition
res <- SimResults(Final, which=1)
head(res$results)
dplyr::tibble(res$condition, res$results)

# obtain empirical bootstrapped CIs during an initial run
# the simulation was completed (necessarily requires save_results = TRUE)
res <- runSimulation(design=Design, replications=1000, boot_method = 'basic',
                    generate=Generate, analyse=Analyse, summarise=Summarise)
res

# alternative bootstrapped CIs that uses saved results via reSummarise().
# Default directory save to:
dirname <- paste0('SimDesign-results_', unname(Sys.info()['nodename']), '/')
res <- reSummarise(summarise=Summarise, dir=dirname, boot_method = 'basic')
res

# remove the saved results from the hard-drive if you no longer want them
SimClean(results = TRUE)

## End(Not run)

#-----
# Example 2: t-test and Welch test when varying sample size, group sizes, and SDs

# skeleton functions to be saved and edited
SimFunctions()

## Not run:
# in real-world simulations it's often better/easier to save
# these functions directly to your hard-drive with
SimFunctions('my-simulation')

```

```

## End(Not run)

#### Step 1 --- Define your conditions under study and create design data.frame

Design <- createDesign(sample_size = c(30, 60, 90, 120),
                      group_size_ratio = c(1, 4, 8),
                      standard_deviation_ratio = c(.5, 1, 2))

Design

#~~~~~
#### Step 2 --- Define generate, analyse, and summarise functions

Generate <- function(condition, fixed_objects) {
  N <- condition$sample_size      # could use Attach() to make objects available
  grs <- condition$group_size_ratio
  sd <- condition$standard_deviation_ratio
  if(grs < 1){
    N2 <- N / (1/grs + 1)
    N1 <- N - N2
  } else {
    N1 <- N / (grs + 1)
    N2 <- N - N1
  }
  group1 <- rnorm(N1)
  group2 <- rnorm(N2, sd=sd)
  dat <- data.frame(group = c(rep('g1', N1), rep('g2', N2)), DV = c(group1, group2))
  dat
}

Analyse <- function(condition, dat, fixed_objects) {
  welch <- t.test(DV ~ group, dat)$p.value
  independent <- t.test(DV ~ group, dat, var.equal=TRUE)$p.value

  # In this function the p values for the t-tests are returned,
  # and make sure to name each element, for future reference
  ret <- nc(welch, independent)
  ret
}

Summarise <- function(condition, results, fixed_objects) {
  #find results of interest here (e.g., alpha < .1, .05, .01)
  ret <- EDR(results, alpha = .05)
  ret
}

#~~~~~
#### Step 3 --- Collect results by looping over the rows in design

# first, test to see if it works
res <- runSimulation(design=Design, replications=2,
                   generate=Generate, analyse=Analyse, summarise=Summarise)

res

```

```

## Not run:
# complete run with 1000 replications per condition
res <- runSimulation(design=Design, replications=1000, parallel=TRUE,
                    generate=Generate, analyse=Analyse, summarise=Summarise)
res
View(res)

## save final results to a file upon completion, and play a beep when done
runSimulation(design=Design, replications=1000, parallel=TRUE, filename = 'mysim',
             generate=Generate, analyse=Analyse, summarise=Summarise, beep=TRUE)

## same as above, but send a notification via Pushbullet upon completion
library(RPushbullet) # read-in default JSON file
runSimulation(design=Design, replications=1000, parallel=TRUE, filename = 'mysim',
             generate=Generate, analyse=Analyse, summarise=Summarise,
             notification = 'complete')

## Submit as RStudio job (requires job package and active RStudio session)
job::job({
  res <- runSimulation(design=Design, replications=100,
                    generate=Generate, analyse=Analyse, summarise=Summarise)
}, title='t-test simulation')
res # object res returned to console when completed

## Debug the generate function. See ?browser for help on debugging
## Type help to see available commands (e.g., n, c, where, ...),
## ls() to see what has been defined, and type Q to quit the debugger
runSimulation(design=Design, replications=1000,
             generate=Generate, analyse=Analyse, summarise=Summarise,
             parallel=TRUE, debug='generate')

## Alternatively, place a browser() within the desired function line to
## jump to a specific location
Summarise <- function(condition, results, fixed_objects) {
  #find results of interest here (e.g., alpha < .1, .05, .01)
  browser()
  ret <- EDR(results[,nms], alpha = .05)
  ret
}

## The following debugs the analyse function for the
## second row of the Design input
runSimulation(design=Design, replications=1000,
             generate=Generate, analyse=Analyse, summarise=Summarise,
             parallel=TRUE, debug='analyse-2')

#####
## EXTRA: To run the simulation on a user-define cluster, use the following setup (not run)
## Network linked via ssh (two way ssh key-paired connection must be
## possible between master and slave nodes)
##

```

```

## Define IP addresses, including primary IP
primary <- '192.168.2.20'
IPs <- list(
  list(host=primary, user='phil', ncore=8),
  list(host='192.168.2.17', user='phil', ncore=8)
)
spec <- lapply(IPs, function(IP)
  rep(list(list(host=IP$host, user=IP$user)), IP$ncore))
spec <- unlist(spec, recursive=FALSE)

cl <- parallel::makeCluster(type='PSOCK', master=primary, spec=spec)
res <- runSimulation(design=Design, replications=1000, parallel = TRUE,
  generate=Generate, analyse=Analyse, summarise=Summarise, cl=cl)

## Using parallel='future' to allow the future framework to be used instead
library(future) # future structure to be used internally
plan(multisession) # specify different plan (default is sequential)

res <- runSimulation(design=Design, replications=100, parallel='future',
  generate=Generate, analyse=Analyse, summarise=Summarise)
head(res)

# The progressr package is used for progress reporting with futures. To redefine
# use progressr::handlers() (see below)
library(progressr)
with_progress(res <- runSimulation(design=Design, replications=100, parallel='future',
  generate=Generate, analyse=Analyse, summarise=Summarise))
head(res)

# re-define progressr's bar (below requires cli)
handlers(handler_pbccl(
  adjust = 1.0,
  complete = function(s) cli::bg_red(cli::col_black(s)),
  incomplete = function(s) cli::bg_cyan(cli::col_black(s))
))

with_progress(res <- runSimulation(design=Design, replications=100, parallel='future',
  generate=Generate, analyse=Analyse, summarise=Summarise))

# reset future computing plan when complete (good practice)
plan(sequential)

#####

##### Post-analysis: Analyze the results via functions like lm() or SimAnova(), and create
##### tables(dplyr) or plots (ggplot2) to help visualize the results.
##### This is where you get to be a data analyst!

library(dplyr)
res %>% summarise(mean(welch), mean(independent))
res %>% group_by(standard_deviation_ratio, group_size_ratio) %>%
  summarise(mean(welch), mean(independent))

```

```

# quick ANOVA analysis method with all two-way interactions
SimAnova( ~ (sample_size + group_size_ratio + standard_deviation_ratio)^2, res,
  rates = TRUE)

# or more specific ANOVAs
SimAnova(independent ~ (group_size_ratio + standard_deviation_ratio)^2,
  res, rates = TRUE)

# make some plots
library(ggplot2)
library(tidyr)
dd <- res %>%
  select(group_size_ratio, standard_deviation_ratio, welch, independent) %>%
  pivot_longer(cols=c('welch', 'independent'), names_to = 'stats')
dd

ggplot(dd, aes(factor(group_size_ratio), value)) + geom_boxplot() +
  geom_abline(intercept=0.05, slope=0, col = 'red') +
  geom_abline(intercept=0.075, slope=0, col = 'red', linetype='dotted') +
  geom_abline(intercept=0.025, slope=0, col = 'red', linetype='dotted') +
  facet_wrap(~stats)

ggplot(dd, aes(factor(group_size_ratio), value, fill = factor(standard_deviation_ratio))) +
  geom_boxplot() + geom_abline(intercept=0.05, slope=0, col = 'red') +
  geom_abline(intercept=0.075, slope=0, col = 'red', linetype='dotted') +
  geom_abline(intercept=0.025, slope=0, col = 'red', linetype='dotted') +
  facet_grid(stats~standard_deviation_ratio) +
  theme(legend.position = 'none')

## End(Not run)

```

---

rValeMaurelli

*Generate non-normal data with Vale & Maurelli's (1983) method*


---

## Description

Generate multivariate non-normal distributions using the third-order polynomial method described by Vale & Maurelli (1983). If only a single variable is generated then this function is equivalent to the method described by Fleishman (1978).

## Usage

```

rValeMaurelli(
  n,
  mean = rep(0, nrow(sigma)),
  sigma = diag(length(mean)),
  skew = rep(0, nrow(sigma)),

```



```

    kurt = rep(0, nrow(sigma))
  )

```

### Arguments

|       |  |
|-------|--|
| n     | number of samples to draw  |
| mean  | a vector of k elements for the mean of the variables                   |
| sigma | desired k x k covariance matrix between bivariate non-normal variables |
| skew  | a vector of k elements for the skewness of the variables               |
| kurt  | a vector of k elements for the kurtosis of the variables               |

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

- Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248
- Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953
- Fleishman, A. I. (1978). A method for simulating non-normal distributions. *Psychometrika*, 43, 521-532.
- Vale, C. & Maurelli, V. (1983). Simulating multivariate nonnormal distributions. *Psychometrika*, 48(3), 465-471.

### Examples

```

set.seed(1)

# univariate with skew
nonnormal <- rValeMaurelli(10000, mean=10, sigma=5, skew=1, kurt=3)
# psych::describe(nonnormal)

# multivariate with skew and kurtosis
n <- 10000
r12 <- .4
r13 <- .9
r23 <- .1
cor <- matrix(c(1,r12,r13,r12,1,r23,r13,r23,1),3,3)
sk <- c(1.5,1.5,0.5)
ku <- c(3.75,3.5,0.5)

nonnormal <- rValeMaurelli(n, sigma=cor, skew=sk, kurt=ku)
# cor(nonnormal)
# psych::describe(nonnormal)

```

Serlin2000

*Empirical detection robustness method suggested by Serlin (2000)***Description**

Hypothesis test to determine whether an observed empirical detection rate, coupled with a given robustness interval, statistically differs from the population value. Uses the methods described by Serlin (2000) as well to generate critical values (similar to confidence intervals, but define a fixed window of robustness). Critical values may be computed without performing the simulation experiment (hence, can be obtained a priori).

**Usage**

```
Serlin2000(p, alpha, delta, R, CI = 0.95)
```

**Arguments**

|       |   |
|-------|---|
| p     | (optional) a vector containing the empirical detection rate(s) to be tested. Omitting this input will compute only the CV1 and CV2 values, while including this input will perform a one-sided hypothesis test for robustness |
| alpha | Type I error rate (e.g., often set to .05)  |
| delta | (optional) symmetric robustness interval around alpha (e.g., a value of .01 when alpha = .05 would test the robustness window .04-.06)  |
| R     | number of replications used in the simulation   |
| CI    | confidence interval for alpha as a proportion. Default of 0.95 indicates a 95% interval   |

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

- Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)
- Serlin, R. C. (2000). Testing for Robustness in Monte Carlo Studies. *Psychological Methods*, 5, 230-240.
- Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. [doi:10.1080/10691898.2016.1246953](https://doi.org/10.1080/10691898.2016.1246953)

**Examples**

```
# Cochran's criteria at alpha = .05 (i.e., 0.5 +- .01), assuming N = 2000
Serlin2000(p = .051, alpha = .05, delta = .01, R = 2000)

# Bradley's liberal criteria given p = .06 and .076, assuming N = 1000
Serlin2000(p = .060, alpha = .05, delta = .025, R = 1000)
Serlin2000(p = .076, alpha = .05, delta = .025, R = 1000)

# multiple p-values
Serlin2000(p = c(.05, .06, .07), alpha = .05, delta = .025, R = 1000)

# CV values computed before simulation performed
Serlin2000(alpha = .05, R = 2500)
```

SFA

*Surrogate Function Approximation via the Generalized Linear Model***Description**

Given a simulation that was executed with `runSimulation`, potentially with the argument `store_results = TRUE` to store the unsummarised analysis results, fit a surrogate function approximation (SFA) model to the results and (optionally) perform a root-solving step to solve a target quantity. See Schoemann et al. (2014) for details.

**Usage**

```
SFA(
  results,
  formula,
  family = "binomial",
  b = NULL,
  design = NULL,
  CI = 0.95,
  interval = NULL,
  ...
)

## S3 method for class 'SFA'
print(x, ...)
```

**Arguments**

|                      |  |
|----------------------|--|
| <code>results</code> | data returned from <code>runSimulation</code> . This can be the original results object or the extracted results stored when using <code>store_results = TRUE</code> included to store the analysis results. |
| <code>formula</code> | formula to specify for the regression model  |

|          |  |
|----------|--|
| family   | character vector indicating the family of GLMs to use (see <a href="#">family</a> )  |
| b        | (optional) Target quantity to use for root solving given the fitted surrogate function (e.g., find sample size associated with SFA implied power of .80)                                   |
| design   | (optional) data.frame object containing all the information relevant for the surrogate model (passed to newdata in <a href="#">predict</a> ) with an NA value in the variable to be solved |
| CI       | advertised confidence interval of SFA prediction around solved target  |
| interval | interval to be passed to <a href="#">uniroot</a> if not specified then the lowest and highest values from results for the respective variable will be used                                 |
| ...      | additional arguments to pass to <a href="#">glm</a>  |
| x        | an object of class SFA   |

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)

Schoemann, A. M., Miller, P., Pornprasertmanit, S., and Wu, W. (2014). Using Monte Carlo simulations to determine power and sample size for planned missing designs. *International Journal of Behavioral Development*, SAGE Publications, 38, 471-479.

**See Also**

[runSimulation](#), [SimSolve](#)

**Examples**

```
## Not run:

# create long Design object to fit surrogate over
Design <- createdesign(N = 100:500,
                      d = .2)

Design

#~~~~~
#### Step 2 --- Define generate, analyse, and summarise functions

Generate <- function(condition, fixed_objects) {
  Attach(condition)
  group1 <- rnorm(N)
  group2 <- rnorm(N, mean=d)
  dat <- data.frame(group = gl(2, N, labels=c('G1', 'G2')),
                   DV = c(group1, group2))
  dat
}
```

```

Analyse <- function(condition, dat, fixed_objects) {
  p <- c(p = t.test(DV ~ group, dat, var.equal=TRUE)$p.value)
  p
}

Summarise <- function(condition, results, fixed_objects) {
  ret <- EDR(results, alpha = .05)
  ret
}

#~~~~~
#### Step 3 --- Estimate power over N

# Use small number of replications given range of sample sizes
## note that due to the lower replications disabling the
## RAM printing will help reduce overhead

sim <- runSimulation(design=Design, replications=10,
                    generate=Generate, analyse=Analyse,
                    summarise=Summarise, store_results=TRUE, save=FALSE,
                    progress=FALSE, control=list(print_RAM=FALSE))

sim

# total of 4010 replication
sum(sim$REPLICATIONS)

# use the unsummarised results for the SFA, and include p.values < alpha
sim_results <- SimResults(sim)
sim_results <- within(sim_results, sig <- p < .05)
sim_results

# fitted model
sfa <- SFA(sim_results, formula = sig ~ N)
sfa
summary(sfa)

# plot the observed and SFA expected values
plot(p ~ N, sim, las=1, pch=16, main='Rejection rates with R=10')
pred <- predict(sfa, type = 'response')
lines(sim_results$N, pred, col='red', lty=2)

# fitted model + root-solved solution given f(.) = b,
# where b = target power of .8
design <- data.frame(N=NA, d=.2)
sfa.root <- SFA(sim_results, formula = sig ~ N,
               b=.8, design=design)

sfa.root

# true root
pwr::pwr.t.test(power=.8, d=.2)

```

```
#####
# example with smaller range but higher precision
Design <- createDesign(N = 375:425,
                      d = .2)

Design

sim2 <- runSimulation(design=Design, replications=100,
                     generate=Generate, analyse=Analyse,
                     summarise=Summarise, store_results=TRUE, save=FALSE,
                     progress=FALSE, control=list(print_RAM=FALSE))

sim2
sum(sim2$REPLICATIONS) # more replications in total

# use the unsummarised results for the SFA, and include p.values < alpha
sim_results <- SimResults(sim2)
sim_results <- within(sim_results, sig <- p < .05)
sim_results

# fitted model
sfa <- SFA(sim_results, formula = sig ~ N)
sfa
summary(sfa)

# plot the observed and SFA expected values
plot(p ~ N, sim2, las=1, pch=16, main='Rejection rates with R=100')
pred <- predict(sfa, type = 'response')
lines(sim_results$N, pred, col='red', lty=2)

# fitted model + root-solved solution given f(.) = b,
# where b = target power of .8
design <- data.frame(N=NA, d=.2)
sfa.root <- SFA(sim_results, formula = sig ~ N,
               b=.8, design=design, interval=c(100, 500))

sfa.root

# true root
pwr::pwr.t.test(power=.8, d=.2)

#####
# vary multiple parameters (e.g., sample size + effect size) to fit
# multi-parameter surrogate

Design <- createDesign(N = seq(from=10, to=500, by=10),
                      d = seq(from=.1, to=.5, by=.1))

Design

sim3 <- runSimulation(design=Design, replications=50,
                     generate=Generate, analyse=Analyse,
                     summarise=Summarise, store_results=TRUE, save=FALSE,
                     progress=FALSE, control=list(print_RAM=FALSE))

sim3
sum(sim3$REPLICATIONS)
```

```

# use the unsummarised results for the SFA, and include p.values < alpha
sim_results <- SimResults(sim3)
sim_results <- within(sim_results, sig <- p < .05)
sim_results

# additive effects (logit(sig) ~ N + d)
sfa0 <- SFA(sim_results, formula = sig ~ N+d)
sfa0

# multiplicative effects (logit(sig) ~ N + d + N:d)
sfa <- SFA(sim_results, formula = sig ~ N*d)
sfa

# multiplicative better fit (sample size interacts with effect size)
anova(sfa0, sfa, test = "LRT")
summary(sfa)

# plot the observed and SFA expected values
library(ggplot2)
sim3$pred <- predict(sfa, type = 'response', newdata=sim3)
ggplot(sim3, aes(N, p, color = factor(d))) +
  geom_point() + geom_line(aes(y=pred)) +
  facet_wrap(~factor(d))

# fitted model + root-solved solution given f(.) = b,
# where b = target power of .8
design <- data.frame(N=NA, d=.2)
sfa.root <- SFA(sim_results, formula = sig ~ N * d,
               b=.8, design=design, interval=c(100, 500))
sfa.root

# true root
pwr::pwr.t.test(power=.8, d=.2)

# root prediction where d *not* used in original data
design <- data.frame(N=NA, d=.25)
sfa.root <- SFA(sim_results, formula = sig ~ N * d,
               b=.8, design=design, interval=c(100, 500))
sfa.root

# true root
pwr::pwr.t.test(power=.8, d=.25)

## End(Not run)

```

## Description

Given the results from a simulation with `runSimulation` form an ANOVA table (without p-values) with effect sizes based on the eta-squared statistic. These results provide approximate indications of observable simulation effects, therefore these ANOVA-based results are generally useful as exploratory rather than inferential tools.

## Usage

```
SimAnova(formula, dat, subset = NULL, rates = TRUE)
```

## Arguments

|                      |   |
|----------------------|---|
| <code>formula</code> | an R formula generally of a form suitable for <code>lm</code> or <code>aov</code> . However, if the dependent variable (left side of the equation) is omitted then all the dependent variables in the simulation will be used and the result will return a list of analyses                   |
| <code>dat</code>     | an object returned from <code>runSimulation</code> of class 'SimDesign'   |
| <code>subset</code>  | an optional argument to be passed to <code>subset</code> with the same name. Used to subset the results object while preserving the associated attributes   |
| <code>rates</code>   | logical; does the dependent variable consist of rates (e.g., returned from <code>ECR</code> or <code>EDR</code> )? Default is <code>TRUE</code> , which will use the logit of the DV to help stabilize the proportion-based summary statistics when computing the parameters and effect sizes |

## Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

## References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. [doi:10.1080/10691898.2016.1246953](https://doi.org/10.1080/10691898.2016.1246953)

## Examples

```
data(BF_sim)

# all results (not usually good to mix Power and Type I results together)
SimAnova(alpha.05.F ~ (groups_equal + distribution)^2, BF_sim)

# only use anova for Type I error conditions
SimAnova(alpha.05.F ~ (groups_equal + distribution)^2, BF_sim, subset = var_ratio == 1)

# run all DVs at once using the same formula
SimAnova(~ groups_equal * distribution, BF_sim, subset = var_ratio == 1)
```



---

|          |   |
|----------|---|
| SimCheck | <i>Check for missing files in array simulations</i> |
|----------|---|

---

**Description**

Given the saved files from a [runArraySimulation](#) remote evaluation check whether all .rds files have been saved. If missing the missing row condition numbers will be returned.

**Usage**

```
SimCheck(dir = NULL, files = NULL, min = 1L, max = NULL)
```

**Arguments**

|       |   |
|-------|---|
| dir   | character vector input indicating the directory containing the .rds files (see files)                         |
| files | vector of file names referring to the saved simulation files. E.g. c('mysim-1.rds', 'mysim-2.rds', ...)       |
| min   | minimum number after the '-' delimitator. Default is 1  |
| max   | maximum number after the '-' delimitator. If not specified is extracted from the attributes in the first file |

**Value**

returns an invisible TRUE if all the files are present and FALSE otherwise

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)

**See Also**

[runArraySimulation](#), [SimCollect](#)

**Examples**

```
## Not run:  
  
# if files are in mysimfiles/ directory  
SimCheck('mysimfiles')  
  
# specifying files explicitly
```

```
setwd('mysimfiles/')
SimCheck(files=dir())
```

```
## End(Not run)
```

---

|          |  |
|----------|--|
| SimClean | <i>Removes/cleans files and folders that have been saved</i> |
|----------|--|

---

### Description

This function is mainly used in pilot studies where results and datasets have been temporarily saved by [runSimulation](#) but should be removed before beginning the full Monte Carlo simulation (e.g., remove files and folders which contained bugs/biased results).

### Usage

```
SimClean(
  ...,
  dirs = NULL,
  temp = TRUE,
  results = FALSE,
  seeds = FALSE,
  save_details = list()
)
```

### Arguments

|              |   |
|--------------|---|
| ...          | one or more character objects indicating which files to remove. Used to remove .rds files which were saved with <a href="#">saveRDS</a> or when using the save and filename inputs to <a href="#">runSimulation</a> |
| dirs         | a character vector indicating which directories to remove   |
| temp         | logical; remove the temporary file saved when passing save = TRUE?  |
| results      | logical; remove the .rds results files saved when passing save_results = TRUE?  |
| seeds        | logical; remove the seed files saved when passing save_seeds = TRUE?  |
| save_details | a list pertaining to information about how and where files were saved (see the corresponding list in <a href="#">runSimulation</a> )  |

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

## References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

## See Also

[runSimulation](#)

## Examples

```
## Not run:

# remove file called 'results.rds'
SimClean('results.rds')

# remove default temp file
SimClean()

# remove customized saved-results directory called 'mydir'
SimClean(results = TRUE, save_details = list(save_results_dirname = 'mydir'))

## End(Not run)
```

---

SimCollect

*Collapse separate simulation files into a single result*

---

## Description

This function collects and aggregates the results from SimDesign's [runSimulation](#) into a single objects suitable for post-analyses, or combines all the saved results directories and combines them into one. This is useful when results are run piece-wise on one node (e.g., 500 replications in one batch, 500 again at a later date, though be careful about the [set.seed](#) use as the random numbers will tend to correlate the more it is used) or run independently across different nodes/computing cores (e.g., see [runArraySimulation](#)).

## Usage

```
SimCollect(
  dir = NULL,
  files = NULL,
  filename = NULL,
  select = NULL,
  check.only = FALSE,
  target.reps = NULL,
```

```
warning_details = FALSE,
error_details = TRUE
)

aggregate_simulations(...)
```

### Arguments

|                              |  |
|------------------------------|--|
| <code>dir</code>             | a character vector pointing to the directory name containing the <code>.rds</code> files. All <code>.rds</code> files in this directory will be used after first checking their status with <a href="#">SimCheck</a> . For greater specificity use the <code>files</code> argument   |
| <code>files</code>           | a character vector containing the names of the simulation's final <code>.rds</code> files.   |
| <code>filename</code>        | (optional) name of <code>.rds</code> file to save aggregate simulation file to. If not specified then the results will only be returned in the R console.  |
| <code>select</code>          | a character vector indicating columns to variables to select from the <code>SimExtract(what='results')</code> information. This is mainly useful when RAM is an issue given simulations with many stored estimates. Default includes the results objects in their entirety, though to omit all internally stored simulation results pass the character <code>'NONE'</code> |
| <code>check.only</code>      | logical; for larger simulations file sets, such as those generated by <a href="#">runArraySimulation</a> , return the design conditions that do not satisfy the <code>target.reps</code> and throw warning if files are unexpectedly missing   |
| <code>target.reps</code>     | (optional) number of replications to check against to evaluate whether the simulation files returned the desired number of replications. If missing, the highest detected value from the collected set of replication information will be used   |
| <code>warning_details</code> | logical; include the aggregate of the warnings to be extracted via <a href="#">SimExtract</a> ?  |
| <code>error_details</code>   | logical; include the aggregate of the errors to be extracted via <a href="#">SimExtract</a> ?  |
| <code>...</code>             | not used   |

### Value

returns a `data.frame/tibble` with the (weighted) average/aggregate of the simulation results

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

- Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the `SimDesign` Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248
- Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

### See Also

[runSimulation](#), [runArraySimulation](#), [SimCheck](#)

**Examples**

```

## Not run:

setwd('my_working_directory')

## run simulations to save the .rds files (or move them to the working directory)
# seeds1 <- genSeeds(design)
# seeds2 <- genSeeds(design, old.seeds=seeds1)
# ret1 <- runSimulation(design, ..., seed=seeds1, filename='file1')
# ret2 <- runSimulation(design, ..., seed=seeds2, filename='file2')

# saves to the hard-drive and stores in workspace
final <- SimCollect(files = c('file1.rds', 'file2.rds'))
final

# If filename not included, can be extracted from results
# files <- c(SimExtract(ret1, 'filename'), SimExtract(ret2, 'filename'))
# final <- SimCollect(files = files)

#####
# Example where each row condition is repeated, evaluated independently,
# and later collapsed into a single analysis object

# Each condition repeated four times (hence, replications
# should be set to desired.reps/4)
Design <- createDesign(mu = c(0,5),
                      N = c(30, 60))
Design

# assume the N=60 takes longer, and should be spread out across more arrays
Design_long <- expandDesign(Design, c(2,2,4,4))
Design_long

replications <- c(rep(50, 4), rep(25,8))
data.frame(Design_long, replications)

#-----

Generate <- function(condition, fixed_objects) {
  dat <- with(condition, rnorm(N, mean=mu))
  dat
}

Analyse <- function(condition, dat, fixed_objects) {
  ret <- c(mean=mean(dat), SD=sd(dat))
  ret
}

Summarise <- function(condition, results, fixed_objects) {
  ret <- colMeans(results)
  ret
}

```

```

}

#-----

# create directory to store all final simulation files
dir.create('sim_files/')

iseed <- genSeeds()

# distribute jobs independently
sapply(1:nrow(Design_long), \(i) {
  runArraySimulation(design=Design_long, replications=replications,
                    generate=Generate, analyse=Analyse, summarise=Summarise,
                    arrayID=i, dirname='sim_files/', filename='job', iseed=iseed)
}) |> invisible()

# check that all replications satisfy target
SimCollect('sim_files/', check.only = TRUE)

# this would have been returned were the target.rep supposed to be 1000
SimCollect('sim_files/', check.only = TRUE, target.reps=1000)

# aggregate into single object
sim <- SimCollect('sim_files/')
sim

SimClean(dir='sim_files/')

## End(Not run)

```

---

 SimDesign

---

*Structure for Organizing Monte Carlo Simulation Designs*


---

## Description

Structure for Organizing Monte Carlo Simulation Designs

## Details

Provides tools to help organize Monte Carlo simulations in R. The package controls the structure and back-end of Monte Carlo simulations by utilizing a general generate-analyse-summarise strategy. The functions provided control common simulation issues such as re-simulating non-convergent results, support parallel back-end computations with proper random number generation within each simulation condition, save and restore temporary files, aggregate results across independent nodes, and provide native support for debugging. The primary function for organizing the simulations is `runSimulation`, while for array jobs submitting to HPC clusters (e.g., SLURM) see `runArraySimulation` and the associated package vignettes.

For an in-depth tutorial of the package please refer to Chalmers and Adkins (2020; [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)). For an earlier didactic presentation of the package users can refer to Sigal and

Chalmers (2016; [doi:10.1080/10691898.2016.1246953](https://doi.org/10.1080/10691898.2016.1246953)). Finally, see the associated wiki on Github (<https://github.com/philchalmers/SimDesign/wiki>) for other tutorial material, examples, and applications of SimDesign to real-world simulations.

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. [doi:10.1080/10691898.2016.1246953](https://doi.org/10.1080/10691898.2016.1246953)

---

SimExtract

*Function to extract extra information from SimDesign objects*

---

### Description

Function used to extract any error or warnings messages, the seeds associated with any error or warning messages, and any analysis results that were stored in the final simulation object.

### Usage

```
SimExtract(object, what, fuzzy = TRUE, append = TRUE)
```

### Arguments

|        |  |
|--------|--|
| object | object returned from <a href="#">runSimulation</a>   |
| what   | character indicating what information to extract. Possible inputs include 'errors' to return a tibble object containing counts of any error messages, 'warnings' to return a data.frame object containing counts of any warning messages, 'seeds' for the specified random number generation seeds, 'Random.seeds' for the complete list of .Random.seed states across replications (only stored when <code>runSimulation(..., control = list(store_Random.seeds=TRUE))</code> ), 'error_seeds' and 'warning_seeds' to extract the associated .Random.seed values associated with the ERROR/WARNING messages, 'results' to extract the simulation results if the option <code>store_results</code> was passed to <a href="#">runSimulation</a> , 'filename' and 'save_results_dirname' for extracting the saved file/directory name information (if used), and 'summarise' if the <a href="#">Summarise</a> definition returned a named list rather than a named numeric vector.<br>Note that 'warning_seeds' are not stored automatically in simulations and require passing <code>store_warning_seeds = TRUE</code> to <a href="#">runSimulation</a> . |

|        |  |
|--------|--|
| fuzzy  | logical; use fuzzy string matching to reduce effectively identical messages? For example, when attempting to invert a matrix the error message " <i>System is computationally singular: reciprocal condition number = 1.92747e-17</i> " and " <i>System is computationally singular: reciprocal condition number = 2.15321e-16</i> " are effectively the same, and likely should be reported in the same columns of the extracted output |
| append | logical; append the design conditions when extracting error/warning messages?  |

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

### Examples

```
## Not run:

Generate <- function(condition, fixed_objects) {
  int <- sample(1:10, 1)
  if(int > 5) warning('GENERATE WARNING: int greater than 5')
  if(int == 1) stop('GENERATE ERROR: integer is 1')
  rnorm(5)
}

Analyse <- function(condition, dat, fixed_objects) {
  int <- sample(1:10, 1)
  if(int > 5) warning('ANALYSE WARNING: int greater than 5')
  if(int == 1) stop('ANALYSE ERROR: int is 1')
  c(ret = 1)
}

Summarise <- function(condition, results, fixed_objects) {
  mean(results)
}

res <- runSimulation(replications = 100, seed=1234, verbose=FALSE,
                    generate=Generate, analyse=Analyse, summarise=Summarise)

res

SimExtract(res, what = 'errors')
SimExtract(res, what = 'warnings')
seeds <- SimExtract(res, what = 'error_seeds')
seeds[,1:3]
```



```
# replicate a specific error for debugging (type Q to exit debugger)
res <- runSimulation(replications = 100, load_seed=seeds[,1], debug='analyse',
                    generate=Generate, analyse=Analyse, summarise=Summarise)

## End(Not run)
```

---

SimFunctions

*Template-based generation of the Generate-Analyse-Summarise functions*

---

## Description

This function prints template versions of the required Design and Generate-Analyse-Summarise functions for SimDesign to run simulations. Templated output comes complete with the correct inputs, class of outputs, and optional comments to help with the initial definitions. Use this at the start of your Monte Carlo simulation study. Following the definition of the SimDesign template file please refer to detailed the information in [runSimulation](#) for how to edit this template to make a working simulation study.

## Usage

```
SimFunctions(
  filename = NULL,
  dir = getwd(),
  save_structure = "single",
  extra_file = FALSE,
  nAnalyses = 1,
  nGenerate = 1,
  summarise = TRUE,
  comments = FALSE,
  openFiles = TRUE,
  spin_header = TRUE,
  SimSolve = FALSE
)
```

## Arguments

|          |   |
|----------|---|
| filename | a character vector indicating whether the output should be saved to two respective files containing the simulation design and the functional components, respectively. Using this option is generally the recommended approach when beginning to write a Monte Carlo simulation |
| dir      | the directory to write the files to. Default is the working directory   |

|                |   |
|----------------|---|
| save_structure | character indicating the number of files to break the simulation code into when filename is included (default is 'single' for one file). When save_structure = 'double' the output is saved to two separate files containing the functions and design definitions, and when save_structure = 'all' the generate, analyse, summarise, and execution code area all saved into separate files. The purpose for this structure is because multiple structured files often makes organization and debugging slightly easier larger Monte Carlo simulations, though in principle all files could be stored into a single R script |
| extra_file     | logical; should and extra file be saved containing user-defined functions or objects? Default is FALSE  |
| nAnalyses      | number of analysis functions to create (default is 1). Increasing the value of this argument when independent analysis are being performed allows function definitions to be better partitioned and potentially more modular  |
| nGenerate      | number of generate functions to create (default is 1). Increase the value of this argument when when the data generation functions are very different and should be isolated from each other (otherwise, if there is much in common between the generate steps, the default of 1 should be preferred). Otherwise, if nGenerate == 0 then no generate function will be provided and instead this data-generation step can be defined in the analysis function(s) (only recommended for smaller simulations)  |
| summarise      | include summarise function? Default is TRUE   |
| comments       | logical; include helpful comments? Default is FALSE   |
| openFiles      | logical; after files have been generated, open them in your text editor (e.g., if Rstudio is running the scripts will open in a new tab)?   |
| spin_header    | logical; include a basic knitr::spin header to allow the simulation to be knitted? Default is TRUE. For those less familiar with spin documents see <a href="https://bookdown.org/yihui/rm">https://bookdown.org/yihui/rm</a> for further details   |
| SimSolve       | logical; should the template be generated that is intended for a SimSolve implementation? Default is FALSE  |

### Details

The recommended approach to organizing Monte Carlo simulation files is to first save the template generated by this function to the hard-drive by passing a suitable filename argument (which, if users are interacting with R via the RStudio IDE, will also open the template file after it has been saved). For larger simulations, two separate files could also be used (achieved by changing `out.files`), and may be easier for debugging/sourcing the simulation code; however, this is a matter of preference and does not change any functionality in the package.

### Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

### References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. [doi:10.20982/tqmp.16.4.p248](https://doi.org/10.20982/tqmp.16.4.p248)

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

### See Also

[runSimulation](#)

### Examples

```
SimFunctions()
SimFunctions(comments = TRUE) #with helpful comments

## Not run:

# write output files to a single file with comments
SimFunctions('mysim', comments = TRUE)

# Multiple analysis functions for optional partitioning
SimFunctions(nAnalyses = 2)
SimFunctions(nAnalyses = 3)

# Multiple analysis + generate functions
SimFunctions(nAnalyses = 2, nGenerate=2)

# save multiple files for the purpose of designing larger simulations
# (also include extra_file for user-defined objects/functions)
SimFunctions('myBigSim', save_structure = 'all',
  nAnalyses = 3, nGenerate=2, extra_file = TRUE)

## End(Not run)
```

---

SimResults

*Function to read in saved simulation results*

---

### Description

If [runSimulation](#) was passed the flag `save_results = TRUE` then the row results corresponding to the design object will be stored to a suitable sub-directory as individual `.rds` files. While users could use [readRDS](#) directly to read these files in themselves, this convenience function will read the desired rows in automatically given the returned object from the simulation. Can be used to read in 1 or more `.rds` files at once (if more than 1 file is read in then the result will be stored in a list).

### Usage

```
SimResults(obj, which, prefix = "results-row", wd = getwd())
```

**Arguments**

|        |  |
|--------|--|
| obj    | object returned from <code>runSimulation</code> where <code>save_results = TRUE</code> or <code>store_results</code> was used. If the former then the remaining function arguments can be useful for reading in specific files |
| which  | a numeric vector indicating which rows should be read in. If missing, all rows will be read in   |
| prefix | character indicating prefix used for stored files  |
| wd     | working directory; default is found with <code>getwd</code> .  |

**Value**

the returned result is either a nested list (when `length(which) > 1`) or a single list (when `length(which) == 1`) containing the simulation results. Each read-in result refers to a list of 4 elements:

`condition` the associate row (ID) and conditions from the respective design object

`results` the object with returned from the `analyse` function, potentially simplified into a matrix or `data.frame`

`errors` a table containing the message and number of errors that caused the generate-analyse steps to be rerun. These should be inspected carefully as they could indicate validity issues with the simulation that should be noted

`warnings` a table containing the message and number of non-fatal warnings which arose from the analyse step. These should be inspected carefully as they could indicate validity issues with the simulation that should be noted

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

**Examples**

```
## Not run:

# store results (default behaviour)
sim <- runSimulation(..., store_results = TRUE)
SimResults(sim)

# store results to drive if RAM issues are present
obj <- runSimulation(..., save_results = TRUE)

# row 1 results
```

```
row1 <- SimResults(obj, 1)

# rows 1:5, stored in a named list
rows_1to5 <- SimResults(obj, 1:5)

# all results
rows_all <- SimResults(obj)

## End(Not run)
```

---

SimShiny

*Generate a basic Monte Carlo simulation GUI template*

---

## Description

This function generates suitable stand-alone code from the shiny package to create simple web-interfaces for performing single condition Monte Carlo simulations. The template generated is relatively minimalistic, but allows the user to quickly and easily edit the saved files to customize the associated shiny elements as they see fit.

## Usage

```
SimShiny(filename = NULL, dir = getwd(), design, ...)
```

## Arguments

|          |   |
|----------|---|
| filename | an optional name of a text file to save the server and UI components (e.g., 'mysimGUI.R'). If omitted, the code will be printed to the R console instead        |
| dir      | the directory to write the files to. Default is the working directory   |
| design   | design object from <code>runSimulation</code>   |
| ...      | arguments to be passed to <code>runSimulation</code> . Note that the design object is not used directly, and instead provides options to be selected in the GUI |

## Author(s)

Phil Chalmers <rphilip.chalmers@gmail.com>

## References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

**See Also**[runSimulation](#)**Examples**

```
## Not run:

Design <- createDesign(sample_size = c(30, 60, 90, 120),
                      group_size_ratio = c(1, 4, 8),
                      standard_deviation_ratio = c(.5, 1, 2))

Generate <- function(condition, fixed_objects) {
  N <- condition$sample_size
  grs <- condition$group_size_ratio
  sd <- condition$standard_deviation_ratio
  if(grs < 1){
    N2 <- N / (1/grs + 1)
    N1 <- N - N2
  } else {
    N1 <- N / (grs + 1)
    N2 <- N - N1
  }
  group1 <- rnorm(N1)
  group2 <- rnorm(N2, sd=sd)
  dat <- data.frame(group = c(rep('g1', N1), rep('g2', N2)), DV = c(group1, group2))
  dat
}

Analyse <- function(condition, dat, fixed_objects) {
  welch <- t.test(DV ~ group, dat)
  ind <- t.test(DV ~ group, dat, var.equal=TRUE)

  # In this function the p values for the t-tests are returned,
  # and make sure to name each element, for future reference
  ret <- c(welch = welch$p.value, independent = ind$p.value)
  ret
}

Summarise <- function(condition, results, fixed_objects) {
  #find results of interest here (e.g., alpha < .1, .05, .01)
  ret <- EDR(results, alpha = .05)
  ret
}

# test that it works
# Final <- runSimulation(design=Design, replications=5,
#                       generate=Generate, analyse=Analyse, summarise=Summarise)

# print code to console
SimShiny(design=Design, generate=Generate, analyse=Analyse,
        summarise=Summarise, verbose=FALSE)
```

```

# save shiny code to file
SimShiny('app.R', design=Design, generate=Generate, analyse=Analyse,
         summarise=Summarise, verbose=FALSE)

# run the application
shiny::runApp()
shiny::runApp(launch.browser = TRUE) # in web-browser

## End(Not run)

```

---

 SimSolve

*One Dimensional Root (Zero) Finding in Simulation Experiments*


---

### Description

This function provides a stochastic root-finding approach to solving specific quantities in simulation experiments (e.g., solving for a specific sample size to meet a target power rate) using the Probabilistic Bisection Algorithm with Bolstering and Interpolations (ProBABLI; Chalmers, accepted). The structure follows the steps outlined in [runSimulation](#), however portions of the design input are taken as variables to be estimated rather than fixed, and the constant  $b$  is required in order to solve the root equation  $f(x) - b = 0$ . Stochastic root search is terminated based on the successive behavior of the  $x$  estimates. For even greater advertised accuracy with ProBABLI, termination criteria can be based on the width of the advertised predicting interval (via `predCI.tol`) or by specifying how long the investigator is willing to wait for the final estimates (via `wait.time`, where longer wait times lead to progressively better accuracy in the final estimates).

### Usage

```

SimSolve(
  design,
  interval,
  b,
  generate,
  analyse,
  summarise,
  replications = list(burnin.iter = 15L, burnin.reps = 100L, max.reps = 500L,
    min.total.reps = 9000L, increase.by = 10L),
  integer = TRUE,
  formula = y ~ poly(x, 2),
  family = "binomial",
  parallel = FALSE,
  cl = NULL,
  save = TRUE,
  resume = TRUE,
  method = "ProBABLI",
  wait.time = NULL,
  ncores = parallelly::availableCores(omit = 1L),

```

```

type = ifelse(.Platform$OS.type == "windows", "PSOCK", "FORK"),
maxiter = 100L,
check.interval = TRUE,
verbose = TRUE,
control = list(),
predCI = 0.95,
predCI.tol = NULL,
...
)

## S3 method for class 'SimSolve'
summary(object, tab.only = FALSE, reps.cutoff = 300, ...)

## S3 method for class 'SimSolve'
plot(x, y, ...)

```

### Arguments

|              |   |
|--------------|---|
| design       | a tibble or data.frame object containing the Monte Carlo simulation conditions to be studied, where each row represents a unique condition and each column a factor to be varied (see also <a href="#">createDesign</a> ). However, exactly one column of this object must be specified with NA placeholders to indicate that the missing value should be solved via the stochastic optimizer   |
| interval     | a vector of length two, or matrix with <code>nrow(design)</code> and two columns, containing the end-points of the interval to be searched. If a vector then the interval will be used for all rows in the supplied design object   |
| b            | a single constant used to solve the root equation $f(x) - b = 0$  |
| generate     | generate function. See <a href="#">runSimulation</a>  |
| analyse      | analysis function. See <a href="#">runSimulation</a>  |
| summarise    | summary function that returns a single number corresponding to a function evaluation $f(x)$ in the equation $f(x) = b$ to be solved as a root $f(x) - b = 0$ . Unlike in the standard <code>runSimulation()</code> definitions this input is required. For further information on this function specification, see <a href="#">runSimulation</a>  |
| replications | a named list or vector indicating the number of replication to use for each design condition per PBA iteration. By default the input is a list with the arguments <code>burnin.iter = 15L</code> , specifying the number of burn-in iterations to used, <code>burnin.reps = 100L</code> to indicate how many replications to use in each burn-in iteration, <code>max.reps = 500L</code> to prevent the replications from increasing higher than this number, <code>min.total.reps = 9000L</code> to avoid termination when very few replications have been explored (lower bound of the replication budget), and <code>increase.by = 10L</code> to indicate how many replications to increase after the burn-in stage. Unless otherwise specified these defaults will be used, but can be overwritten by explicit definition (e.g., <code>replications = list(increase.by = 25L)</code> )<br><br>Vector inputs can specify the exact replications for each iterations. As a general rule, early iterations should be relatively low for initial searches to avoid unnecessary computations for locating the approximate root, though the number of |



|           |  |
|-----------|--|
|           | replications should gradually increase to reduce the sampling variability as the PBA approaches the root.  |
| integer   | logical; should the values of the root be considered integer or numeric? If TRUE then bolstered directional decisions will be made in the pba function based on the collected sampling history throughout the search   |
| formula   | regression formula to use when interpolate = TRUE. Default fits an orthogonal polynomial of degree 2   |
| family    | family argument passed to <code>glm</code> . By default the 'binomial' family is used, as this function defaults to power analysis setups where isolated results passed to summarise will return 0/1s, however other families should be used had summarise returned something else (e.g., if solving for a particular standard error then a 'gaussian' family would be more appropriate).<br>Note that if individual results from the analyse steps should not be used (i.e., only the aggregate from summarise is meaningful) then set control = list(summarise.reg_data = TRUE) to override the default behavior, thereby using only the aggregate information and weights |
| parallel  | for parallel computing for slower simulation experiments (see <a href="#">runSimulation</a> for details)   |
| cl        | see <a href="#">runSimulation</a>  |
| save      | logical; store temporary file in case of crashes. If detected in the working directory will automatically be loaded to resume (see <a href="#">runSimulation</a> for similar behaviour)  |
| resume    | logical; if a temporary SimDesign file is detected should the simulation resume from this location? Keeping this TRUE is generally recommended, however this should be disabled if using SimSolve within <a href="#">runSimulation</a> to avoid reading improper save states   |
| method    | optimizer method to use. Default is the stochastic root-finder 'ProBAbLI', but can also be the deterministic options 'Brent' (which uses the function <a href="#">uniroot</a> ) or 'bisection' (for the classical bisection method). If using deterministic root-finders then replications must either equal a single constant to reflect the number of replication to use per deterministic iteration or be a vector of length maxiter to indicate the replications to use per iteration  |
| wait.time | (optional) argument passed to <a href="#">PBA</a> to indicate the time to wait (specified in minutes) per row in the Design object rather than using pre-determined termination criteria based on the estimates. For example, if three three conditions were defined in Design, and wait.time="5", then the total search time till terminate after 15 minutes regardless of independently specified termination criteria in control. Note that maxiter is still used alongside wait.time, therefore this should be increased as well (e.g., to maxiter = 1000)   |
| ncores    | see <a href="#">runSimulation</a>  |
| type      | type of cluster object to define. If type used in plot then can be 'density' to plot the density of the iteration history after the burn-in stage, 'iterations' for a bubble plot with inverse replication weights. If not specified then the default PBA plots are provided (see <a href="#">PBA</a> )  |
| maxiter   | the maximum number of iterations (default 100)   |

|                             |  |
|-----------------------------|--|
| <code>check.interval</code> | logical; should an initial check be made to determine whether $f(\text{interval}[1L])$ and $f(\text{interval}[2L])$ have opposite signs? If FALSE, the specified interval is assumed to contain a root, where $f(\text{interval}[1]) < 0$ and $f(\text{interval}[2]) > 0$ . Default is TRUE  |
| <code>verbose</code>        | logical; print information to the console?   |
| <code>control</code>        | a list of the algorithm control parameters. If not specified, the defaults described below are used.<br><code>tol</code> tolerance criteria for early termination (.1 for integer = TRUE searches; .00025 for non-integer searches)<br><code>rel.tol</code> relative tolerance criteria for early termination (default .0001)<br><code>k.success</code> number of consecutive tolerance success given <code>rel.tol</code> and <code>tol</code> criteria. Consecutive failures add -1 to the counter (default is 3)<br><code>bolster</code> logical; should the PBA evaluations use bolstering based on previous evaluations? Default is TRUE, though only applicable when <code>integer = TRUE</code><br><code>interpolate.R</code> number of replications to collect prior to performing the interpolation step (default is 3000 after accounting for data exclusion from <code>burnin.iter</code> ). Setting this to 0 will disable any interpolation computations<br><code>include_reps</code> logical; include a column in the condition elements to indicate how many replications are currently being evaluated? Mainly useful when further precision tuning within each ProBABLI iteration is desirable (e.g., for bootstrapping). Default is FALSE<br><code>summarise.reg_data</code> logical; should the aggregate results from Summarise (along with its associated weights) be used for the interpolation steps, or the raw data from the Analyse step? Set this to TRUE when the individual results from Analyse give less meaningful information |
| <code>predCI</code>         | advertised confidence interval probability for final model-based prediction of target <code>b</code> given the root input estimate. Returned as an element in the <code>summary()</code> list output   |
| <code>predCI.tol</code>     | (optional) rather than relying on the changes between successive estimates (default), if the predicting CI is consistently within this supplied tolerance input range then terminate. This provides termination behaviour based on the predicted precision of the root solutions rather than their stability history, and therefore can be used to obtain estimates with a particular level of advertised accuracy. For example, when solving for a sample size value ( <code>N</code> ) if the solution associated with <code>b = .80</code> requires that the advertised 95 is consistently between <code>[.795, .805]</code> then <code>predCI.tol = .01</code> to indicate this tolerance range  |
| <code>...</code>            | additional arguments to be pasted to <a href="#">PBA</a>   |
| <code>object</code>         | object of class 'SimSolve'   |
| <code>tab.only</code>       | logical; print only the (reduce) table of estimates?   |
| <code>reps.cutoff</code>    | integer indicating the rows to omit from output if the number of replications do not reach this value  |
| <code>x</code>              | object of class 'SimSolve'   |
| <code>y</code>              | design row to plot. If omitted defaults to 1   |

**Details**

Root finding is performed using a progressively bolstered version of the probabilistic bisection algorithm (PBA) to find the associated root given the noisy simulation objective function evaluations. Information is collected throughout the search to make more accurate predictions about the associated root via interpolation. If interpolations fail, then the last iteration of the PBA search is returned as the best guess.

**Value**

the filled-in design object containing the associated lower and upper interval estimates from the stochastic optimization

**Author(s)**

Phil Chalmers <rphilip.chalmers@gmail.com>

**References**

Chalmers, R. P. (in press). Solving Variables with Monte Carlo Simulation Experiments: A Stochastic Root-Solving Approach. *Psychological Methods*. DOI: 10.1037/met0000689

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

**See Also**

[SFA](#)

**Examples**

```
## Not run:

#####
## A Priori Power Analysis
#####

# GOAL: Find specific sample size in each group for independent t-test
# corresponding to a power rate of .8
#
# For ease of the setup, assume the groups are the same size, and the mean
# difference corresponds to Cohen's d values of .2, .5, and .8
# This example can be solved numerically using the pwr package (see below),
# though the following simulation setup is far more general and can be
# used for any generate-analyse combination of interest

# SimFunctions(SimSolve=TRUE)

#### Step 1 --- Define your conditions under study and create design.data.frame.
#### However, use NA placeholder for sample size as it must be solved,
#### and add desired power rate to object
```

```

Design <- createDesign(N = NA,
                      d = c(.2, .5, .8),
                      sig.level = .05)
Design # solve for NA's

#~~~~~
#### Step 2 --- Define generate, analyse, and summarise functions

Generate <- function(condition, fixed_objects) {
  Attach(condition)
  group1 <- rnorm(N)
  group2 <- rnorm(N, mean=d)
  dat <- data.frame(group = gl(2, N, labels=c('G1', 'G2')),
                    DV = c(group1, group2))
  dat
}

Analyse <- function(condition, dat, fixed_objects) {
  p <- t.test(DV ~ group, dat, var.equal=TRUE)$p.value
  p
}

Summarise <- function(condition, results, fixed_objects) {
  # Must return a single number corresponding to f(x) in the
  # root equation f(x) = b

  ret <- c(power = EDR(results, alpha = condition$sig.level))
  ret
}

#~~~~~
#### Step 3 --- Optimize N over the rows in design

### (For debugging) may want to see if simulation code works as intended first
### for some given set of inputs
# runSimulation(design=createDesign(N=100, d=.8, sig.level=.05),
#               replications=10, generate=Generate, analyse=Analyse,
#               summarise=Summarise)

# Initial search between N = [10,500] for each row using the default
# integer solver (integer = TRUE). In this example, b = target power
solved <- SimSolve(design=Design, b=.8, interval=c(10, 500),
                  generate=Generate, analyse=Analyse,
                  summarise=Summarise)

solved
summary(solved)
plot(solved, 1)
plot(solved, 2)
plot(solved, 3)

# also can plot median history and estimate precision
plot(solved, 1, type = 'history')

```

```

plot(solved, 1, type = 'density')
plot(solved, 1, type = 'iterations')

# verify with true power from pwr package
library(pwr)
pwr.t.test(d=.2, power = .8) # sig.level/alpha = .05 by default
pwr.t.test(d=.5, power = .8)
pwr.t.test(d=.8, power = .8)

# use estimated N results to see how close power was
N <- solved$N
pwr.t.test(d=.2, n=N[1])
pwr.t.test(d=.5, n=N[2])
pwr.t.test(d=.8, n=N[3])

# with rounding
N <- ceiling(solved$N)
pwr.t.test(d=.2, n=N[1])
pwr.t.test(d=.5, n=N[2])
pwr.t.test(d=.8, n=N[3])

### failing analytic formula, confirm results with more precise
### simulation via runSimulation()
### (not required, if accuracy is important then ProBABLI should be run longer)
# csolved <- solved
# csolved$N <- ceiling(solved$N)
# confirm <- runSimulation(design=csolved, replications=10000, parallel=TRUE,
#                           generate=Generate, analyse=Analyse,
#                           summarise=Summarise)
# confirm

# Similarly, terminate if the prediction interval is consistently predicted
# to be between [.795, .805]. Note that maxiter increased as well
solved_predCI <- SimSolve(design=Design, b=.8, interval=c(10, 500),
                        generate=Generate, analyse=Analyse, summarise=Summarise,
                        maxiter=200, predCI.tol=.01)

solved_predCI
summary(solved_predCI) # note that predCI.b are all within [.795, .805]

N <- solved_predCI$N
pwr.t.test(d=.2, n=N[1])
pwr.t.test(d=.5, n=N[2])
pwr.t.test(d=.8, n=N[3])

# Alternatively, and often more realistically, wait.time can be used
# to specify how long the user is willing to wait for a final estimate.
# Solutions involving more iterations will be more accurate,
# and therefore it is recommended to run the ProBABLI root-solver as long
# the analyst can tolerate if the most accurate estimates are desired.
# Below executes the simulation for 5 minutes for each condition up
# to a maximum of 1000 iterations, terminating based on whichever occurs first

solved_5min <- SimSolve(design=Design, b=.8, interval=c(10, 500),

```

```

        generate=Generate, analyse=Analyse, summarise=Summarise,
        wait.time="5", maxiter=1000)
solved_5min
summary(solved_5min)

# use estimated N results to see how close power was
N <- solved_5min$N
pwr.t.test(d=.2, n=N[1])
pwr.t.test(d=.5, n=N[2])
pwr.t.test(d=.8, n=N[3])

#-----

#####
## Sensitivity Analysis
#####

# GOAL: solve effect size d given sample size and power inputs (inputs
# for root no longer required to be an integer)

# Generate-Analyse-Summarise functions identical to above, however
# Design input includes NA for d element
Design <- createDesign(N = c(100, 50, 25),
                      d = NA,
                      sig.level = .05)
Design # solve for NA's

#-----
#### Step 2 --- Define generate, analyse, and summarise functions (same as above)

#-----
#### Step 3 --- Optimize d over the rows in design
# search between d = [.1, 2] for each row

# In this example, b = target power
# note that integer = FALSE to allow smooth updates of d
solved <- SimSolve(design=Design, b = .8, interval=c(.1, 2),
                  generate=Generate, analyse=Analyse,
                  summarise=Summarise, integer=FALSE)

solved
summary(solved)
plot(solved, 1)
plot(solved, 2)
plot(solved, 3)

# plot median history and estimate precision
plot(solved, 1, type = 'history')
plot(solved, 1, type = 'density')
plot(solved, 1, type = 'iterations')

# verify with true power from pwr package
library(pwr)

```

```

pwr.t.test(n=100, power = .8)
pwr.t.test(n=50, power = .8)
pwr.t.test(n=25, power = .8)

# use estimated d results to see how close power was
pwr.t.test(n=100, d = solved$d[1])
pwr.t.test(n=50, d = solved$d[2])
pwr.t.test(n=25, d = solved$d[3])

### failing analytic formula, confirm results with more precise
### simulation via runSimulation() (not required; if accuracy is important
### PROBABLI should longer)
# confirm <- runSimulation(design=solved, replications=10000, parallel=TRUE,
#                           generate=Generate, analyse=Analyse,
#                           summarise=Summarise)
# confirm

#-----

#####
## Criterion Analysis
#####

# GOAL: solve Type I error rate (alpha) given sample size, effect size, and
# power inputs (inputs for root no longer required to be an integer). Only useful
# when Type I error is less important than achieving the desired 1-beta (power)

Design <- createDesign(N = 50,
                      d = c(.2, .5, .8),
                      sig.level = NA)
Design # solve for NA's

# all other function definitions same as above

# search for alpha within [.0001, .8]
solved <- SimSolve(design=Design, b = .8, interval=c(.0001, .8),
                  generate=Generate, analyse=Analyse,
                  summarise=Summarise, integer=FALSE)

solved
summary(solved)
plot(solved, 1)
plot(solved, 2)
plot(solved, 3)

# plot median history and estimate precision
plot(solved, 1, type = 'history')
plot(solved, 1, type = 'density')
plot(solved, 1, type = 'iterations')

# verify with true power from pwr package
library(pwr)
pwr.t.test(n=50, power = .8, d = .2, sig.level=NULL)

```

```

pwr.t.test(n=50, power = .8, d = .5, sig.level=NULL)
pwr.t.test(n=50, power = .8, d = .8, sig.level=NULL)

# use estimated alpha results to see how close power was
pwr.t.test(n=50, d = .2, sig.level=solved$sig.level[1])
pwr.t.test(n=50, d = .5, sig.level=solved$sig.level[2])
pwr.t.test(n=50, d = .8, sig.level=solved$sig.level[3])

### failing analytic formula, confirm results with more precise
### simulation via runSimulation() (not required; if accuracy is important
### PROBABLI should just be run longer)
# confirm <- runSimulation(design=solved, replications=10000, parallel=TRUE,
#                           generate=Generate, analyse=Analyse,
#                           summarise=Summarise)
# confirm

## End(Not run)

```

---

|           |  |
|-----------|--|
| Summarise | <i>Summarise simulated data using various population comparison statistics</i> |
|-----------|--|

---

### Description

This collapses the simulation results within each condition to composite estimates such as RMSE, bias, Type I error rates, coverage rates, etc. See the See Also section below for useful functions to be used within Summarise.

### Usage

```
Summarise(condition, results, fixed_objects)
```

### Arguments

|               |   |
|---------------|---|
| condition     | a single row from the design input from <a href="#">runSimulation</a> (as a <code>data.frame</code> ), indicating the simulation conditions   |
| results       | a tibble data frame (if <code>Analyse</code> returned a named numeric vector of any length) or a list (if <code>Analyse</code> returned a list or multi-rowed <code>data.frame</code> ) containing the analysis results from <a href="#">Analyse</a> , where each cell is stored in a unique row/list element |
| fixed_objects | object passed down from <a href="#">runSimulation</a>   |

### Value

for best results should return a named numeric vector or `data.frame` with the desired meta-simulation results. Named list objects can also be returned, however the subsequent results must be extracted via [SimExtract](#)



## References

Chalmers, R. P., & Adkins, M. C. (2020). Writing Effective and Reliable Monte Carlo Simulations with the SimDesign Package. *The Quantitative Methods for Psychology*, 16(4), 248-280. doi:10.20982/tqmp.16.4.p248

Sigal, M. J., & Chalmers, R. P. (2016). Play it again: Teaching statistics with Monte Carlo simulation. *Journal of Statistics Education*, 24(3), 136-156. doi:10.1080/10691898.2016.1246953

## See Also

[bias](#), [RMSE](#), [RE](#), [EDR](#), [ECR](#), [MAE](#), [SimExtract](#)

## Examples

```
## Not run:

summarise <- function(condition, results, fixed_objects) {

  #find results of interest here (alpha < .1, .05, .01)
  lessthan.05 <- EDR(results, alpha = .05)

  # return the results that will be appended to the design input
  ret <- c(lessthan.05=lessthan.05)
  ret
}

## End(Not run)
```

---

timeFormater

*Format time string to suitable numeric output*

---

## Description

Format time input string into suitable numeric output metric (e.g., seconds). Input follows the SBATCH utility specifications. Accepted time formats include "minutes", "minutes:seconds", "hours:minutes:seconds", "days-hours", "days-hours:minutes" and "days-hours:minutes:seconds".

## Usage

```
timeFormater(time, output = "sec")
```

## Arguments

|        |  |
|--------|--|
| time   | a character string to be formatted. If a numeric vector is supplied then this will be interpreted as seconds.                          |
| output | type of numeric output to convert time into. Currently supported are 'sec' for seconds (default), 'min' for minutes, 'hour', and 'day' |

## Details

For example, `max_time = "60"` indicates a maximum time of 60 minutes, `max_time = "03:00:00"` a maximum time of 3 hours, `max_time = "4-12"` a maximum of 4 days and 12 hours, and `max_time = "2-02:30:00"` a maximum of 2 days, 2 hours and 30 minutes.

## Examples

```
# Test cases (outputs in seconds)
timeFormater("4-12")          # day-hours
timeFormater("4-12:15")      # day-hours:minutes
timeFormater("4-12:15:30")   # day-hours:minutes:seconds

timeFormater("30")           # minutes
timeFormater("30:30")        # minutes:seconds
timeFormater("4:30:30")      # hours:minutes:seconds

# output in hours
timeFormater("4-12", output = 'hour')
timeFormater("4-12:15", output = 'hour')
timeFormater("4-12:15:30", output = 'hour')

timeFormater("30", output = 'hour')
timeFormater("30:30", output = 'hour')
timeFormater("4:30:30", output = 'hour')
```

# Index

- \* **data**
  - BF\_sim, 11
  - BF\_sim\_alternative, 12
- \* **package**
  - SimDesign, 118
- abs, 14
- add\_missing(addMissing), 3
- addMissing, 3, 30
- aggregate\_simulations(SimCollect), 115
- Analyse, 5, 8, 29, 32, 88–90, 96, 98, 136
- AnalyseIf, 6, 7, 98
- aov, 112
- as.vector, 57
- Attach, 9, 30, 98
- attach, 9
  
- BF\_sim, 11, 12
- BF\_sim\_alternative, 11, 12
- bias, 13, 55, 72, 137
- boot\_predict(bootPredict), 15
- bootPredict, 15
- Bradley1978, 18, 27
- browser, 90
  
- c, 48, 49
- cat, 53
- CC, 20
- clusterSetRNGStream, 21, 92
- clusterSetRNGSubStream, 21
- colMeans, 22
- colSDs(colVars), 21
- colVars, 21
- commandArgs, 35
- cor, 20
- createDesign, 22, 23, 28, 29, 88, 96, 98, 128
  
- dunif, 51
  
- ECR, 18, 19, 24, 27, 112, 137
- EDR, 18, 19, 26, 27, 112, 137
  
- ERR (EDR), 27
- expand.grid, 22
- expandDesign, 23, 28, 81, 83
  
- family, 108
- future, 87, 91, 97
  
- gc, 95
- gen\_seeds, 92
- gen\_seeds(genSeeds), 33
- Generate, 5, 6, 10, 29, 88, 96, 98
- GenerateIf, 31, 88
- genSeeds, 33, 81–83
- getArrayID, 35, 82, 83
- getwd, 124
- glm, 108, 129
- grepl, 40, 42, 43
  
- integrate, 37
- IRMSE, 36
  
- library, 89
- lm, 16, 112
  
- MAE, 38, 137
- makeCluster, 91
- manageMessages, 39, 43
- manageWarnings, 6, 41, 42, 54, 93, 98
- message, 53
- MSRSE, 47
  
- nc, 48
- nextRNGStream, 83
- nextRNGSubStream, 82
  
- optimize, 61
  
- parallel, 87
- PBA, 50, 76, 77, 129–131
- plan, 91
- plot.PBA (PBA), 50

- plot.RM (RobbinsMonro), 76
- plot.SimSolve (SimSolve), 127
- predict, 108
- print.Design (createDesign), 22
- print.PBA (PBA), 50
- print.RM (RobbinsMonro), 76
- print.SFA (SFA), 107
- print.SimDesign (runSimulation), 87
- quiet, 39, 41, 43, 53
- RAB, 54
- rbind.SimDesign, 55
- RD, 57
- RE, 58, 137
- readRDS, 123
- rejectionSampling, 60
- require, 89
- reSummarise, 63, 88, 91, 98
- rHeadrick, 29, 30, 66
- rint, 68
- rinvWishart, 69
- rmgh, 29, 30, 70
- RMSD (RMSE), 71
- RMSE, 14, 37, 59, 71, 137
- rmvnorm, 73
- rmvt, 74
- RNG, 21
- RobbinsMonro, 52, 76
- RSE, 78
- rtruncate, 79
- runArraySimulation, 28, 29, 33, 35, 81, 87, 94, 95, 98, 113, 115, 116, 118
- runSimulation, 5–8, 10, 16, 23, 28–32, 55, 56, 64, 69, 74, 75, 81–83, 87, 95, 107, 108, 112, 114–116, 118, 119, 121, 123–129, 136
- rValeMaurelli, 29, 30, 104
- sample, 68
- sample.int, 68
- saveRDS, 97, 114
- Serlin2000, 19, 106
- set.seed, 92, 115
- SFA, 107, 131
- SimAnova, 96, 98, 111
- SimCheck, 83, 113, 116
- SimClean, 98, 114
- SimCollect, 28, 29, 83, 98, 113, 115
- SimDesign, 118
- SimDesign-package (SimDesign), 118
- SimExtract, 88, 90, 94, 97, 98, 116, 119, 136, 137
- SimFunctions, 96, 98, 121
- SimResults, 90, 91, 98, 123
- SimShiny, 98, 125
- SimSolve, 52, 108, 122, 127
- stop, 6, 30
- subset, 112
- Summarise, 79, 88, 96, 98, 119, 136
- summary.SimDesign (runSimulation), 87
- summary.SimSolve (SimSolve), 127
- suppressWarnings, 42, 43
- timeFormater, 51, 83, 95, 137
- try, 6, 30
- tryCatch, 43
- uniroot, 50, 52, 76, 77, 108, 129
- unname, 14, 18, 20, 21, 25, 27, 38, 47, 55, 57, 59, 72, 79
- Vectorize, 60