

Package: kwb.qmra (via r-universe)

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Title QMRA (quantitative microbial risk assessment)

Version 0.3.0

Description QMRA for water supply systems.

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URL <http://github.com/KWB-R/kwb.qmra/>

BugReports <https://github.com/KWB-R/kwb.qmra/issues>

Depends R (>= 4.0.0)

Imports dplyr (>= 1.0.2), EnvStats (>= 2.3.1), ggplot2 (>= 3.3.2), jsonlite (>= 1.7.2), kwb.utils (>= 0.7.0), magrittr (>= 1.5), plyr (>= 1.8.6), readr (>= 1.4.0), readxl (>= 1.3.1), rlang (>= 0.4.8), rmarkdown (>= 2.4), rvest (>= 0.3.6), sfsmisc (>= 1.1.7), shiny (>= 1.5.0), stringr (>= 1.4.0), tidyverse (>= 1.1.2), xml2 (>= 1.3.2)

Suggests ggrepel (>= 0.8.2), knitr (>= 1.30), qmra.db (>= 0.10.0), remotes (>= 2.2.0), sessioninfo (>= 1.1.1), stringi (>= 1.5.3), testthat (>= 3.0.0), tibble (>= 3.0.4)

VignetteBuilder knitr

Remotes github::kwb-r/kwb.utils@v0.7.0, github::kwb-r/qmra.db@v0.10.0

Encoding UTF-8

LazyLoad yes

RoxygenNote 7.1.1

Config/testthat.edition 3

Repository <https://kwb-r.r-universe.dev>

RemoteUrl <https://github.com/KWB-R/kwb.qmra>

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```
backcalc_infectionRisk_perDay
```

Backcalculate risk: infection (using optimise() function)

Description

Based on exposure days per year and target infection risk per year the acceptable daily infection risk is backcalculated

Usage

```
backcalc_infectionRisk_perDay(  
  target_infectionRisk_perYear = 1/10000,  
  exposures_daysPerYear = 1:365  
)
```

Arguments

```
target_infectionRisk_perYear  
  target infection risk per per year (default: 1/10000)  
exposures_daysPerYear  
  exposure days per year (default: 1 to 365)
```

Value

acceptable daily infection risk for given exposures per year and target infection risk per year

```
calc_health_risk
```

Risk calculation: health

Description

Risk calculation: health

Usage

```
calc_health_risk(  
  infectionRisk_perYear = 9.5 * 10^-4,  
  infection_to_illness = 0.7,  
  diseaseBurden_dalyPerCase = 1.5 * 10^-3,  
  fraction_population = 1,  
  target_dalyPerYearPerPerson = 1/1e+06  
)
```

Arguments

infectionRisk_perYear
 as retrieved by `calc_infection_risk()$events$infectionRisk_perYear` or user defined input (default: 9.5×10^{-4} infections/year)

infection_to_illness
 probability of illness given infection (default: 0.7, i.e. 70 percent illness probability giving infection)

diseaseBurden_dalyPerCase
 disabled adjusted life years per case. Value depends on multiple parameters such as illness type (default: 1.5×10^{-3})

fraction_population
 fraction of population at risk of getting the illness (default: 1, i.e. 100 percent of population can possibly get the illness, worst-case assumption assuming no prior immunization!)

target_dalyPerYearPerPerson
 target disabled adjusted life years (DALY) per person per year (default: 1/1000000 DALY per person per year, WHO standard)

Value

data frame with input parameters and calculated health risk outputs

See Also

[calc_infection_risk](#) for infection risk input

`calc_infection_risk` *Risk calculation: infection*

Description

Risk calculation: infection

Usage

```
calc_infection_risk(
  inflow_orgPerLitre = 10,
  treatment_logRemoval = 5.8,
  exposure_daysPerYear = 365,
  doseresponse_modelType = "dr.expo",
  waterConsumption_LitrePerDay = 1,
  target_infectionRisk_perYear = 1/10000,
  ...
)
```

Arguments

inflow_orgPerLitre
 concentration of microbiological parameter in inflow to water treatment plant
 (default: 10 Org/L)

treatment_logRemoval
 reduction of microbiological concentration in water treatment plant (default: 5.8 log)

exposure_daysPerYear
 exposure days per year (default: 365)

doseresponse_modelType
 dose response model to be used: "dr.expo" for exponential or "dr.betapoisson" for beta-poisson model

waterConsumption_LitrePerDay
 daily water consumption (default: 1 L/day)

target_infectionRisk_perYear
 NOT IMPLEMENTED YET!!!! target infection risk per per year. Only used if one of the following input parameters (i.e. "inflow_orgPerLitre", "treatment_logRemoval", "exposure_daysPerYear", "waterConsumption_LitrePerDay") is not defined (default: 1/10000).

... additional parameters used for dose response modelling. Depends on used dose-response model

Value

list with input parameters and calculated infection risk

See Also

[dr.expo](#) for exponential or [dr.betapoisson](#) for beta-poisson dose-response model

config_default_json *config_dummy_json:* *dummy configuration for kwb.qmra::opencpu_simulate_risk() in JSON format*

Description

JSON input configuration generated with R script (data-raw/config_json.R)

Usage

```
data(config_dummy_json)
```

Format

A json list with all configuration parameters used by [opencpu_simulate_risk](#), which are passed to [simulate_risk](#)

Examples

```
data("config_dummy_json")
config_dummy_json
```

config_dummy_json

config_default_json: default configuration developed by Christoph Sprenger for kwb.qmra::opencpu_simulate_risk() in JSON format

Description

JSON input configuration generated with R script (data-raw/config_json.R)

Usage

```
data(config_default_json)
```

Format

A json list with all configuration parameters used by [opencpu_simulate_risk](#), which are passed to simulate_risk

Examples

```
data("config_default_json")
config_default_json
```

config_read

Config: read configuration

Description

Config: read configuration

Usage

```
config_read(
  confDir = system.file("extdata/configs/dummy", package = "kwb.qmra")
)
```

Arguments

confDir	directory to read configuration files (Default: system.file('extdata/configs/dummy', package = 'kwb.qmra'))
---------	---

Value

stores configuration in R list structure

`config_write` *Config: create configuration*

Description

Config: create configuration

Usage

```
config_write(config, confName = "dummy", confDir = tempdir(), zipFiles = TRUE)
```

Arguments

<code>config</code>	a configuration as retrieved by config_read()
<code>confName</code>	name of configuration
<code>confDir</code>	directory to save the configuration files (Default: tempdir())
<code>zipFiles</code>	should also zipfile be created in folder confDir (default: TRUE)

Value

writes configuration in confDir subfolder defined in confName

`config_write_dummy` *Config: create dummy configuration*

Description

Config: create dummy configuration

Usage

```
config_write_dummy(  
  confDir = system.file("extdata/configs/dummy", package = "kwb.qmra")  
)
```

Arguments

<code>confDir</code>	directory to save the dummy configuration files (Default: system.file('extdata/config', package = 'kwb.qmra'))
----------------------	--

Value

writes dummy configuration in confDir

create_random_distribution
Create random distribution

Description

Create random distribution

Usage

```
create_random_distribution(
  type = "uniform",
  number_of_repeating = 1,
  number_of_events = 365,
  value = 10,
  min = 10,
  max = 1000,
  percent_within_minmax = 0.9,
  min_zero = 0.01,
  log10_min = default_min(min, max, min_zero, f = log10),
  log10_max = default_max(max, min_zero * 10, f = log10),
  log10_mean = (log10_min + log10_max)/2,
  log10_sdev = abs((log10_max - log10_mean)/get_percentile(percent_within_minmax)),
  mean = (default_min(min, max, min_zero) + default_max(max, 10 * min_zero))/2,
  sdev = abs((default_max(max, 10 * min_zero) -
    mean)/get_percentile(percent_within_minmax)),
  meanlog = mean(log(default_min(min, max, min_zero) + default_max(max, 10 *
    min_zero))/2),
  sdlog = abs(sd(c(default_min(min, max, min_zero, f = log), default_max(max, 10 *
    min_zero, f = log)))),
  mode = (default_min(min, max, min_zero) + default_max(max, 10 * min_zero))/2,
  debug = TRUE
)
```

Arguments

type	"uniform" calls runif(), "log10_uniform" calls 10^runif(number_of_events, log10_min, log10_max), "triangle" calls EnvStats::rtri(), "lognorm" calls rlnorm(), "norm" calls rnorm() and "log10_norm" calls 10^rnorm(number_of_events, mean = log10_mean, sdev = log10_sdev), (default: "uniform")
number_of_repeating	how often should the random distribution with the same parameters be generated (default: 1)
number_of_events	number of events
value	constant value (no random number), gets repeated number_of_events times (if 'type' = 'value')

min	minimum value (default: 10), only used if 'type' is "runif" or "triangle"
max	maximum value (default: 1000), only used if 'type' is "runif" or "triangle"
percent_within_minmax	percent of data point within min/max (default: 0.9 i.e. 90 percent)
min_zero	only used if 'type' is "log10_uniform" or "log10_norm", "norm" or "lognorm" and "min" value equal zero. In this case the zero is replaced by this value (default: 0.01), see also default_min
log10_min	minimum value (default: default_min(min, max, min_zero, f = log10)), only used if 'type' is "log10_uniform" or "log10_norm"
log10_max	maximum value (default: ifelse(max > 0, log10(max), log10_zero_threshold)), only used if 'type' is "log10_uniform" or "log10_norm"
log10_mean	mean value (default: (log10_min + log10_max)/2), only used if 'type' is "log10_norm"
log10_sdev	standard deviation (default: abs((log10_max - log10_mean) / get_percentile(0.95))), only used if 'type' is "log10_norm"
mean	mean value (default: (default_min(min, max, min_zero) / default_max(max, 10*min_zero)) / 2), only used if 'type' is "norm"
sdev	standard deviation (default: abs((default_max(max, 10*min_zero) - mean) / get_percentile(0.95))), only used if 'type' is "norm"
meanlog	log mean value (default: mean(log((min + max) / 2))), only used if 'type' is "lognorm"
sdlog	standard deviation (default: abs(sd(c(default_min(min, max, min_zero, f = log))))), only used if 'type' is "lognorm"
mode	(default: default_min(min, max, min_zero) + default_max(max, 10 * min_zero) / 2), only used if 'type' is "triangle"
debug	print debug information (default: TRUE)

Value

list with parameters of user defined random distribution and corresponding values

See Also

for random triangle see [rtri](#), for default min/max see [default_min](#), [default_max](#) and [get_percentile](#)

default_max

*Default Max***Description**

Default Max

Usage

```
default_max(org_max, new_max, f = c)
```

Arguments

<code>org_max</code>	original maximum value
<code>new_max</code>	new maximum value for replacement
<code>f</code>	function apply on "org_max" to transform to the correct dimension (e.g. "log" in case of "lognorm" or "log10" in case of log10_norm) (default: <code>c()</code>)

Examples

```
default_max(org_max = 2, new_max = 0.01, f = log10)
default_max(org_max = 0, new_max = 0.01, f = log10)
default_max(org_max = 2, new_max = 0.01, f = log)
default_max(org_max = 0, new_max = 0.01, f = log)
```

default_min*Default Min***Description**

Default Min

Usage

```
default_min(org_min, org_max, new_min, f = c)
```

Arguments

<code>org_min</code>	original minimum value
<code>org_max</code>	original maximum value
<code>new_min</code>	new minimum value for replacement
<code>f</code>	function apply on org_min and new_min to transform to the correct dimension (e.g. "log" in case of "lognorm" or "log10" in case of log10_norm), (default: <code>c()</code>)

Examples

```
default_min(org_min = 2, org_max = 100, new_min = 0.01, f = log10)
default_min(org_min = 0, org_max = 100, new_min = 0.01, f = log10)
default_min(org_min = 2, org_max = 100, new_min = 0.01, f = log)
default_min(org_min = 0, org_max = 100, new_min = 0.01, f = log)
```

distribution_repeater *Helper function: distribution repeater*

Description

Helper function: distribution repeater

Usage

```
distribution_repeater(  
  number_of_repeating = 10,  
  number_of_events = 365,  
  func,  
  ...  
)
```

Arguments

number_of_repeating	how often should the random distribution with the same parameters be generated (default: 1)
number_of_events	number of events
func	distribution function to be repeated (e.g. runif, rlnorm, rnorm)
...	further parameters passed to func

Value

data.frame with columns repeatID, eventID and values

Examples

```
distribution_repeater(  
  number_of_repeating = 2,  
  number_of_events = 10,  
  func = runif,  
  min = 1,  
  max = 10  
)
```

dr.betapoisson*Dose-response model: beta-poisson***Description**

Dose-response model: beta-poisson

Usage

```
dr.betapoisson(
  dose = sfsmisc::lseq(from = 1, to = 10^10, length = 1000),
  alpha = 0.328,
  N50 = 5430
)
```

Arguments

dose	vector of dose data (default: sfsmisc::lseq(from = 0.1, to = 10^10, length = 1000))
alpha	alpha (default: 3.28E-01)
N50	N50 (default: 5.43E+03)

Value

tibble

dr.db_download*Download dose-response model database from QMRAwiki***Description**

Download dose-response model database from QMRAwiki

Usage

```
dr.db_download(fromInternet = FALSE)
```

Arguments

fromInternet	download from internet (default: FALSE), if FALSE import from local copy
--------------	--

Value

tibble for different microbial parameters

Source

http://qmrawiki.canr.msu.edu/index.php?title=Table_of_Recommended_Best-Fit_Parameters

dr.db_model

*Generate table with different doses for dr.db_download()***Description**

Generate table with different doses for dr.db_download()

Usage

```
dr.db_model(
  dr.db = dr.db_download(),
  dose = sfsmisc::lseq(from = 1, to = 10^10, length = 1000)
)
```

Arguments

dr.db	as retrieved by dr.db_download(), default: dr.db_download()
dose	vector of dose data (default: sfsmisc::lseq(from=0.1, to = 10^10, length = 1000))

Value

tibble

dr.expo

*Dose-response model: exponential***Description**

Dose-response model: exponential

Usage

```
dr.expo(dose = sfsmisc::lseq(from = 1, to = 10^10, length = 1000), k = 0.572)
```

Arguments

dose	vector of dose data (default: sfsmisc::lseq(from = 0.1, to = 10^10, length = 1000))
k	k-value (default: 5.72E-01)

Value

tibble

`generate_random_values`

Create random distribution based on configuration file

Description

Create random distribution based on configuration file

Usage

```
generate_random_values(
    config,
    number_of_repeating = 1,
    number_of_events,
    debug = TRUE
)
```

Arguments

<code>config</code>	as retrieved by <code>config_read()</code>
<code>number_of_repeating</code>	how often should the random distribution with the same parameters be generated (default: 1)
<code>number_of_events</code>	number of events
<code>debug</code>	print debug information (default: TRUE)

Value

list random distributions based on configuration file

`get_percentile`

Helper function: get percentile

Description

Helper function: get percentile

Usage

```
get_percentile(percent_within_minmax = 0.9)
```

Arguments

<code>percent_within_minmax</code>	percent of data point within min/max (default: 0.9 i.e. 90 percent)
------------------------------------	---

Examples

```
get_percentile(0.9)  
get_percentile(0.95)
```

opencpu_config_read *OpenCPU wrapper: import config from CSV and convert to JSON*

Description

OpenCPU wrapper: import config from CSV and convert to JSON

Usage

```
opencpu_config_read(  
  confDir = system.file("extdata/configs/dummy", package = "kwb.qmra")  
)
```

Arguments

confDir	directory to read configuration files (Default: system.file('extdata/configs/dummy', package = 'kwb.qmra'))
---------	---

Value

stores configuration in JSON format

Examples

```
### Example json config file  
config_json <- kwb.qmra::opencpu_config_read()  
head(config_json)
```

opencpu_simulate_risk *OpenCPU wrapper: run risk calculation and convert results to JSON format*

Description

OpenCPU wrapper: run risk calculation and convert results to JSON format

Usage

```
opencpu_simulate_risk(
  config = config_read(),
  usePoisson = TRUE,
  debug = TRUE,
  lean = TRUE
)
```

Arguments

config	config object as retrieved by <code>kwb.qmra::config_read()</code>
usePoisson	should a poisson process (see function <code>dose_perEvent()</code>) be used to calculate the <code>dose_perEvent</code> (TRUE) or just the <code>exposure_perEvent</code> column (FALSE), (default: TRUE)
debug	print debug information (default: TRUE)
lean	if TRUE, a "lean" version of this function is called, see <code>kwb.qmra:::simulate_risk_lean</code> , (default: TRUE)

Value

JSON list with parameters of user defined random distribution and corresponding values

Examples

```
### Example simulation run
## Read from JSON
config_json <- kwb.qmra::opencpu_config_read()
config <- jsonlite::fromJSON(config_json)
## Optionally directly import from CSVs
# config <- kwb.qmra::config_read()
risk <- kwb.qmra::opencpu_simulate_risk(config)
risk_json <- jsonlite::toJSON(risk, pretty = TRUE)
writeLines(text = risk_json, "risk.json")
```

`plot_doseresponse` *plot reduction*

Description

plotting

Usage

```
plot_doseresponse(risk)
```

Arguments

risk list as retrieved by simulate_risk()

Value

ggplot for reduction

plot_effluent *plot effluent*

Description

plotting

Usage

`plot_effluent(risk)`

Arguments

risk list as retrieved by simulate_risk()

Value

ggplot for effluent

plot_event_dalys *plot dalys_per_event*

Description

plotting

Usage

`plot_event_dalys(risk)`

Arguments

risk list as retrieved by simulate_risk()

Value

ggplot for dalys_per_event

plot_event_dose *plot dose per event*

Description

plotting

Usage

```
plot_event_dose(risk)
```

Arguments

risk list as retrieved by simulate_risk()

Value

ggplot for dose per event

plot_event_exposure *plot exposure per event*

Description

plotting

Usage

```
plot_event_exposure(risk)
```

Arguments

risk list as retrieved by simulate_risk()

Value

ggplot for exposure per event

plot_event_illnessProb
plot illness probability

Description

plotting

Usage

`plot_event_illnessProb(risk)`

Arguments

`risk` list as retrieved by `simulate_risk()`

Value

ggplot for illness probability

plot_event_infectionProb
plot infection probability

Description

plotting

Usage

`plot_event_infectionProb(risk)`

Arguments

`risk` list as retrieved by `simulate_risk()`

Value

ggplot for infection probability

plot_event_volume *plot volume per event*

Description

plotting volume

Usage

```
plot_event_volume(risk)
```

Arguments

risk list as retrieved by simulate_risk()

Value

ggplot for volume per event

plot_inflow *plot inflow*

Description

plotting

Usage

```
plot_inflow(risk)
```

Arguments

risk list as retrieved by simulate_risk

Value

ggplot for inflow

`plot_reduction` *plot reduction*

Description

plotting

Usage

```
plot_reduction(risk)
```

Arguments

`risk` list as retrieved by `simulate_risk()`

Value

ggplot for reduction

`plot_total_dalys` *plot total DALYs*

Description

plotting

Usage

```
plot_total_dalys(risk, labelling = FALSE, title = "", tolerance = 1e-06)
```

Arguments

`risk` list as retrieved by `simulate_risk()`

`labelling` if TRUE labels with absolute DALYs will be plotted (default: FALSE)

`title` title for plot (default: "")

`tolerance` acceptable tolerance level of risk (default: 1E-6)

Value

ggplot for total DALYs

```
plot_total_illnessProb  
    plot total illness probability
```

Description

plotting

Usage

```
plot_total_illnessProb(risk, tolerance = 1e-04)
```

Arguments

risk	list as retrieved by simulate_risk()
tolerance	acceptable tolerance level of risk (default: 1E-4)

Value

ggplot for total illness probability

```
plot_total_infectionProb  
    plot total infection probability
```

Description

plotting

Usage

```
plot_total_infectionProb(risk, tolerance = 1e-04)
```

Arguments

risk	list as retrieved by simulate_risk()
tolerance	acceptable tolerance level of risk (default: 1E-4)

Value

ggplot for total infection probability

poisson_dose

Helper function: poisson distribution based on exposure per event

Description

Helper function: poisson distribution based on exposure per event

Usage

```
poisson_dose(exposure_perEvent)
```

Arguments

exposure_perEvent
exposed organisms per event

Value

dose per event based on poisson process

report_workflow

Create report (not working for "shiny" reports)

Description

Create report (not working for "shiny" reports)

Usage

```
report_workflow(  
  confDirs = system.file("extdata/configs/", package = "kwb.qmra"),  
  report_template_dir = system.file("extdata/report", package = "kwb.qmra"),  
  report_template_name = "workflow.Rmd",  
  report_output_dir = NULL,  
  openReport = TRUE  
)
```

Arguments

confDirs directory containing subdirectory/ies of QMRA configurations default: system.file("extdata/configs/", package = "kwb.qmra")
report_template_dir report template directory (default: system.file("extdata/report", package = "kwb.qmra"))
report_template_name default: "workflow.Rmd"

```

report_output_dir
  directory where report should be saved, if NULL report_template_dir is used
  (default: NULL)
openReport    open report in browser default: TRUE

```

Value

generate html report

<code>risk_dummy_json</code>	<i>risk_dummy_json:</i> example risk_json object returned by <code>kwb.qmra::opencpu_simulate_risk()</code>
------------------------------	--

Description

JSON risk result object generated with R script (data-raw/risk_json.R)

Usage

```
data(risk_dummy_json)
```

Format

A json list with all risk results generated by [opencpu_simulate_risk](#), using the dataset ‘config_dummy_json’

Examples

```

data("risk_dummy_json")
risk <- jsonlite:::fromJSON(risk_dummy_json)
### only show "stats" elements (skip "events" and "total" due to much data)
risk_stats <- list(stats_total = risk$stats_total,
                    stats_logremoval = risk$stats_logremoval)
jsonlite:::toJSON(risk_stats, pretty = TRUE)

```

<code>run_app</code>	<i>Run shiny app</i>
----------------------	----------------------

Description

Run shiny app

Usage

```

run_app(
  appDir = system.file("extdata/shiny", package = "kwb.qmra"),
  launch.browser = TRUE,
  ...
)

```

Arguments

appDir	directory of shiny app (default: system.file("extdata/shiny", package = "kwb.qmra"))
launch.browser	If true, the system's default web browser will be launched automatically after the app is started. Defaults to true in interactive sessions only. This value of this parameter can also be a function to call with the application's URL.
...	additional parameters passed to shiny::runApp

simulate_exposure *Simulate: exposure*

Description

Simulate: exposure

Usage

```
simulate_exposure(config, debug = TRUE)
```

Arguments

config	as retrieved by config_read()
debug	print debug information (default: TRUE)

Value

list with parameters of user defined exposure scenario (number of events and volumes per Event)

simulate_inflow *Simulate: inflow*

Description

Simulate: inflow

Usage

```
simulate_inflow(config, debug = TRUE, lean = FALSE)
```

Arguments

config	as retrieved by config_read()
debug	print debug information (default: TRUE)
lean	if TRUE, only the "events" are returned (in a reduced version, i.e. without column PathogenName and with all ID columns being of class integer), otherwise a list with the events in element events" and the corresponding parameters in element paras. The default is FALSE, i.e. events and parameters are returned in a list.

Value

list with parameters of user defined random distribution and corresponding values

`simulate_risk`

Simulate: risk

Description

Simulate: risk

Usage

```
simulate_risk(config, usePoisson = TRUE, debug = TRUE, lean = FALSE)
```

Arguments

<code>config</code>	as retrieved by <code>config_read()</code>
<code>usePoisson</code>	should a poisson process (see function <code>dose_perEvent()</code>) be used to calculate the <code>dose_perEvent</code> (TRUE) or just the <code>exposure_perEvent</code> column (FALSE), (default: TRUE)
<code>debug</code>	print debug information (default: TRUE)
<code>lean</code>	if TRUE, a "lean" version of this function is called, see <code>kwb.qmra::simulate_risk_lean</code>

Value

list with parameters of user defined random distribution and corresponding values

`simulate_treatment`

Simulate: treatment

Description

Simulate: treatment

Usage

```
simulate_treatment(config, wide = FALSE, debug = TRUE, lean = FALSE)
```

Arguments

<code>config</code>	as retrieved by <code>config_read()</code>
<code>wide</code>	if TRUE results will be converted to wide format (default: FALSE)
<code>debug</code>	print debug information (default: TRUE)
<code>lean</code>	lean (default: FALSE)

Value

list with parameters of user defined random distribution and corresponding values

`who_getTreatment` *Treatment: get WHO reductions*

Description

Treatment: get WHO reductions

Usage

`who_getTreatment()`

Value

list with \$untidy and \$tidy data frames

See Also

http://apps.who.int/iris/bitstream/10665/44584/1/9789241548151_eng.pdf#page=162

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