

# Package ‘RIbench’

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**Type** Package

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**Title** Benchmark Suite for Indirect Methods for RI Estimation

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**Depends** R (>= 3.3.0)

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RColorBrewer

**Suggests** knitr, rmarkdown

## Description

The provided benchmark suite enables the automated evaluation and comparison of any existing and novel indirect method for reference interval ('RI') estimation in a systematic way. Indirect methods take routine measurements of diagnostic tests, containing pathological and non-pathological samples as input and use sophisticated statistical methods to derive a model describing the distribution of the non-pathological samples, which can then be used to derive reference intervals. The benchmark suite contains 5,760 simulated test sets with varying difficulty.

To include any indirect method, a custom wrapper function needs to be provided.

The package offers functions for generating the test sets, executing the indirect method and evaluating the results.

See ?RIbench or vignette(``RIbench\_package'') for a more comprehensive description of the features.

A detailed description and application is described in Ammer T., Schuetzenmeister A., Prokosch H.-U., Zierk J., Rank C.M., Rauh M. ``RIbench: A Proposed Benchmark for the Standardized Evaluation of Indirect Methods for Reference Interval Estimation''. Clinical Chemistry (2022) <[doi:10.1093/clinchem/hvac142](https://doi.org/10.1093/clinchem/hvac142)>.

**License** GPL (>= 3)

**VignetteBuilder** knitr, rmarkdown

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<b>RIBench-package</b>	<i>RIBench: Benchmark Suite for the Standardized Evaluation of Indirect Methods for Reference Interval Estimation</i>
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## Description

RIBench enables the automated evaluation and comparison of any existing and novel indirect method in a systematic way. Indirect methods take routine measurements of diagnostic tests, containing pathological and non-pathological samples as input and use sophisticated statistical methods to derive a model describing the distribution of the non-pathological samples, which can then be used to derive reference intervals. The benchmark suite contains 5,760 simulated data sets with varying difficulty. To include any indirect method, a custom wrapper function needs to be provided. The package offers functions for generating the test sets `generateBiomarkerTestSets`, executing the indirect method `evaluateBiomarkerTestSets` and evaluating the results `evaluateAlgorithmResults`.

## Details

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Type:	Package
Version:	1.0.2
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## Author(s)

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## References

Ammert, T., Schuetzenmeister, A., Prokosch, HU., Zierk, J., Rank, C.M., Rauh, M. RIBench: A Proposed Benchmark for the Standardized Evaluation of Indirect Methods for Reference Interval Estimation. *Clin Chem* (2022) [Accepted, July 12].

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addGrid	<i>Add a grid to an existing plot.</i>
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**Description**

It is possible to use automatically determined grid lines ( $x=NULL$ ,  $y=NULL$ ) or specifying the number of cells  $x = 3$ ,  $y = 4$  as done by `grid`. Additionally, x- and y-locations of grid-lines can be specified, e.g.  $x = 1:10$ ,  $y = seq(0,10,2)$ .

**Usage**

```
addGrid(x = NULL, y = NULL, col = "lightgray", lwd = 1L, lty = 3L)
```

**Arguments**

<code>x</code>	(integer, numeric) single integer specifies number of cells, numeric vector specifies vertical grid-lines
<code>y</code>	(integer, numeric) single integer specifies number of cells, numeric vector specifies horizontal grid-lines
<code>col</code>	(character) color of grid-lines
<code>lwd</code>	(integer) line width of grid-lines
<code>lty</code>	(integer) line type of grid-lines

**Author(s)**

Andre Schuetzenmeister <[andre.schuetzenmeister@roche.com](mailto:andre.schuetzenmeister@roche.com)>

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as.rgb	<i>Convert color-names or RGB-code to possibly semi-transparent RGB-code.</i>
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**Description**

Function takes the name of a color and converts it into the rgb space. Parameter "alpha" allows to specify the transparency within [0,1], 0 meaning completely transparent and 1 meaning completely opaque. If an RGB-code is provided and alpha != 1, the RGB-code of the transparency adapted color will be returned.

**Usage**

```
as.rgb(col = "black", alpha = 1)
```

**Arguments**

- col (character) name of the color to be converted/transformed into RGB-space (code). Only those colors can be used which are part of the set returned by function colors(). Defaults to "black".
- alpha (numeric) value specifying the transparency to be used, 0 = completely transparent, 1 = opaque.

**Value**

RGB-code

**Author(s)**

Andre Schuetzenmeister <andre.schuetzenmeister@roche.com>

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BoxCox

*One-parameter Box-Cox transformation.*

---

**Description**

One-parameter Box-Cox transformation.

**Usage**

BoxCox(x, lambda)

**Arguments**

- x (numeric) data to be transformed
- lambda (numeric) Box-Cox transformation parameter

**Value**

(numeric) vector with Box-Cox transformation of x

**Author(s)**

Andre Schuetzenmeister <andre.schuetzenmeister@roche.com>

<code>computeDirect</code>	<i>Function to simulate the direct method</i>
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## Description

Function to simulate the direct method

## Usage

```
computeDirect(
  N = 120,
  analyte,
  params,
  seed = 123,
  NIter = 10000,
  RIperc = c(0.025, 0.975)
)
```

## Arguments

<code>N</code>	(integer) specifying the number of samples used as sample size for the direct method, default: 120
<code>analyte</code>	(character) specifying the biomarker that is currently simulated
<code>params</code>	(list) of parameters for non-pathological distribution (nonp_mu, nonp_sigma, nonp_lambda, and nonp_shift)
<code>seed</code>	(integer) specifying the seed used for the simulation, default: 123
<code>NIter</code>	(integer) specifying the number of times N samples should be drawn out of the simulated non-pathological distribution (default: 10,000)
<code>RIperc</code>	(numeric) value specifying the percentiles, which define the reference interval

## Value

(data frame) with the estimated reference intervals for NIter iterations

## Author(s)

Tatjana Ammer <tatjana.ammer@roche.com>

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**computePerfMeas***Function for computing performance measurements*

---

## Description

Function for computing performance measurements

## Usage

```
computePerfMeas(  
  analyte,  
  algo,  
  resRIs,  
  subTable,  
  RIperc = c(0.025, 0.975),  
  cutoffZ = 5  
)
```

## Arguments

analyte	(character) specifying current analyzed analyte
algo	(character) specifying used algorithm
resRIs	(data.frame) with all calculated reference intervals
subTable	(data.frame) containing all information about the simulated test sets
RIperc	(numeric) vector specifying the percentiles for the reference interval, default: 0.025 and 0.975
cutoffZ	(numeric) specifying if a cutoff should be used to classify results as implausible and exclude from analysis

## Value

updated data frame with computed performance measures

## Author(s)

Tatjana Ammer <tatjana.ammer@roche.com>

**computePerfMeasAll**      *Function for computing reference intervals for all markers*

### Description

Function for computing reference intervals for all markers

### Usage

```
computePerfMeasAll(analytes, algo, risIn, tableTCs, cutoffZ = 5)
```

### Arguments

analytes	(character) listing all analytes for which the result files should be parsed
algo	(character) specifying used algorithm
risIn	(list) with data frame of all calculated reference intervals
tableTCs	(data.frame) containing all information about the simulated test sets
cutoffZ	(integer) specifying if and if so which cutoff should be used to classify results as implausible (default: 5)

### Value

list with the calculated errors as data frame for each marker

### Author(s)

Tatjana Ammer <tatjana.ammer@roche.com>

**computeRIs**      *Function for computing reference intervals*

### Description

Function for computing reference intervals

### Usage

```
computeRIs(
  analyte,
  algo,
  results,
  tableTCs,
  RIperc = c(0.025, 0.975),
  truncNormal = FALSE
)
```

**Arguments**

analyte	(character) specifying analyte
algo	(character) specifying used algorithm
results	(list) with all calculated results as RWDRI objects
tableTCs	(data frame) containing all information about the simulated test sets
RIperc	(numeric) vector specifying the percentiles for the reference interval, default: 0.025 and 0.975
truncNormal	(logical) specifying if a normal distribution truncated at zero shall be assumed

**Value**

data frame with computed reference intervals

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

---

computeRIsAll

*Function for computing reference intervals for all markers*

---

**Description**

Function for computing reference intervals for all markers

**Usage**

```
computeRIsAll(analytes, algo, resIn, tableTCs, truncNormal = FALSE)
```

**Arguments**

analytes	(character) listing all markers for which the result files should be parsed
algo	(character) specifying used algorithm
resIn	(list) with all calculated results for all markers as RWDRI objects
tableTCs	(data.frame) containing all information about the simulated test sets
truncNormal	(logical) specifying if a normal distribution truncated at zero shall be assumed

**Value**

list with the calculated reference intervals as data frame for each marker

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

`computeRuntimeAll`      *Function to compute runtime statistics for all analytes*

### Description

Function to compute runtime statistics for all analytes

### Usage

```
computeRuntimeAll(analytes, algo, risIn, tableTCs)
```

### Arguments

<code>analytes</code>	(character) listing all analytes for which the result files should be parsed
<code>algo</code>	(character) specifying used algorithm
<code>risIn</code>	(list) with data frame of all calculated reference intervals and runtime
<code>tableTCs</code>	(data.frame) containing all information about the simulated test cases

### Value

(list) with runtime statistics per analyte and data frames with raw runtime overall and per analyte

### Author(s)

Tatjana Ammer <[tatjana.ammer@roche.com](mailto:tatjana.ammer@roche.com)>

`computeSubResults`      *Helper function to compute the subscores for the distribution types and the mentioned categories*

### Description

Helper function to compute the subscores for the distribution types and the mentioned categories

### Usage

```
computeSubResults(
  errorDf,
  tableTCs,
  distCat,
  errorParam,
  catList,
  catLabels,
  perfCombination = "mean"
)
```

**Arguments**

errorDf	(data frame) containing the estimate reference intervals and all computed error measures
tableTCs	(data.frame) containing all information about the simulated test sets
distCat	(character) specifying the distribution category
errorParam	(character) specifying for which error parameter the data frame should be generated
catList	(character) vector containing the categories to split the dataset
catLabels	(character) vector containing the labels that will be used for the categories
perfCombination	(character) specifying if mean (default), median or sum should be computed

**Value**

(data frame) containing the computed subscores

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

defineSubset

*Function for defining a subset that is used for analyzing the computation time and can be used for other subanalyses.*

**Description**

Function for defining a subset that is used for analyzing the computation time and can be used for other subanalyses.

**Usage**

```
defineSubset(tableTCs = NULL, N = 50, seed = 123)
```

**Arguments**

tableTCs	(data frame) describing the pre-defined testcases
N	(integer) describing the number of testcases per biomarker contained in the subset (default: 50)
seed	(integer) specifying the seed used for defining the subset, default: 123

**Value**

(data frame) describing the updated table with all test case definitions.

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

---

**evaluateAlgorithmResults**

*Convenience Function to generate all result plots and calculate the benchmark score*

---

**Description**

Convenience Function to generate all result plots and calculate the benchmark score

**Usage**

```
evaluateAlgorithmResults(
  workingDir = "",
  algoNames = NULL,
  subset = "all",
  evalFolder = "Evaluation",
  withDirect = TRUE,
  withMean = TRUE,
  outline = TRUE,
  errorParam = c("zzDevAbs_Ov", "AbsPercError_Ov", "AbsError_Ov"),
  cutoffZ = 5,
  cols = NULL,
  ...
)
```

**Arguments**

workingDir	(character) specifying the working directory: Plots will be stored in 'workingDir/evalFolder' and results will be used from 'workingDir/Results/algoName/biomarker'
algoNames	(character) vector specifying all algorithms that should be part of the evaluation
subset	(character, numeric, or data.frame) to specify for which subset the algorithm should be evaluated. character options: 'all' (default) for all test sets, a distribution type: 'normal', 'skewed', 'heavilySkewed', 'shifted'; a biomarker: 'Hb', 'Ca', 'FT4', 'AST', 'LACT', 'GGT', 'TSH', 'IgE', 'CRP', 'LDH'; 'Runtime' for runtime analysis subset; numeric option: number of test sets per biomarker, e.g. 10; data.frame: customized subset of table with test set specifications
evalFolder	(character) specifying the name of the ouput directory, Plots will be stored in workingDir/evalFolder, default: 'Evaluation'
withDirect	(logical) indicating whether the direct method should be simulated for comparison (default:TRUE)
withMean	(logical) indicating whether the mean should be plotted as well (default: TRUE)
outline	(logical) indicating whether outliers should be drawn (TRUE, default), or not (FALSE)
errorParam	(character) specifying for which error parameter the data frame should be generated, choose between absolute z-score deviation ("zzDevAbs_Ov"), absolute percentage error ("AbsPercError_Ov"), and absolute error ("AbsError_Ov")

cutoffZ	(integer) specifying if and so which cutoff for the absolute z-score deviation should be used to classify results as implausible and exclude them from the overall benchmark score (default: 5)
cols	(character) vector specifying the colors used for the different algorithms
...	additional arguments to be passed to the method, e.g. "truncNormal" (logical) vector specifying if a normal distribution truncated at zero shall be assumed, can be either TRUE/FALSE or a vector with TRUE/FALSE for each algorithm; "coldDirect" (character) specifying the color used for the direct method, default: "grey" "ylab" (character) specifying the label for the y-axis

**Value**

(data frame) containing the computed benchmark results

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

**Examples**

```
## Not run:
# Ensure that 'generateBiomarkerTestSets()' and 'evaluateBiomarkerTestSets()' is called
# with the same workingDir and for all mentioned algorithms before calling this function.

# first example, evaluation for several algorithms
benchmarkScore <- evaluateAlgorithmResults(workingDir=tempdir(),
algoNames=c("Hoffmann", "TML", "kosmic", "TMC", "refineR"))
# The function will create several plots saved in workingDir/Evaluation.

# second example, evaluation for only one algorithm and a defined subset
benchmarkScore <- evaluateAlgorithmResults(workingDir = tempdir(),
algoNames = "refineR", subset = 'Ca')

# third example, saving the results in a different folder, and setting a different cutoff
# for the absolute z-score deviation
benchmarkScore <- evaluateAlgorithmResults(workingDir = tempdir(), algoNames = "refineR",
subset = 'Ca', cutoffZ = 4, evalFolder = "Eval_Test")

## End(Not run)
```

**evaluateBiomarkerTestSets**

*Wrapper function to evaluate all test sets or a specified subset for a specified algorithm.*

## Description

Wrapper function to evaluate all test sets or a specified subset for a specified algorithm.

## Usage

```
evaluateBiomarkerTestSets(
  workingDir = "",
  algoName = "refineR",
  algoFunction = "findRI",
  libs = "refineR",
  sourceFiles = NULL,
  params = NULL,
  requireDecimals = FALSE,
  requirePercentiles = FALSE,
  subset = "all",
  timeLimit = 14400,
  verbose = TRUE,
  showWarnings = FALSE,
  ...
)
```

## Arguments

workingDir	(character) specifying the working directory: Results will be stored in 'workingDir/Results/algo/biomarker' and data will be used from 'workingDir/Data/biomarker'
algoName	(character) specifying the name of the algorithm that is evaluated
algoFunction	(character) specifying the name of the function needed for estimating RIs
libs	(list) containing all libraries needed for executing the algorithm
sourceFiles	(list) containing all source files needed for executing the algorithm
params	(list) with additional parameters needed for calling algoFunction
requireDecimals	(logical) indicating whether the algorithm needs the number of decimal places (TRUE) or not (FALSE, default)
requirePercentiles	(logical) indicating whether only percentiles and no model is estimated
subset	(character, numeric, or data.frame) to specify for which subset the algorithm should be executed. character options: 'all' (default) for all test sets; a distribution type: 'normal', 'skewed', 'heavilySkewed', 'shifted'; a biomarker: 'Hb', 'Ca', 'FT4', 'AST', 'LACT', 'GGT', 'TSH', 'IgE', 'CRP', 'LDH'; 'Runtime' for runtime analysis subset; numeric option: number of test sets per biomarker, e.g. 10; data.frame: customized subset of table with test set specifications
timeLimit	(integer) specifying the maximum amount of time in seconds allowed to execute one single estimation (default: 14400 sec (4h))
verbose	(logical) indicating if the progress counter should be shown (default: TRUE)
showWarnings	(logical) indicating whether warnings from the call to the indirect method/algorithm should be shown (default: FALSE)

... additional arguments to be passed to the method, e.g. specified in- and output directory ('inputDir', 'outputDir')

### Value

(data frame) containing information about the test sets where the algorithm terminated the R session or failed to report a result

### Author(s)

Tatjana Ammer <tatjana.ammer@roche.com>

### Examples

```
## Not run:
# The evaluation of all test sets can take several hours depending on
#   the computation time of the algorithm.
# Wrapper function for indirect method required, see vignette("RIBench_package")
# Ensure that 'generateBiomarkerTestSets()' is called with the same workingDir
#   before calling this function.

# first generic example
evaluateBiomarkerTestSets(workingDir = tempdir(), algoName = 'myOwnAlgo',
  algoFunction = 'estimateModel', libs = c('myOwnAlgo'),
  sourceFiles = list("C:\\\\Temp\\\\MyAlgoWrapper.R"),
  requireDecimals = FALSE, requirePercentiles = FALSE,
  subset = 'all', timeLimit = 14400)

# second example, evaluation for only 'Calcium' test sets.
progress <- evaluateBiomarkerTestSets(workingDir = tempdir(), algoName = 'myOwnAlgo',
  algoFunction = 'estimateModel', libs = c('myOwnAlgo'), subset = "Ca")

# third example, evaluation for only a subset testsets that follow a skewed distribution.
progress <- evaluateBiomarkerTestSets(workingDir = tempdir(), algoName = 'myOwnAlgo',
  algoFunction = 'estimateModel', libs = c('myOwnAlgo'), subset = "skewed")

# forth example, evaluation for a subset of 3 testsets per biomarker.
progress <- evaluateBiomarkerTestSets(workingDir = tempdir(), algoName = 'myOwnAlgo',
  algoFunction = 'estimateModel', libs = c('myOwnAlgo'), subset = 3)

# fifth example, evaluation for a customized subset with all test sets that have
#   a pathological fraction <= 30%.
testsets <- loadTestsetDefinition()
progress <- evaluateBiomarkerTestSets(workingDir = tempdir(), algoName = 'myOwnAlgo',
  algoFunction = 'estimateModel', libs = c('myOwnAlgo'),
  subset = testsets[testsets$fractionPathol <= 0.3, ] )
```

```

# sixth example, evaluation forwarding additional parameters to the 'algoFunction'
progress <- evaluateBiomarkerTestSets(workingDir = tempdir(), algoName = 'myOwnAlgo',
algoFunction = 'estimateModel', libs = c('myOwnAlgo'),
sourceFiles = list("Test_RIEst_2pBoxCox"), params = list("model='2pBoxCox'"))

# seventh example, evaluation for indirect method that requires the number of
# decimal points as input
evaluateBiomarkerTestSets(workingDir = tempdir(), algoName = 'myOwnAlgo',
algoFunction = 'estimateModelDec', libs = c('myOwnAlgo'),
sourceFiles = "C:\\\\Temp\\\\Test_RIEst_dec.R", requireDecimals = TRUE)

# eigth example, evaluation for indirect method that directly estimates the percentiles
evaluateBiomarkerTestSets(workingDir = tempdir(), algoName="myOwnAlgo",
algoFunction="estimateRIs", libs="myOwnAlgo",
sourceFiles = "C:\\\\Temp\\\\Test_RIEst.R", requirePercentiles=TRUE)

## End(Not run)

```

**formatNumber***Rounding method with trailing zeros.***Description**

Rounding method with trailing zeros.

**Usage**

```
formatNumber(x, digits)
```

**Arguments**

- |        |  |
|--------|--|
| x      | (numeric) value that is rounded                              |
| digits | (integer) indicating the number of decimal places to be used |

**Value**

Rounded value with trailing zeros

**Author(s)**

Christopher Rank <[christopher.rank@roche.com](mailto:christopher.rank@roche.com)>

---

**generateBiomarkerTestSets**

*Convenience function to generate simulated data and save each test set as a separate file*

---

**Description**

Convenience function to generate simulated data and save each test set as a separate file

**Usage**

```
generateBiomarkerTestSets(  
  workingDir = "",  
  subset = "all",  
  rounding = TRUE,  
  verbose = TRUE  
)
```

**Arguments**

workingDir	(character) specifying the working directory where 'workingDir/Data/biomarker' folders will be generated containing the simulated data
subset	(character, numeric, or data.frame) to specify for which subset the data should be generated and the algorithms later applied to. character options: 'all' (default) for all test sets; a distribution type: 'normal', 'skewed', 'heavilySkewed', 'shifted'; a biomarker: 'Hb', 'Ca', 'FT4', 'AST', 'LACT', 'GGT', 'TSH', 'IgE', 'CRP', 'LDH'; 'Runtime' for runtime analysis subset; numeric option: number of test sets per biomarker, e.g. 10; data.frame: customized subset of table with test set specification
rounding	(logical) indicating whether decimal places stated in test set specification should be applied (default, TRUE), if FALSE, data will be rounded to 5 decimal places to mimic unrounded data
verbose	(logical) indicating if the progress counter should be shown (default: TRUE)

**Value**

No return value, instead the data files are generated and saved in the workingDir

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

## Examples

```
## Not run:
workingDir <- "C:\\Temp\\RIbench\\"
generateBiomarkerTestSets(workingDir = workingDir)

## End(Not run)

# example generating a subset of 2 test sets per biomarker
generateBiomarkerTestSets(workingDir = tempdir(), subset = 2)
```

## generateBoxPlotOneAnalyte

*Wrapper function to generate one boxplot for a specified analyte*

## Description

Wrapper function to generate one boxplot for a specified analyte

## Usage

```
generateBoxPlotOneAnalyte(
  errorListAll,
  colList,
  nameList,
  catList,
  catLabels,
  a,
  errorParam,
  outline = TRUE,
  withMean = TRUE,
  withCats = TRUE,
  withDirect = TRUE,
  titlePart = NULL,
  outputDir,
  filenamePart = NULL,
  ylim1 = c(0, 100),
  ylim2 = c(100, 1000),
  ...
)
```

## Arguments

errorListAll (list) containing the overall benchmark results per algorithm

colList	(character) vector specifying the colors used for the different algorithms (should correspond to columns of benchmark results)
nameList	(character) vector specifying the names used in the legend (should correspond to columns of benchmark results), if NULL, colnames will be used
catList	(character) vector specifying the categories for which the boxes should be drawn
catLabels	(character) vector specifying the labels to the associated categories used for the x-axis
a	(character) specifying the analyte for which the boxplot should be generated
errorParam	(character) specifying for which error measure the plot should be generated
outline	(logical) indicating whether outliers should be drawn (TRUE, default), or not (FALSE)
withMean	(logical) indicating whether the mean should be plotted as well (default: TRUE)
withCats	(logical) set to TRUE if categories (e.g. pathological fraction) should be plotted (default: FALSE)
withDirect	(logical) indicating whether the box of the direct method should be elongated to facilitate comparison (default: TRUE)
titlePart	(character) specifying the latter part of the title
outputDir	(character) specifying a output directory
filenamePart	(character) specifying a filename for the plot
ylim1	(numeric) vector specifying the limits in y-direction for the first granular scale
ylim2	(numeric) vector specifying the limits in y-direction for the second less detailed scale
...	additional arguments passed forward to other functions

**Value**

No return value. Instead, a plot is generated.

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

generateBoxplotsDistTypes

*Wrapper function to generate all boxplots for the specified distribution types split by defined categories*

**Description**

Wrapper function to generate all boxplots for the specified distribution types split by defined categories

**Usage**

```
generateBoxplotsDistTypes(
  errorListAll,
  colList,
  nameList,
  catList,
  catLabels,
  errorParam = "zzDevAbs_0v",
  outline = TRUE,
  withMean = TRUE,
  withDirect = TRUE,
  withCats = TRUE,
  titlePart = NULL,
  outputDir = NULL,
  filenamePart = NULL,
  ylim1Vec = NULL,
  ylim2Vec = NULL,
  yticks1Vec = NULL,
  yticks2Vec = NULL,
  ...
)
```

**Arguments**

errorListAll	(list) containing the overall benchmark results per algorithm
colList	(character) vector specifying the colors used for the different algorithms (should correspond to columns of benchmark results)
nameList	(character) vector specifying the names used in the legend (should correspond to columns of benchmark results), if NULL, colnames will be used
catList	(character) vector specifying the categories for which the boxes should be drawn
catLabels	(character) vector specifying the labels to the associated categories used for the x-axis
errorParam	(character) specifying for which error measure the plot should be generated
outline	(logical) indicating whether outliers should be drawn (TRUE, default), or not (FALSE)
withMean	(logical) indicating whether the mean should be plotted as well (default: TRUE)
withDirect	(logical) indicating whether the box of the direct method should be elongated to facilitate comparison (default:TRUE)
withCats	(logical) set to TRUE if categories (e.g. pathological fraction) should be plotted (default: FALSE)
titlePart	(character) specifying the latter part of the title
outputDir	(character) specifying a output directory
filenamePart	(character) specifying a filename for the plot
ylim1Vec	(numeric) vector specifying the limits in y-direction for the first granular scale

ylim2Vec	(numeric) vector specifying the limits in y-direction for the second less detailed scale
yTicks1Vec	(numeric) vector specifying the ticks in y-direction for the first granular scale
yTicks2Vec	(numeric) vector specifying the ticks in y-direction for the second less detailed scale
...	additional arguments passed forward to other functions

**Value**

No return value. Instead, a plot is generated.

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

**generateBoxplotsMultipleCats**

*Wrapper function to generate all boxplots for the specified analytes split by defined categories*

**Description**

Wrapper function to generate all boxplots for the specified analytes split by defined categories

**Usage**

```
generateBoxplotsMultipleCats(
  analytes,
  errorListAll,
  colList,
  nameList,
  category = c("fractionPathol", "fractionPathol_cum", "N", "N_cum", "OvFreq",
              "OvFreq_cum"),
  catList = NULL,
  catLabels = NULL,
  errorParam = "zzDevAbs_Ov",
  outline = TRUE,
  withMean = TRUE,
  withDirect = TRUE,
  withCats = TRUE,
  titlePart = NULL,
  outputDir = NULL,
  filenamePart = NULL,
  ylim1Vec = NULL,
  ylim2Vec = NULL,
  yticks1Vec = NULL,
```

```

  yticks2Vec = NULL,
  ...
)

```

## Arguments

analytes	(character) vector specifying for which analytes the plots should be generated
errorListAll	(named list) containing the overall benchmark results per algorithm (names of list elements should be the names of the algorithms)
colList	(character) vector specifying the colors used for the different algorithms (should correspond to columns of benchmark results)
nameList	(character) vector specifying the names used in the legend (should correspond to columns of benchmark results), if NULL, colnames will be used
category	(character) defining the category used for creating the subsets. All defined sub-features are used for the categorization. Choose from "fractionPathol" (default), "N", or "OvFreq", individual or cumulative ("_cum"); if category is set this will be used to define catList and catLabels
catList	(character) vector specifying the categories for which the boxes should be drawn
catLabels	(character) vector specifying the labels to the associated categories used for the x-axis
errorParam	(character) specifying for which error measure the plot should be generated
outline	(logical) indicating whether outliers should be drawn (TRUE, default), or not (FALSE)
withMean	(logical) indicating whether the mean should be plotted as well (default: TRUE)
withDirect	(logical) indicating whether the box of the direct method should be elongated to facilitate comparison (default:TRUE)
withCats	(logical) set to TRUE if categories (e.g. pathological fraction) should be plotted (default: FALSE)
titlePart	(character) specifying the latter part of the title
outputDir	(character) specifying an output directory
filenamePart	(character) specifying a filename for the plot
ylim1Vec	(numeric) vector specifying the limits in y-direction for the first granular scale
ylim2Vec	(numeric) vector specifying the limits in y-direction for the second less detailed scale
yticks1Vec	(numeric) vector specifying the ticks in y-direction for the first granular scale
yticks2Vec	(numeric) vector specifying the ticks in y-direction for the second less detailed scale
...	additional arguments passed forward to other functions

## Value

No return value. Instead, a plot is generated.

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

---

generateDataFiles	<i>Generate simulated data with one start seed for each biomarker and save each test set as separate file</i>
-------------------	---

---

**Description**

Generate simulated data with one start seed for each biomarker and save each test set as separate file

**Usage**

```
generateDataFiles(  
  tableTCs = NULL,  
  outputDir = NULL,  
  rounding = TRUE,  
  verbose = TRUE  
)
```

**Arguments**

tableTCs	(data.frame) containing all information about the simulated test cases
outputDir	(character) specifying the output directory where the data files should be written to
rounding	(logical) indicating whether decimal places stated in tableTCs should be applied (default, TRUE), if FALSE, data will be rounded to 5 decimal places to mimic unrounded data
verbose	(logical) indicating if the progress counter should be shown (default: TRUE)

**Value**

No return value, instead the data files are generated

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

---

generateMD5

*Generate an MD5 hash sum for any R object.*

---

### Description

Generate an MD5 hash sum for any R object.

### Usage

```
generateMD5(x)
```

### Arguments

x	(object) any R object.
---	------------------------

### Value

(character) MD5 hash sum of the input object.

### Author(s)

Christopher Rank <christopher.rank@roche.com>

---

generateScatterplotsAll

*Wrapper function to generate scatterplots for the specified analytes*

---

### Description

Wrapper function to generate scatterplots for the specified analytes

### Usage

```
generateScatterplotsAll(
  analytes,
  errorListAll,
  colList = NULL,
  nameList,
  tableTCs,
  errorParam = "zzDevAbs",
  withColorCat = NULL,
  titlePart = NULL,
  outputDir = NULL,
  filenamePart = NULL,
  ylim = NULL,
  xlim = NULL,
```

```
  xlab = NULL,  
  ylab = NULL,  
  ...  
)
```

## Arguments

analytes	(character) vector specifying for which analytes the scatterplot should be generated
errorListAll	(list) containing the overall benchmark results per algorithm
colList	(character) vector specifying the colors used for the different algorithms (should correspond to columns of benchmark results)
nameList	(character) vector specifying the names used in the legend (should correspond to columns of benchmark results), if NULL, colnames will be used
tableTCs	(data frame) containing all test case information
errorParam	(character) specifying for which error measure the plot should be generated
withColorCat	(character) indicating if plot should be colored according to the pathological fraction ("fractionPathol"), sample size ("N"), or "overlapPatholLeft", "overlapPatholRight"
titlePart	(character) specifying the latter part of the title
outputDir	(character) specifying a output directory
filenamePart	(character) specifying a filename for the plot
ylim	(numeric) vector specifying the limits in y-direction for the first granular scale
xlim	(numeric) vector specifying the limits in y-direction for the second less detailed scale
xlab	(character) specifying x-axis label
ylab	(character) specifying y-axis label
...	additional arguments passed forward to other functions

## Value

No return value. Instead, a plot is generated.

## Author(s)

Tatjana Ammer <tatjana.ammer@roche.com>

`getBenchmarkResults`    *Computing benchmark table with the mean overall results.*

## Description

Computing benchmark table with the mean overall results.

## Usage

```
getBenchmarkResults(
  errorList,
  nameVec,
  tableTCs,
  errorParam = "zzDevAbsCutoff_0v",
  cutoffZ = 5,
  catList = c("fractionPathol <= 0.20 & N <= 5000",
             "fractionPathol <= 0.20 & N > 5000", "fractionPathol > 0.20 & N <= 5000",
             "fractionPathol > 0.20 & N > 5000"),
  catLabels = c("lowPlowN", "lowPhighN", "highPlowN", "highPhighN"),
  perfCombination = c("mean", "median", "sum")
)
```

## Arguments

<code>errorList</code>	(list) containing the the computed errors for the different (indirect) methods/algorithms
<code>nameVec</code>	(character) vector specifying the names of the different (indirect) methods/algorithms
<code>tableTCs</code>	(data.frame) containing all information about the simulated test sets
<code>errorParam</code>	(character) specifying for which error parameter the data frame should be generated
<code>cutoffZ</code>	(integer) specifying if and if so which cutoff for the absolute z-score deviation should be used to classify results as implausible and exclude them from the overall benchmark score (default: 5)
<code>catList</code>	(character) vector containing the categories to split the dataset
<code>catLabels</code>	(character) vector containing the labels that will be used for the categories
<code>perfCombination</code>	(character) specifying which measure should be used to compute the overall benchmark score; choose from "mean" (default), "median", or "sum"

## Value

(data frame) containing the computed benchmark results

## Author(s)

Tatjana Ammer <[tatjana.ammer@roche.com](mailto:tatjana.ammer@roche.com)>

---

getRI	<i>Method to calculate reference intervals (percentiles) for objects of class 'RWDRI'</i>
-------	---

---

## Description

Method to calculate reference intervals (percentiles) for objects of class 'RWDRI'

## Usage

```
getRI(  
  x,  
  RIperc = c(0.025, 0.975),  
  CIprop = 0.95,  
  pointEst = c("fullDataEst", "medianBS", "meanBS"),  
  truncNormal = FALSE,  
  Scale = c("original", "transformed")  
)
```

## Arguments

x	(object) of class 'RWDRI'
RIperc	(numeric) value specifying the percentiles, which define the reference interval
CIprop	(numeric) value specifying the central region for estimation of confidence intervals
pointEst	(character) specifying the point estimate determination: (1) using the full dataset ("fullDataEst"), (2) calculating the median from all bootstrap samples ("medianBS"), (2) works only if NBootstrap > 0 (3) calculating the mean from all bootstrap samples ("meanBS"), (3) works only if NBootstrap > 0
truncNormal	(logical) specifying if a normal distribution truncated at zero shall be assumed
Scale	(character) specifying if percentiles are calculated on the original scale ("Or") or the transformed scale ("Tr")

## Value

(data.frame) with columns for percentile, point estimate and confidence intervals.

## Author(s)

Christopher Rank <[christopher.rank@roche.com](mailto:christopher.rank@roche.com)>, Tatjana Ammer <[tatjana.ammer@roche.com](mailto:tatjana.ammer@roche.com)>

`getRIsAllwithoutModel` *Function for retrieving reference intervals if directly computed*

### Description

Function for retrieving reference intervals if directly computed

### Usage

```
getRIsAllwithoutModel(analytes, algo, resIn, tableTCs)
```

### Arguments

<code>analytes</code>	(character) listing all markers for which the result files should be parsed
<code>algo</code>	(character) specifying used algorithm
<code>resIn</code>	(list) with all calculated results for all markers as RWDRI objects
<code>tableTCs</code>	(data.frame) containing all information about the simulated test sets

### Value

list with the calculated reference intervals as data frame for each marker

### Author(s)

Tatjana Ammer <[tatjana.ammer@roche.com](mailto:tatjana.ammer@roche.com)>

`getRuntime` *Helper function to compute runtime statistics*

### Description

Helper function to compute runtime statistics

### Usage

```
getRuntime(x, analyte)
```

### Arguments

<code>x</code>	(data.frame) with one column specifying the Runtime
<code>analyte</code>	(character) specifying current analyzed marker

### Value

(data.frame) containing runtime statistics (min, mean, median, max)

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

---

getSubset

*Function to group the data according to a specified feature.*

---

**Description**

The feature can either be the pathological fraction, the sample size or the overlap (category) individually or cumulative (\_cum). For an individualized categorisation see getSubsetForDefinedCats.

**Usage**

```
getSubset(  
  subsetDef,  
  distType = FALSE,  
  tableTCs,  
  errorList,  
  category = c("fractionPathol", "fractionPathol_cum", "N", "N_cum", "OvFreq",  
             "OvFreq_cum"),  
  restrict = NULL  
)
```

**Arguments**

subsetDef	(character) listing either the analytes or distribution types for which the result files should be parsed
distType	(logical) indicating if parameter subsetDef refers to analytes (FALSE, default) or distribution types (TRUE)
tableTCs	(data.frame) containing all information about the simulated test sets
errorList	(list) containing for each method the table with the computed error measurements
category	(character) defining the category used for creating the subsets. All defined sub-features are used for the categorization. Choose from "fractionPathol" (default), "N", or "OvFreq", individual or cumulative ("_cum")
restrict	(character) indicating whether test sets should be filtered according to specified restriction, default NULL, e.g. fractionPathol <= 0.30

**Value**

(list) containing the performance measurements grouped according to specified subset definition and categories

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

---

`getSubsetForDefinedCats`

*Function to group the data according to a specified feature.*

---

## Description

Function to group the data according to a specified feature.

## Usage

```
getSubsetForDefinedCats(
  subsetDef,
  distType = FALSE,
  tableTCs,
  errorList,
  catList = NULL,
  catLabels = NULL,
  restrict = NULL
)
```

## Arguments

<code>subsetDef</code>	(character) listing either the analytes or distribution types for which the result files should be parsed
<code>distType</code>	(logical) indicating if 'subsetDef' refers to analytes (FALSE, default) or distribution types (TRUE)
<code>tableTCs</code>	(data.frame) containing all information about the simulated test sets
<code>errorList</code>	(list) containing the table with the computed error measurements
<code>catList</code>	(list) containing the categories to split the dataset
<code>catLabels</code>	(list) containing the labels that will be used for the categories
<code>restrict</code>	(character) indicating whether testcases should be filtered according to specified restriction, default NULL, e.g. fractionPathol <= 0.30

## Value

(list) containing the performance measurements grouped according to specified subset definition and categories

## Author(s)

Tatjana Ammer <tatjana.ammer@roche.com>

---

`invBoxCox`

*Inverse of the one-parameter Box-Cox transformation.*

---

## Description

Inverse of the one-parameter Box-Cox transformation.

## Usage

```
invBoxCox(x, lambda)
```

## Arguments

<code>x</code>	(numeric) data to be transformed
<code>lambda</code>	(numeric) Box-Cox transformation parameter

## Value

(numeric) vector with inverse Box-Cox transformation of `x`

## Author(s)

Andre Schuetzenmeister <[andre.schuetzenmeister@roche.com](mailto:andre.schuetzenmeister@roche.com)>

---

---

`loadTestsetDefinition` *Convenience function to load the table with the information about the pre-defined test sets*

---

## Description

Convenience function to load the table with the information about the pre-defined test sets

## Usage

```
loadTestsetDefinition()
```

## Value

(data frame) containing the pre-defined parameter combinations to generate the simulations

## Author(s)

Tatjana Ammer <[tatjana.ammer@roche.com](mailto:tatjana.ammer@roche.com)>

## Examples

```
testsets <- loadTestsetDefinition()
str(testsets)
```

**mergeAnalytes**

*Function to combine analytes for defined categories*

## Description

The feature can either be the pathological fraction, the sample size or the overlap (category) individually or cumulative (\_cum). For a individualized categorisation see `getSubsetForDefinedCats`.

## Usage

```
mergeAnalytes(
  tableTCs,
  errorList,
  catList = NULL,
  catLabels = NULL,
  distTypes = TRUE
)
```

## Arguments

<code>tableTCs</code>	(data.frame) containing all information about the simulated test sets
<code>errorList</code>	(list) containing for each method the table with the computed error measurements
<code>catList</code>	(list) containing the categories to split the dataset
<code>catLabels</code>	(list) containing the labels that will be used for the categories
<code>distTypes</code>	(logical) indicating if 'catList' refers to analytes (FALSE, default) or distribution types (TRUE)

## Value

(list) containing the merged performance measurements grouped according to specified category

## Author(s)

Tatjana Ammer <[tatjana.ammer@roche.com](mailto:tatjana.ammer@roche.com)>

---

mergeSummaryErrors      *Helper function to combine all computed summary errors*

---

## Description

Helper function to combine all computed summary errors

## Usage

```
mergeSummaryErrors(  
  errorList,  
  nameVec,  
  errorParam = "MedianAbsPercErrorOV",  
  cutoffZ = FALSE  
)
```

## Arguments

errorList	(list) of the error lists for the different methods for which the summary errors should be combined
nameVec	(character) vector specifying the names of the methods
errorParam	(character) specifying for which error parameter the data frame should be generated
cutoffZ	(logical) indicating if a cutoff was set, needed for CRP case

## Value

(data frame) containing the summary errors per analyte per method

## Author(s)

Tatjana Ammer <tatjana.ammer@roche.com>

---

plotBarplot      *Plot method for generating a barplot out of the benchmark results*

---

## Description

Plot method for generating a barplot out of the benchmark results

**Usage**

```
plotBarplot(
  benchmarkRes,
  perDistType = FALSE,
  colList,
  nameList = NULL,
  withLabels = FALSE,
  withHorizLines = FALSE,
  title = NULL,
  xlim = NULL,
  xlab = "Mean of Absolute Z-Score Deviations",
  outputDir = NULL,
  filename = NULL,
  ...
)
```

**Arguments**

<code>benchmarkRes</code>	(data frame) containing the overall benchmark results
<code>perDistType</code>	(logical) indicating if one overall plot should be generated or if it should be separated by the distribution type
<code>colList</code>	(character) vector specifying the colors used for the different algorithms (should correspond to columns of benchmark results)
<code>nameList</code>	(character) vector specifying the names used in the legend (should correspond to columns of benchmark results), if <code>NULL</code> , <code>colnames</code> will be used
<code>withLabels</code>	(logical) indicating whether the corresponding values should be plotted as well (default: <code>FALSE</code> )
<code>withHorizLines</code>	(logical) indicating whether horizontal lines should be plotted for a better visual separation of the different categories (default: <code>FALSE</code> )
<code>title</code>	(character) specifying plot title
<code>xlim</code>	(numeric) vector specifying the limits in x-direction
<code>xlab</code>	(character) specifying the x-axis label
<code>outputDir</code>	(character) specifying a output directory
<code>filename</code>	(character) specifying a filename for the plot
<code>...</code>	additional arguments passed forward to other functions

**Value**

No return value. Instead, a plot is generated.

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

---

<code>plotBoxplot</code>	<i>Plot method for generating a boxplot of the benchmark results</i>
--------------------------	--

---

## Description

Plot method for generating a boxplot of the benchmark results

## Usage

```
plotBoxplot(
  errorList,
  colList,
  nameList,
  outline = TRUE,
  withMean = TRUE,
  withCats = FALSE,
  withDirect = TRUE,
  title = "",
  outputDir = NULL,
  filename = NULL,
  ylim1 = c(0, 100),
  ylim2 = c(100, 1000),
  ...
)
```

## Arguments

<code>errorList</code>	containing the overall benchmark results
<code>colList</code>	(character) vector specifying the colors used for the different algorithms (should correspond to columns of benchmark results)
<code>nameList</code>	(character) vector specifying the names used in the legend (should correspond to columns of benchmark results), if <code>NULL</code> , <code>colnames</code> will be used
<code>outline</code>	(logical) indicating whether outliers should be drawn ( <code>TRUE</code> , default), or not ( <code>FALSE</code> )
<code>withMean</code>	(logical) indicating whether the mean should be plotted as well (default: <code>TRUE</code> )
<code>withCats</code>	(logical) set to <code>TRUE</code> if categories (e.g. pathological fraction) should be plotted (default: <code>FALSE</code> )
<code>withDirect</code>	(logical) indicating whether the box of the direct method should be elongated to facilitate comparison (default: <code>TRUE</code> )
<code>title</code>	(character) specifying plot title
<code>outputDir</code>	(character) specifying a output directory
<code>filename</code>	(character) specifying a filename for the plot
<code>ylim1</code>	(numeric) vector specifying the limits in y-direction for the first granular scale
<code>ylim2</code>	(numeric) vector specifying the limits in y-direction for the second less detailed scale
<code>...</code>	additional arguments passed forward to other functions

**Value**

No return value. Instead, a plot is generated.

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

**plotScatterplot**      *Plot method for generating a scatterplot*

**Description**

Plot method for generating a scatterplot

**Usage**

```
plotScatterplot(
  errorList,
  colList,
  nameList,
  withColor = NULL,
  cats = NULL,
  title = "",
  outputDir = NULL,
  filename = NULL,
  xlim = NULL,
  ylim = NULL,
  xlab = NULL,
  ylab = NULL,
  ...
)
```

**Arguments**

<code>errorList</code>	(data frame) containing the overall benchmark results
<code>colList</code>	(character) vector specifying the colors used for the different algorithms (should correspond to columns of benchmark results)
<code>nameList</code>	(character) vector specifying the names used in the legend (should correspond to columns of benchmark results), if NULL, colnames will be used
<code>withColor</code>	(character) indicating if plot should be colored according to pathological fraction, sample size or pathological overlap left / right
<code>cats</code>	(character) specifying the category labels
<code>title</code>	(character) specifying plot title
<code>outputDir</code>	(character) specifying a output directory
<code>filename</code>	(character) specifying a filename for the plot

xlim	(numeric) vector specifying the limits in y-direction for the first granular scale
ylim	(numeric) vector specifying the limits in y-direction for the second less detailed scale
xlab	(character) specifying x-axis label
ylab	(character) specifying y-axis label
...	additional arguments passed forward to other functions

**Value**

No return value. Instead, a plot is generated.

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

**print.RWDRI**

*Standard print method for objects of class 'RWDRI'*

**Description**

Standard print method for objects of class 'RWDRI'

**Usage**

```
## S3 method for class 'RWDRI'
print(
  x,
  RIperc = c(0.025, 0.975),
  CIprop = 0.95,
  pointEst = c("fullDataEst", "medianBS", "meanBS"),
  truncNormal = FALSE,
  ...
)
```

**Arguments**

x	(object) of class 'RWDRI'
RIperc	(numeric) value specifying the percentiles, which define the reference interval
CIprop	(numeric) value specifying the central region for estimation of confidence intervals
pointEst	(character) specifying the point estimate determination: (1) using the full dataset ("fullDataEst"), (2) calculating the median from the bootstrap samples ("medianBS"), (2) works only if NBootstrap > 0 (3) calculating the mean from the bootstrap samples ("meanBS"), (3) works only if NBootstrap > 0
truncNormal	(logical) specifying if a normal distribution truncated at zero shall be assumed
...	additional arguments passed forward to other functions.

**Value**

No return value. Instead, a summary is printed.

**Author(s)**

Christopher Rank <christopher.rank@roche.com>

**progressInd**

*Function for setting up the progress indicator.*

**Description**

Function for setting up the progress indicator.

**Usage**

```
progressInd(value, maxValue, nCharMsg = 0)
```

**Arguments**

<b>value</b>	(integer) indicating the current number
<b>maxValue</b>	(integer) indicating the maximum number
<b>nCharMsg</b>	(integer) indicating the number of characters the message already has

**Value**

(character) returning generated progress message

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

**readResultFiles**

*Function for reading in the result files for one marker*

**Description**

Function for reading in the result files for one marker

**Usage**

```
readResultFiles(analyte, algo, path = NULL, tableTCs = NULL)
```

**Arguments**

analyte	(character) specifying analyte
algo	(character) specifying used algorithm
path	(character) specifying path to Results directories
tableTCs	(data frame) containing all information about the simulated test sets

**Value**

list with calculated results as RWDRI objects

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

---

readResultFilesAll     *Function for reading all results files.*

---

**Description**

Function for reading all results files.

**Usage**

```
readResultFilesAll(analytes, algo, baseDir = NULL, inputDir = NULL, tableTCs)
```

**Arguments**

analytes	(character) listing all analytes for which the result files should be parsed
algo	(character) specifying used algorithm
baseDir	(character) specifying the baseDir: Results will be used from baseDir/Results/algo/marker if baseDir is set, inputDir will be ignored; if baseDir is NULL, the current working directory will be used
inputDir	(character) specifying path directly to Results directories
tableTCs	(data frame) containing all information about the simulated test sets

**Value**

list with all calculated results as RWDRI objects for each marker

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

**readResultsAndComputeErrors**

*Function to read the result files and compute performance measures to create customized plots afterwards*

**Description**

Function to read the result files and compute performance measures to create customized plots afterwards

**Usage**

```
readResultsAndComputeErrors(
  workingDir = getwd(),
  algoName = NULL,
  subset = "all",
  cutoffZ = 5,
  ...
)
```

**Arguments**

workingDir	(character) specifying the working directory: Plots will be stored in workingDir/evalFolder and results will be used from workingDir/Results/algoName/biomarker;
algoName	(character) vector specifying one algorithm for which the performance measures should be evaluated
subset	(character, numeric, or data.frame) to specify for which subset the algorithm should be executed. character options: 'all' (default) for all test sets, a distribution type: 'normal', 'skewed', 'heavilySkewed', 'shifted'; a biomarker: 'Hb', 'Ca', 'FT4', 'AST', 'LACT', 'GGT', 'TSH', 'IgE', 'CRP', 'LDH'; 'runtime' for runtime analysis subset; numeric option: number of test sets per biomarker, e.g. 10; data.frame: customized subset of table with test set specifications
cutoffZ	(integer) specifying if and if so which cutoff for the absolute z-score deviation should be used to classify results as implausible and exclude them from the overall benchmark score (default: 5)
...	additional arguments to be passed to the method truncNormal (logical) specifying if a normal distribution truncated at zero shall be assumed

**Value**

(list) with (data frame) and a (list) with the computed performance measures

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

---

**restrictSet***Function to get error subsets for defined category and restriction.*

---

**Description**

Function to get error subsets for defined category and restriction.

**Usage**

```
restrictSet(overallCat, tableTCs, errorList, distType = TRUE, restrict = NULL)
```

**Arguments**

overallCat	(list) containing the categories to split the dataset
tableTCs	(data.frame) containing all information about the simulated test sets
errorList	(list) containing for each method the table with the computed error measurements
distType	(logical) indicating if 'overallCat' refers to analytes (FALSE, default) or distribution types (TRUE)
restrict	(character) indicating whether testcases should be filtered according to specified restriction, default NULL, e.g. fractionPathol <= 0.30

**Value**

(list) containing the merged performance measurements grouped according to specified category

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

---

**runDirectMethod***Convenience function to simulate the direct method for the specified subset*

---

**Description**

Convenience function to simulate the direct method for the specified subset

**Usage**

```
runDirectMethod(tableTCs = NULL, N = 120, cutoffZ = 5)
```

**Arguments**

tableTCs	(data frame) containing the pre-defined parameter combinations to generate the simulations
N	(integer) specifying the number of samples used as sample size for the direct method, default: 120
cutoffZ	(numeric) specifying if a cutoff should be used to classify results as implausible and exclude from analysis

**Value**

(data frame) with computed performance measures

**Author(s)**

Tatjana Ammer <tatjana.ammer@roche.com>

**Examples**

```
# example to run direct method only for test sets for hemoglobin (Hb)
testsets <- loadTestsetDefinition()
directRes <- runDirectMethod(tableTCs = testsets[testsets$Analyte == "Hb", ], N = 120, cutoffZ = 5)
```

**runTC\_usingRscript**      *Function for running test sets per algorithm per marker with calling Rscript for each test set*

**Description**

Function for running test sets per algorithm per marker with calling Rscript for each test set

**Usage**

```
runTC_usingRscript(
  biomarker = NULL,
  algoName = "myOwnAlgo",
  algoFunction = "estimateModel",
  sourceFiles = NULL,
  libs = NULL,
  params = NULL,
  decimals = FALSE,
  ris = FALSE,
  RIperc = c(0.025, 0.975),
  tableTCs = NULL,
  outputDir = NULL,
```

```

  inputDir = NULL,
  timeLimit = 14400,
  subsetDef = "all",
  verbose = TRUE,
  showWarnings = FALSE,
  ...
)

```

## Arguments

biomarker	(character) specifying the biomarker for which the algorithm should calculate RIs
algoName	(character) specifying the name of the algorithm that is evaluated
algoFunction	(character) specifying the name of the function needed for estimating RIs
sourceFiles	(list) containing all source files needed for executing the algorithm
libs	(list) containing all libraries needed for executing the algorithm
params	(list) with additional parameters needed for calling algoFunction
decimals	(logical) indicating whether the algorithm needs the number of decimal places (TRUE) or not (FALSE, default)
ris	(logical) indicating whether only percentiles and no model is estimated
RIperc	(numeric) value specifying the percentiles, which define the reference interval
tableTCs	(data.frame) with the information about the simulated test sets
outputDir	(character) specifying the outputDir: Results will be stored in outputDir/Results/algo/biomarker
inputDir	(character) specifying the inputDir: Data files should be stored in inputDir/Data/biomarker
timeLimit	(integer) specifying the maximum amount of time in seconds allowed to execute one single estimation (default: 14400 sec (4h))
subsetDef	(character) describing the specified subset of all test sets the algorithm is applied to, used for naming the progress file
verbose	(logical) indicating if the progress counter should be shown (default: TRUE)
showWarnings	(logical) indicating whether warnings from the call to the indirect method/algorithm should be shown (default: FALSE)
...	additional arguments to be passed to the method

## Value

(data frame) containing information about the test sets where the algorithm terminated the R session or failed to report a result

## Author(s)

Tatjana Ammer <tatjana.ammer@roche.com>

---

<code>setupDirStructure</code>	<i>Convenience function to set up the directory structure used for storing data and results.</i>
--------------------------------	--

---

## Description

Convenience function to set up the directory structure used for storing data and results.

## Usage

```
setupDirStructure(
  outputDir = NULL,
  onlyData = FALSE,
  onlyResults = FALSE,
  tableTCs = NULL,
  algoName = NULL
)
```

## Arguments

<code>outputDir</code>	(character) specifying the base output directory. From here, Data/biomarker and Result/algoName/biomarker directories are generated
<code>onlyData</code>	(logical) if set to TRUE, only the biomarker subdirectories are generated, name of output directory is used as it is (default:FALSE)
<code>onlyResults</code>	(logical) if set to TRUE, only the algoName/biomarker subdirectories are generated, name of output directory is used as it is (default:FALSE)
<code>tableTCs</code>	(data frame) containing the pre-defined parameter combinations to generate the simulations
<code>algoName</code>	(character) specifying the name of the algorithm used for creating the subdirectory

## Value

No return value. Instead, the directory structure is set up.

## Author(s)

Tatjana Ammer <tatjana.ammer@roche.com>

---

writeResFile	<i>Helper function to write result file when time out occured or R session terminated</i>
--------------	---

---

## Description

Helper function to write result file when time out occured or R session terminated

## Usage

```
writeResFile(  
  algoName,  
  biomarker,  
  N = 0,  
  error = NULL,  
  runtime = NULL,  
  filename = NULL,  
  outputDir = NULL  
)
```

## Arguments

algoName	(character) specifying the name of the algorithm that is evaluated
biomarker	(character) specifying the biomarker for which the algorithm should calculate RIIs
N	(numeric) specifying the number of input data points
error	(character) specifying the type of error (e.g. timeout, RSessionTerminated)
runtime	(numeric) specifying the computation time up until the error occured
filename	(character) specifying the filename for which the algorithm failed
outputDir	(character) specifying the outputDir: Data files should be stored in outputDir/Data/biomarker and Results will be stored in outputDir/Results/algo/biomarker

## Author(s)

Tatjana Ammer <tatjana.ammer@roche.com>

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