Package ‘COINr’

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Type Package

Title Composite Indicator Construction and Analysis

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Description A comprehensive high-level package for composite indicator construction and analysis. It is a "development environment" for composite indicators and scoreboards, which includes utilities for construction (indicator selection, denomination, imputation, data treatment, normalisation, weighting and aggregation) and analysis (multivariate analysis, correlation plotting, short cuts for principal component analysis, global sensitivity analysis, and more). A composite indicator is completely encapsulated inside a single hierarchical list called a "COIN". This allows a fast and efficient work flow, as well as making quick copies, testing methodological variations and making comparisons. It also includes many plotting options, both statistical (scatter plots, distribution plots) as well as for presenting results (maps, bar charts, radar charts, and more). Finally, three Shiny apps are available which enable fast data exploration, results presentation, and checking the effects of altering weights.

Full documentation is found in the online book at <https://bluefoxr.github.io/COINrDoc/>, as well as the vignette.

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aggregate

Aggregate indicators

Description

Takes indicator data and a specified structure and hierarchically aggregates according to the structure specified in IndMeta. Uses a variety of aggregation methods as specified by agtype, which can be different for each level of aggregation (see agtype_by_level).

Usage

aggregate(
  COIN,  # COIN object
  agtype = "arith_mean",  # The type of aggregation method. One of either:
    • "arith_mean" - weighted arithmetic mean
    • "median" - weighted median
    • "geom_mean" - weighted geometric mean
    • "harm_mean" - weighted harmonic mean
    • "copeland" - weighted Copeland method
    • "custom" - a custom function - see agfunc
    • "mixed" - a different aggregation method for each level. In this case, aggregation methods are specified as any of the previous options using the agtype_bylevel argument.
  agweights = NULL,  # The weights to use in the aggregation. This can either be: NULL, in which case it will use the weights that were attached to IndMeta and AggMeta in assemble() (if they exist), or A character string which corresponds to a named list of weights stored in .$Parameters$Weights. You can either add these manually or through rew8r(). E.g. entering agweights = "Original" will use the original weights read in on assembly. This is equivalent to agweights = NULL. Or, a data frame of weights to use in the aggregation.
  dset = NULL,  # Which data set (contained in COIN object) to use
  agtype_bylevel = NULL,
  agfunc = NULL,
  avail_limit = NULL,
  out2 = NULL
)

Arguments

COIN: COIN object
agtype: The type of aggregation method. One of either:
  • "arith_mean" - weighted arithmetic mean
  • "median" - weighted median
  • "geom_mean" - weighted geometric mean
  • "harm_mean" - weighted harmonic mean
  • "copeland" - weighted Copeland method
  • "custom" - a custom function - see agfunc
  • "mixed" - a different aggregation method for each level. In this case, aggregation methods are specified as any of the previous options using the agtype_bylevel argument.
agweights: The weights to use in the aggregation. This can either be: NULL, in which case it will use the weights that were attached to IndMeta and AggMeta in assemble() (if they exist), or A character string which corresponds to a named list of weights stored in .$Parameters$Weights. You can either add these manually or through rew8r(). E.g. entering agweights = "Original" will use the original weights read in on assembly. This is equivalent to agweights = NULL. Or, a data frame of weights to use in the aggregation.
dset: Which data set (contained in COIN object) to use
aggregate

agtype_bylevel  A character vector with aggregation types for each level. Note that if this is specified, agtype must be specified as agtype = "mixed", otherwise agtype_by_level will be ignored.

agfunc  A custom function to use for aggregation if agtype = "custom", of the type \( y = f(x, w) \), where \( y \) is a scalar aggregated value and \( x \) and \( w \) are vectors of indicator values and weights respectively. Ensure that NAs are handled (e.g. set na.rm = T) if your data has missing values.

avail_limit  A data availability threshold, below which aggregation returns NA. This parameter is the fraction of data availability needed in a given aggregation group to return an aggregated score. Specified as either NULL (default, aggregation values are always returned if possible) or a value between 0 and 1 (below this value of data availability, NA will be returned). See Details. are ignored during aggregation, so that as long as there is at least one value in an aggregation group

out2  Where to output the results. If "COIN" (default for COIN input), appends to updated COIN, otherwise if "df" outputs to data frame.

Details

This function aggregates indicators according to the index structure specified in IndMeta. It will either use a single aggregation method for all aggregation levels (by specifying agtype) or can use a different aggregation method for each level of the index (see agtype_by_level). Aggregation methods are typically weighted (e.g. weighted means), and the weights for the aggregation are specified using the agweights argument.

By default, this function will aggregate wherever possible - generally this means that if at least one value is available for a given unit inside an aggregation group, it will return an aggregated score. Optionally, you can also specify a data availability threshold which will instead return NA if the data availability (within group and for each unit) falls below the threshold. For example, you may have four indicators inside a group, and you might want to only produce an aggregated score if data availability is at least 50% - this would be specified by avail_limit = 0.5. It is also possible to specify different data availability thresholds for different levels of the index, by specifying avail_limit as a vector which has one value for every aggregation level (the first value gives the threshold for the first aggregation, and so on up to the final level).

Value

An updated COIN containing the new aggregated data set at .$Data$Aggregated.

Examples

# assemble a COIN first
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# normalise the data
ASEM <- normalise(ASEM, dset = "Raw")
# aggregate the data
ASEM <- COINr::aggregate(ASEM, agtype="arith_mean", dset = "Normalised")
# check aggregated data set exists
stopifnot(!is.null(ASEM$Data$Aggregated))
ASEM aggregate metadata

Description

This contains all the metadata for the aggregate groups, including the names, weights and codes. See the ASEM Portal for further information and detailed description of each indicator, and COINr documentation for the formatting of this data set.

Usage

ASEMaggMeta

Format

A data frame with 8 rows and 9 variables:

Source

https://composite-indicators.jrc.ec.europa.eu/asem-sustainable-connectivity/repository

ASEM raw indicator data

Description

A data set containing raw values of indicators for 51 countries, groups and denominators. See the ASEM Portal for further information and detailed description of each indicator, and COINr documentation for the formatting of this data set.

Usage

ASEMIndData

Format

A data frame with 51 rows and 60 variables.

Source

https://composite-indicators.jrc.ec.europa.eu/asem-sustainable-connectivity/repository
**ASEMIndMeta**

---

**ASEM indicator metadata**

**Description**

This contains all metadata for ASEM indicators, including names, weights, directions, etc. See the ASEM Portal for further information and detailed description of each indicator, and COINr documentation for the formatting of this data set.

**Usage**

ASEMIndMeta

**Format**

A data frame with 49 rows and 9 variables

**Source**

https://bluefoxr.github.io/COINrDoc/coins-the-currency-of-coinr.html#aggregation-metadata

---

**assemble**  

*Build COIN object*

**Description**

This takes the raw data provided by the user and puts it into a list format (COIN object) that is recognised by COINr. It also checks whether there are any syntax errors in the data provided. Optionally, you can exclude or include indicators using the include and exclude arguments. Note that if an indicator is specified in BOTH include and exclude, it will be excluded.

**Usage**

assemble(
    IndData,
    IndMeta,
    AggMeta,
    include = NULL,
    exclude = NULL,
    preagg = NULL,
    use_year = NULL,
    impute_latest = FALSE
)
Arguments

- **IndData**: A data frame of indicator data.
- **IndMeta**: A data frame containing auxiliary information for each indicator.
- **AggMeta**: A data frame specifying the names and weights of each aggregation group.
- **include**: Optional argument specifying a subset of indicator codes to include (default all indicators included).
- **exclude**: Optional argument specifying a subset of indicator codes to exclude (default none excluded).
- **preagg**: Set to TRUE if you want to assemble a COIN using pre-aggregated data (typically for ex-post analysis).
- **use_year**: If IndData includes a Year column, and there are multiple observations for each unit (one per year), this can be set to a target year or years. For example, setting use_year = 2020 will filter IndData to only include points from 2020. Setting to use_year = c(2019, 2020) will return a list of COINs. Set use_year = "all" to return a COIN for all years where data is available. Keep in mind that a COIN represents a single year of data.
- **impute_latest**: Logical: if TRUE, imputes missing data points using most recent value from previous years. If FALSE (default) simply extracts the data frame as is. This only works if !is.null(use_year) and there are previous years of data available (before use_year). Currently does not support imputation using future values or interpolation.

Details

A "COIN" is an S3 class which is a structured list of indicator data, metadata, results and methodology which is used throughout COINr. COINs are a convenient way to store all variables relating to the composite indicator in a single named object. This keeps the workspace tidy, but also allows fast and concise calls to functions, as well as copying COINs to introduce methodological variations, and enables complex operations such as global sensitivity analysis (see sensitivity()).

If panel data is input to this function, it will output a tibble of COINs (see use_year). This feature is currently under development and more support will be included for these tibbles of COINs in COINr over time.

For general information on COINs see the COINr vignette as well as the relevant chapter in the COINr online documentation.

For details on copying, adjusting and comparing COINs see the COINr chapter on adjustments and comparisons.

Value

A "COIN" S3 class object (list) formatted to the specifications of COINr. If the input is panel data and use_year is set to return multiple years, this function returns a tibble of COINs, indexed by the year. This latter feature is new and currently under development.
**BoxCox**

**Examples**

```r
# build the ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
```

---

**BoxCox**  
*Box Cox transformation*

**Description**

Simple Box Cox, with no optimisation of lambda. See [COINr online documentation](https://coinnr.r-forge.r-project.org/) for more details.

**Usage**

```r
BoxCox(x, lambda, makepos = TRUE)
```

**Arguments**

- `x`  
  A vector or column of data to transform
- `lambda`  
  The lambda parameter of the Box Cox transform
- `makepos`  
  If `TRUE` (default) makes all values positive by subtracting the minimum and adding 1.

**Value**

A vector of length `length(x)` with transformed values.

**See Also**

- `treat()` Outlier treatment

**Examples**

```r
# get a column of data with outliers
x <- ASEMIndData$Tariff
# Apply Box Cox
xBox <- BoxCox(x, lambda = 2)
# plot one against the other
plot(x, xBox)
```
### build_ASEM

*Builds ASEM example*

**Description**

A short cut function for building the ASEM COIN. This builds the ASEM COIN up to and including aggregated results. See the ASEM Portal for the underlying data set and online documentation for more information on COINs.

**Usage**

```r
build_ASEM()
```

**Value**

COIN class object (a list) of the ASEM index, as well as information printed to the console - see `assemble()`.

**Examples**

```r
# Build the ASEM COIN
ASEM <- build_ASEM()
```

### checkData

*Detailed unit data check and screener by data availability*

**Description**

Gives detailed tables of data availability, and optionally screens units based on a data availability threshold and presence of zeros. Units can be optionally "forced" to be included or excluded, making exceptions for the data availability threshold.

**Usage**

```r
checkData(
  COIN,
  dset = NULL,
  ind_thresh = NULL,
  zero_thres = NULL,
  unit_screen = "none",
  Force = NULL,
  out2 = "COIN"
)
```
checkData

Arguments

COIN The COIN object
dset The data set to be checked/screened
ind_thresh A data availability threshold used for flagging low data and screening units if unit_screen != "none". Default 0.66. Specify as a fraction.
zero_thresh As ind_thresh but for non-zero values. Defaults to 0.05, i.e. it will flag any units with less than 5% non-zero values (equivalently more than 95% zero values).
unit_screen Specifies whether and how to screen units based on data availability or zero values.
  • If set to "none" (default), does not screen any units.
  • If set to "byNA", screens units with data availability below ind_thresh
  • If set to "byzeros", screens units with non-zero values below zero_thresh
  • If set to "byNAandzeros", screens units based on either of the previous two criteria being true.
  • If you simply want to force a unit or units to be excluded (without any other screening), use the Force argument and set unit_screen = TRUE. unit_screen != "none" outputs a new data set .$Data$Screened.
Force A data frame with any additional countries to force inclusion or exclusion. First column is "UnitCode". Second column "Status" either "Include" or "Exclude" for each country to force.
out2 Where to output the results. If "COIN" (default for COIN input), appends to updated COIN, otherwise if "list" outputs to data frame.

Details

The two main criteria of interest are NA values, and zeros. The summary table gives percentages of NA values for each unit, across indicators, and percentage zero values (as a percentage of non-NA values). Each unit is flagged as having low data or too many zeros based on thresholds.

This function currently only supports COINS as inputs, not data frames.

Value

An updated COIN with data frames showing missing data in .$Analysis, and if unit_screen != "none" outputs a new data set .$Data$Screened. If out2 = "list" wraps missing data stats and screened data set into a list.

Examples

# build ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# return stats to the COIN, plus screened data set, return to list
ScreenedData <- checkData(ASEM, dset = "Raw", unit_screen = "byNA", ind_thresh = 0.9, out2 = "list")
# See which units were removed
print(ScreenedData$RemovedUnits)
coin2Excel  
*Write a COIN to Excel*

**Description**

Takes a COIN and writes all main data tables and other things to an Excel file. This uses the 'openxlsx' package.

**Usage**

coin2Excel(COIN, fname = "COINresults.xlsx")

**Arguments**

- **COIN**: A COIN object
- **fname**: The file name to write to, as a character string

**Value**

An Excel workbook with each table on a separate named tab.

**Examples**

```r
## Here we write a COIN to Excel, but this is done to a temporary directory
## to avoid "polluting" the working directory when running automatic tests.
## In a real case, set fname to a directory of your choice.
ASEM <- build_ASEM()
# write to Excel in temporary directory
coin2Excel(ASEM, fname = paste0(tempdir(), "\ASEM_results.xlsx"))
# spreadsheet is at:
print(paste0(tempdir(), "\ASEM_results.xlsx"))
# now delete temporary file to keep things tidy in testing
unlink(paste0(tempdir(),"\ASEM_results.xlsx"))
```

COINToolIn  
*Import data directly from COIN Tool*

**Description**

This provides a direct interface for reading a COIN Tool input deck and converting it to COINr. You need to provide a COIN Tool file, with the "Database" sheet properly compiled.

**Usage**

COINToolIn(fname, makecodes = FALSE, oldtool = FALSE)
Arguments

fname  The file name and path to read, e.g. "C:/Documents/COINToolFile.xlsx".
makecodes  Logical: if TRUE, will generate short indicator codes based on indicator names, otherwise if FALSE, will use COIN Tool indicator codes "Ind.01", etc. Currently only does this for indicators, not aggregation groups.
oldtool  Logical: if TRUE, compatible with old COIN Tool (pre-release, early 2019 or earlier). There are some minor differences on where the elements are found.

Details

The COIN Tool is an Excel-based tool for building composite indicators. See COINr online documentation for more details and an example.

Value

A list containing:

- .$IndData  A data frame of imported indicator data to be input as the IndData argument in assemble().
- .$IndMeta  A data frame of imported indicator metadata to be input as the IndMeta argument in assemble().
- .$AggMeta  A data frame of imported aggregation metadata to be input as the AggMeta argument in assemble().

See Also

- coin2Excel() Export COIN contents to Excel

Examples

```r
## Not run:
## This example downloads a COIN Tool spreadsheet containing example data, saves it to a temporary directory, unzips, and reads into R. Finally it assembles it into a COIN.

# Make temp zip filename in temporary directory
tmpz <- tempfile(fileext = ".zip")
# Download an example COIN Tool file to temporary directory
# NOTE: the download.file() command may need its "method" option set to a specific value depending on the platform you run this on. You can also choose to download/unzip this file manually.
# Unzip
CTpath <- unzip(tmpz, exdir = tempdir())
# Read COIN Tool into R
l <- COINToolIn(CTpath, makecodes = TRUE)
# We can finish by assembling it
CT_exampleCOIN <- assemble(l$IndData, l$IndMeta, l$AggMeta)
```
coin_win

## Description
To be used inside `treat()` to avoid repetitions. Winsorises a numerical vector of data.

#### Usage

```r
coin_win(
  icol,
  winmax,
  winchange = TRUE,
  t_skew = 2,
  t_kurt = 3.5,
  icode = NULL
)
```

#### Arguments

- **icol**: The vector of data to Winsorise.
- **winmax**: The maximum number of points to Winsorise for each indicator. If NA, will keep Winsorising until skewness and kurtosis thresholds achieved (but it is likely this will cause errors).
- **winchange**: Logical: if TRUE, Winsorisation can change direction from one iteration to the next. Otherwise if FALSE (default), no change.
- **t_skew**: Absolute skew threshold (default 2).
- **t_kurt**: Kurtosis threshold (default 3.5).
- **icode**: The indicator name - used for error messages in `treat()`.

#### Details
Outliers are identified according to skewness and kurtosis thresholds. The algorithm attempts to reduce the absolute skew and kurtosis by successively Winsorising points up to a specified limit. The process is detailed in the COINr online documentation.

#### Value
A list containing:
- `.icol` the vector of treated data
- `.imax` the indices of elements of the vector that were Winsorised as high values
- `.imin` the indices of elements of the vector that were Winsorised as low values
- `.winz` the total number of Winsorised points
See Also

- `treat()` Outlier treatment

Examples

```r
# get a column of data with outliers
x <- ASEMIndData$Tariff
# Winsorise up to five points
winlist <- coin_win(x, winmax = 5)
# check the differences
data.frame(
  Orig = x,
  Treated = winlist$icol,
  Changes = ifelse(x == winlist$icol, "Same", "Treated"))
```

---

**colourTable**

*Conditionally formatted table*

Description

Given a data frame, generates a conditionally-formatted html table using reactable. This function is used by `iplotTable()`. It is a quick wrapper for `reactable::reactable`.

Usage

```r
colourTable(
  df,
  freeze1 = TRUE,
  sortcol = NULL,
  sortorder = "desc",
  searchable = TRUE,
  pagesize = 10,
  cell_colours = NULL,
  reverse_colours = FALSE
)
```

Arguments

- `df` A data frame to be displayed as a table.
- `freeze1` If TRUE (default), freezes the first column. This may be for example the unit name or code.
- `sortcol` A column name to sort the table by. Defaults to first numerical column. Set to "none" to disable.
- `sortorder` Either "desc" for sorted column to be sorted from high to low (default) or "asc" for the opposite.
compareDF

searchable If TRUE, includes a search box
pagesize The number of rows to display on each page.
cell_colours A character vector of at least two colour codes (e.g. Hex codes) to use for the colour palette. Should be in order of low to high values. Defaults to a simple green palette of c("#eef4", "#358554"). See grDevices::colorRamp() for more info.
reverse_colours If TRUE, reverses the colour map - useful for rank tables where lowest numbers mean high scores.

Value
An interactive table generated by reactable.

See Also
- iplotTable() Interactive table of indicator data (from a COIN)

Examples

```r
# some random data
df <- as.data.frame(matrix(runif(12), 3, 4))
# a names column
df <- cbind(Rnames = letters[1:3], df)
# round it
df <- roundDF(df)
# make a table
colourTable(df)
```

Description
A custom function for comparing two data frames of indicator data, to see whether they match up, at a specified number of significant figures.

Usage

```r
compareDF(df1, df2, matchcol, sigfigs = 5)
```

Arguments

df1 A data frame
df2 Another data frame
matchcol A common column name that is used to match row order. E.g. this might be UnitCode.
sigfigs The number of significant figures to use for matching numerical columns
Details

This function compares numerical and non-numerical columns to see if they match. Rows and columns can be in any order. The function performs the following checks:

- Checks that the two data frames are the same size
- Checks that column names are the same, and that the matching column has the same entries
- Checks column by column that the elements are the same, after sorting according to the matching column

It then summarises for each column whether there are any differences, and also what the differences are, if any.

This is intended to cross-check results. For example, if you run something in COINr and want to check indicator results against external calculations.

Value

A list with comparison results. List contains:

- `.Same`: overall summary: if TRUE the data frames are the same according to the rules specified, otherwise FALSE.
- `.Details`: details of each column as a data frame. Each row summarises a column of the data frame, saying whether the column is the same as its equivalent, and the number of differences, if any. In case the two data frames have differing numbers of columns and rows, or have differing column names or entries in matchcol, `.Details` will simply contain a message to this effect.
- `.Differences`: a list with one entry for every column which contains different entries. Differences are summarised as a data frame with one row for each difference, reporting the value from df1 and its equivalent from df2.

Examples

```r
# take a sample of indicator data (including the UnitCode column)
data1 <- ASEMIndData[c(2,12:15)]
# copy the data
data2 <- data1
# make a change: replace one value in data2 by NA
data2[1,2] <- NA
# compare data frames
compareDF(data1, data2, matchcol = "UnitCode")
```
### Description
Takes two COINs, and generates a rank comparison between specified indicator/aggregates. COINs must share at least some common unit codes, and the indicator selected by `isel`.

### Usage
```
compTable(
  COIN1,
  COIN2,
  dset = "Raw",
  isel,
  COINnames = NULL,
  sort_by = "AbsRankChange"
)
```

### Arguments
- **COIN1**: First COIN
- **COIN2**: Second COIN
- **dset**: The data set of interest
- **isel**: The indicator/column of interest
- **COINnames**: An optional character vector of the names of `COIN1` and `COIN2`, to be used in the table headers.
- **sort_by**: If "RankCOIN1", sorts by the indicator values of `COIN1`, if "RankCOIN2", sorts by `COIN2`, if "RankChange", sorts by rank change, and if "AbsRankChange" sorts by absolute rank change.

### Value
A data frame with ranks and rank changes between two COINs.

### See Also
- `compTableMulti()` Comparison table between multiple COINs

### Examples
```
ASEM <- build_ASEM()
# Make a copy
ASEMAltNorm <- ASEM
# Edit .$Method
ASEMAltNorm$Method$normalise$ntype <- "borda"
```
# Regenerate
ASEMAltNorm <- regen(ASEMAltNorm, quietly = TRUE)
# compare
CT <- compTable(ASEM, ASEMAltNorm, dset = "Aggregated", isel = "Index")

---

## compTableMulti

### Description

Takes multiple COINs (two or more), and generates a rank comparison for a single indicator or aggregate.

### Usage

```
compTableMulti(
  COINs,
  dset = "Aggregated",
  isel = "Index",
  tabtype = "Ranks",
  ibase = 1,
  sort_table = TRUE,
  extra_cols = NULL
)
```

### Arguments

- **COINs**: A list of COINs
- **dset**: The data set to extract the indicator from (must be present in each COIN). Default "Aggregated".
- **isel**: Code of the indicator or aggregate to extract from each COIN (must be present in the specified data set of each COIN). Default "Index".
- **tabtype**: The type of table to generate - "Ranks", "Diffs", or "AbsDiffs".
- **ibase**: The index of the COIN to use as a base comparison
- **sort_table**: If TRUE, sorts by the base COIN (ibase) (default).
- **extra_cols**: A character vector of any extra columns to include from the COIN referenced by ibase. For example, this could include group columns.

### Value

Rank comparison table as a data frame

### See Also

- `compTable()` Comparison table between two COINs
Examples

```r
ASEM <- build_ASEM()
# Make a copy
ASEMAltNorm <- ASEM
# Edit .$Method
ASEMAltNorm$Method$normalise$ntype <- "borda"
# Regenerate
ASEMAltNorm <- COINr::regen(ASEMAltNorm, quietly = TRUE)
# compare
ctable <- compTableMulti(list(ASEM, ASEMAltNorm), dset = "Aggregated", isel = "Index")

# add more COINs to the list to see more cols in the table...
```

copeland  

Copeland scores

Description

Aggregates a data frame of indicator values into a single column using the Copeland method. This function is used inside `aggregate()`, and calls `outrankMatrix()`.

Usage

```r
copeland(ind_data, w = NULL)
```

Arguments

- `ind_data` A data frame or matrix of indicator data, with observations as rows and indicators as columns. No other columns should be present (e.g. label columns).
- `w` A vector of weights, which should have length equal to `ncol(ind_data)`. Weights are relative and will be re-scaled to sum to 1. If `w` is not specified, defaults to equal weights.

Value

Numeric vector of Copeland scores.

Examples

```r
# get a sample of a few indicators
ind_data <- ASEMIndData[12:16]
# calculate outranking matrix
cop_results <- copeland(ind_data)
# check output
stopifnot(length(cop_results$Scores) == nrow(ind_data))
```
corrweightscat  

Scatter plot of correlations against weights

Description

Plots correlations on the x axis and weights on the y axis. Allows a selected highlighted point and a line showing low correlation boundary. This function is intended for use inside `rew8r()`.

Usage

corrweightscat(
  dat,
  facet = FALSE,
  acvar = NULL,
  linesw = FALSE,
  locorval = NULL,
  hicorval = NULL
)

Arguments

dat  Data frame with first col indicator codes, second is weights, third is correlations
facet Logical: if TRUE creates subplots.
acvar Active variable to highlight (one of the indicator codes)
linesw Whether to plot a vertical line showing low correlation boundary
locorval x value of low correlation line
hicorval x value of high correlation line

Details

Since this plot is really only intended for use inside `rew8r()` no example is provided.

Value

A scatter plot generated using plotly, also outputs event data (the clicked indicator).

See Also

- `rew8r()` Interactive app for adjusting weights and seeing effects on correlations
- `getCorr()` Get correlations between indicators/levels
denominate

Denominate indicator data sets

Description

Denominates (divides) indicators by other "denominator" indicators that are either input here or were attached as "Den_*" columns of IndData when assembling the COIN. This function can work either on COINs or on data frames.

Usage

denominate(
  obj,
  dset = NULL,
  specby = NULL,
  denomby = NULL,
  scaledenoms = NULL,
  denominators = NULL,
  out2 = "COIN"
)

Arguments

obj
COIN object or a data frame of indicator data to be denominated. If a data frame, must include a UnitCode column.

dset
The data set to denominate (only if COIN used as input)

specby
Selects the source of the specifications for denomination.

  • If "metadata", uses the denominator column in .$metadata.

  • If "user", takes a character vector of denominator codes (one for each indicator, with NA for indicators that should not be denominated, and in the same order as the indicators).

denomby
Character vector specifying which denominators to use for each indicator. Only used if specby = "user". For indicators with no denomination, set elements to NA. Elements must be column names of denominators.

scaledenoms
This allows the possibility to scale denominators if needed. For example, if GDP is a denominator and is measured in dollars, dividing will create very small numbers (order 1e-10 and smaller) which could cause problems with numerical precision. This should be a named list of the form e.g. list(Den_GDP = 1e-9), where the name is the denominator to be scaled, and the entry is a factor to multiply the denominator values by. In the example, this would multiply GDP values by 1e-9, which (if the original values are in dollars) would scale them to billions of dollars. The list can include more than one entry, corresponding to any denominators that are present.
denominate

**denominators**
A data frame of denominator data. Columns should be denominator data, with column names corresponding to entries in denomby. This must also include a UnitCode column to match units (ordering is unimportant, this is done inside the function). Ensure that the unit codes correspond to the unit codes in the indicator data.

**out2**
Where to output the results. If "COIN" (default for COIN input), appends to updated COIN, otherwise if "df" outputs to data frame.

**Details**

Typically, the aim here is to convert extensive (size-related) variables into intensive variables (comparable between units of different sizes). There is also the option scaledenoms to scale denominators to avoid very small or very large numbers resulting.

This function expects that denominators$UnitCode contains all unit codes found in the data frame to be denominated. Unused unit codes (rows) in denominators will be ignored. Note that some national-level denominator data is available inside COINr at COINr::WorldDenoms.

See online documentation for further details and examples.

**Value**

If out2 = COIN and obj is a COIN, returns an updated COIN object, with new dataset .$Data$Denominated of denominated indicators. Otherwise returns a data frame of denominated indicator data.

**See Also**

- WorldDenoms A data set of some common national-level denominators.

**Examples**

```r
# assemble ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# denote using specs present on assembly
ASEM <- denominate(ASEM, dset = "Raw")

# OR, use function on data frame
# Get a sample of indicator data (note must be indicators plus a "UnitCode" column)
IndData <- ASEMIndData[cbind("UnitCode", "Goods", "Services", "FDI")]
# Also get some denominator data
Denoms <- ASEMIndData[cbind("UnitCode", "Den_Pop", "Den_GDP")]
# Denominate one by the other
IndDataDenom <- denominate(IndData, denomby = c("Den_GDP", NA, "Den_Pop"), denominators = Denoms)
```
**effectiveWeight**  
*Effective weights*

**Description**
This calculates the effective weights of each element in the indicator hierarchy. This is useful for understanding e.g. the true weight of each indicator in the framework and is also used in `plotframework()`.

**Usage**
```
effectiveWeight(COIN)
```

**Arguments**
- **COIN**
  *COIN object, or list with first entry is the indicator metadata, second entry is the aggregation metadata*

**Value**
A list with effective weights, as well as a data frame with labels and parents for the sunburst plot.

**Examples**
```
# build ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# get effective weights
effwts <- effectiveWeight(ASEM)
```

---

**extractYear**  
*Impute panel data*

**Description**
Given a data frame of the IndData format, with a Year column, imputes any missing data using the latest available year. This function is used inside `assemble()`.

**Usage**
```
extractYear(use_year, IndData, impute_latest = FALSE)
```
extractYear

Arguments

use_year The year of data to extract and impute.
IndData A data frame of indicator data, containing a Year column and with multiple observations for each unit code.
impute_latest Logical: if TRUE, imputes missing data points using most recent value from previous years. If FALSE (default) simply extracts the data frame as is.

Details

This expects a data frame in the IndData format, i.e. it should at least have a UnitCode column, and a Year column, as well as other columns that are to be imputed. It also presumes that there are multiple observations for each unit code, i.e. one per year. It then searches for any missing values in the target year, and replaces them with the equivalent points from previous years. It will replace using the most recently available point.

Value

A list containing:

- $.IndDataImp: An IndData format data frame from the specified year (use_year), with missing data imputed using previous years (where possible).
- $.DataYears: A data frame in the same format as IndData, where each entry shows which year each data point came from. Points where there was no missing data will have use_year, imputed points will have the corresponding year used to impute, and any points in $.IndDataImp which are still NA will be NA.
- $.ImpTable: A data frame where each row is a point that was successfully imputed. This is a filtered and arranged version of $.DataYears that focuses only on the imputed points.
- $.NImputed: The number of imputed points.

See Also

- assemble() Assemble a COIN - this function optionally calls extractYear().
- impute() Impute data using other imputation options (not using panel data).

Examples

# artificial example using ASEM data
# We only have one year of data so we copy it and "pretend" that they are from different years
# First, introduce 3 NAs
dat2018 <- ASEMIndData
dat2018[2, 12] <- NA
dat2018[4, 14] <- NA
# Now make copy, pretending it is the previous year
dat2017 <- ASEMIndData
dat2017$Year <- 2017
# This df still has one missing point
dat2017[4, 14] <- NA
Finally we have a 2016 data frame where none of the previous points are missing:

```r
dat2016 <- ASEMIndData
dat2016$Year <- 2016
```

We can now put them together:

```r
IndData <- rbind(dat2018, dat2017, dat2016)
```

And extract the 2018 data, with missing data taken from previous years:

```r
Imp <- extractYear(2018, IndData, impute_latest = TRUE)
```

View which points have been imputed and the years of data used:

```r
Imp$ImpTable
```

---

**geoMean**

*Weighted geometric mean*

### Description

Weighted geometric mean of a vector. NA are skipped by default. This function is used inside `aggregate()`.

### Usage

```r
geoMean(x, w = NULL)
```

### Arguments

- **x**: A numeric vector of positive values.
- **w**: A vector of weights, which should have length equal to `length(x)`. Weights are relative and will be re-scaled to sum to 1. If `w` is not specified, defaults to equal weights.

### Value

The geometric mean, as a numeric value.

### Examples

```r
# a vector of values
x <- 1:10
# a vector of weights
w <- runif(10)
# weighted geometric mean
geoMean(x, w)
```
geoMean_rescaled

Rescaled weighted geometric mean

Description

NOTE this function is not really in use but is kept here for the moment. Not sure it is very useful.

Usage

geoMean_rescaled(x, w = NULL)

Arguments

x  A numeric vector of positive values.

w  A vector of weights, which should have length equal to length(x). Weights are relative and will be re-scaled to sum to 1. If w is not specified, defaults to equal weights.

Details

Weighted geometric mean of a vector. Here, any zero or negative values are automatically dealt with by re-scaling the data to be all positive, i.e. it shifts so that the minimum is equal to 0.1.

Note that this could be better achieved by normalising first. However, following default normalisation between 0 and 100, this function offers a quick way to test the effect of a geometric mean, for example in a sensitivity analysis, and avoids bugs arising.

Value

Rescaled weighted geometric mean, as a numeric value.

Examples

# a vector of values
x <- 1:10
# a vector of weights
w <- runif(10)
# rescaled weighted geometric mean
geoMean_rescaled(x, w)
getCorr

Get different types of correlation matrices

Description

Helper function for getting correlations between indicators. This retrieves subsets of correlation matrices between different aggregation levels, in different formats.

Usage

getCorr(
  COIN,
  dset,
  icodes = NULL,
  aglevs = NULL,
  cortype = "pearson",
  pval = 0.05,
  withparent = TRUE,
  grouplev = NULL
)

Arguments

COIN The COIN object
dset The target data set
icodes An optional list of character vectors where the first entry specifies the indicator/aggregate codes to correlate against the second entry (also a specification of indicator/aggregate codes).
aglevs The aggregation levels to take the two groups of indicators from. See getIn() for details. Defaults to indicator level.
cortype The type of correlation to calculate, either "pearson", "spearman", or "kendall".
pval The significance level for including correlations. Correlations with \( p > pval \) will be returned as NA. Default 0.05. Set to 0 to disable this.
withparent If TRUE, and aglev[1] != aglev[2], will only return correlations of each row with its parent.
grouplev The aggregation level to group correlations by if aglev[1] == aglev[2]. By default, groups correlations into the aggregation level above. Set to 0 to disable grouping and return the full matrix.

Details

Note that this function can only call correlations within the same data set (i.e. only one data set in .$Data).
getCronbach

Value

A data frame of correlation values in long format. Correlations with \( p > pval \) will be returned as NA.

See Also

- plotCorr() Plot correlation matrices of indicator subsets

Examples

```r
# build ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# correlations of indicators in Political pillar
corrs <- getCorr(ASEM, dset = "Raw", icodes = "Political", aglevs = 1)
```

getCronbach

Cronbach’s alpha

Description

Calculates Cronbach’s alpha, a measure of statistical reliability. Cronbach’s alpha is a simple measure of “consistency” of a data set, where a high value implies higher reliability/consistency.

Usage

```r
getCronbach(
  COIN,
  dset = "Raw",
  icodes = NULL,
  aglev = NULL,
  use = "pairwise.complete.obs"
)
```

Arguments

- **COIN**: A COIN or a data frame containing only numerical columns of data.
- **dset**: The data set to check the consistency of.
- **icodes**: Indicator codes if a subset of dset is requested
- **aglev**: The aggregation level to take icodes from (see getIn() for details)
- **use**: Argument to pass to stats::cor to calculate the covariance matrix. Default "pairwise.complete.obs".

Details

This function simply returns Cronbach’s alpha. If you want a lot more details on reliability, the 'psych' package has a much more detailed analysis.
Value

Cronbach alpha as a numerical value.

Examples

```r
# build ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# get Cronbach of indicators in Physical pillar
getCronbach(ASEM, dset = "Raw", icodes = "Physical", aglev = 1)
```

---

**getIn**

Get subsets of indicator data

Description

This function does a number of things that are useful for many COINr functions and operations. First, it checks to see what kind of input object is input. Then, it selects the indicator data according to the specs supplied.

Usage

```r
getIn(
  obj,
  dset = "Raw",
  icodes = NULL,
  aglev = NULL,
  usel = NULL,
  use_group = NULL,
  justnumeric = TRUE
)
```

Arguments

- **obj** An input object. The function can handle either the COIN object, or a data frame. The data frame should have each column as an indicator, and optional columns `UnitCode` and `UnitName` which specify the code (or name) of each unit. Any columns except these latter two will be treated as indicators. Any other type of object will return an error.
- **dset** If input is a COIN object, this specifies which data set in `.Data` to use.
- **icodes** An optional character vector of indicator codes to subset the indicator data. Usefully, can also refer to an aggregation group name, and data will be subsetted accordingly. NOTE does not work with multiple aggregate group names.
- **aglev** The aggregation level to take indicator data from. Integer from 1 (indicator level) to N (top aggregation level, typically the index).
getIn

use1  An optional unit code, or character vector of unit codes to use to filter the data. The returned data will only include rows corresponding to the use1, unless use1 = NULL (default).

use_group  An optional grouping variable and group to filter data from. Of the format list(Group_Var = Group), where Group_Var is a Group_ column that must be present in the selected data set, and Group is a specified group inside that grouping variable. This filters the selected data to only include rows from the specified group. Alternatively, this argument can work in conjunction with use1: if use1 is specified, use_group may be input as a string simply representing a group column. In that case the data will be filtered to include only rows from the group(s) which the use1 belong to. If use1 is specified and use_group is specified as a list, use1 will take precedence and use_group will be ignored.

justnumeric  Logical: if TRUE, removes any non-numeric columns from ind_data_only. Otherwise keeps all except those.

Details

For example, specifying dset = "Raw" and icodes = c("Ind1", "Ind5"), it will return the indicator columns named "Ind1" and "Ind5" (if they exist), in the format described below. icodes can be indicators or aggregation groups, and can call multiple groups.

You can also specify which aggregation level to target, using the aglev argument. See examples below, and in particular the COINr online documentation.

As well as selection of columns, you can also filter to specific rows using unit codes as a reference, and/or grouping variables. This is done using the use1 and use_group arguments. This gives the ability to isolate a unit inside a given group, for example.

getIn() is used by many COINr functions for plotting, accessing and reporting subsets of indicator data.

Value

A list with the following entries:

- .$IndCodes The indicator codes
- .$IndNames The indicator names (if a COIN object is input)
- .$ind_data A data frame of indicator data, according to the input specifications, including any unit codes, names and groups
- .$ind_data_only A data frame, as above, but without unit codes, names, groups.
- .$UnitCodes Unit codes of selected data set.
- .$otype Object type (a string: either "COINobj" or "df").

Examples

# assemble ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# get indicator data from Social pillar
SocData <- getIn(ASEM, dset = "Raw", icodes = "Social", aglev = 1)
# Indicator codes
getPCA

Perform PCA on a COIN

Description

Performs Principle Component Analysis (PCA) on a specified data set and subset of indicators or aggregation groups. Returns weights corresponding to the first principal component, i.e. the weights that maximise the variance explained by the linear combination of indicators.

Usage

getPCA(
  COIN,
  dset = "Raw",
  icodes = NULL,
  aglev = NULL,
  by_groups = TRUE,
  nowarnings = FALSE,
  out2 = "COIN"
)

Arguments

COIN An input object. The function can handle either the COIN object, or a data frame. The data frame should have each column as an indicator, and an optional column "UnitCode" which specifies the code (or name) of each unit. Any other type of object will return an error.

dset If input is a COIN object, this specifies which data set in .$Data to use.

icodes An optional character vector of indicator codes to subset the indicator data. Usefully, can also refer to an aggregation group name, and data will be sub-setted accordingly. NOTE does not work with multiple aggregate group names.

aglev The aggregation level to take indicator data from. Integer from 1 (indicator level) to N (top aggregation level, typically the index).

by_groups If TRUE (default), performs PCA inside each aggregation group inside the specified level. If FALSE, performs a single PCA over all indicators/aggregates in the specified level.

nowarnings If FALSE (default), will give warnings where missing data are found. Set to TRUE to suppress these warnings.

out2 If the input is a COIN object, this controls where to send the output. If "COIN", it sends the results to the COIN object, otherwise if "list", outputs to a separate list.
getResults

Details

Note that getPCA() is simply a quick wrapper for stats::prcomp() which makes PCA on COINs quicker. See COINr online documentation for more details and examples.

Value

If out2 = "COIN", results are appended to the COIN object. Specifically:

- A new set of PCA weights is added to $Parameters$Weights
- A list is added to $Analysis$ containing PCA weights (loadings) of the first principle component, and the output of stats::prcomp, for each aggregation group found in the targeted level. If out2 = "list" the same outputs are contained in a list.

See Also

- stats::prcomp Principle component analysis

Examples

```r
# build ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta,
                 AggMeta = ASEMAggMeta)
# get PCA results for pillar groups inside "Sust" (sustainability) sub-index
# (warnings about missing data are suppressed)
PCAres <- getPCA(ASEM, dset = "Raw", icodes = "Sust",
                 aglev = 1, nowarnings = TRUE, out2 = "list")
# summarise PCA results for Social pillar
summary(PCAres$PCAresults$Social$PCAres)
```

Description

Generates fast results tables, either attached to the COIN or as a data frame.

Usage

```r
getResults(  COIN,
              tab_type = "Summ",
              use = "scores",
              order_by = NULL,
              nround = 2,
              use_group = NULL,
              out2 = "df")
```
Arguments

COIN  The COIN object, or a data frame of indicator data

tab_type  The type of table to generate. Either "Summ" (a single indicator plus rank), "Aggs" (all aggregated scores/ranks above indicator level), or "Full" (all scores/ranks plus all group, denominator columns).

use  Either "scores" (default), "ranks", or "groupranks". For the latter, use_group must be specified.

order_by  A code of the indicator or aggregate to sort the table by. If not specified, defaults to the highest aggregate level, i.e. the index in most cases. If use_group is specified, rows will also be sorted by the specified group.

nround  The number of decimal places to round numerical values to. Defaults to 2.

use_group  An optional grouping variable. If specified, the results table includes this group column, and if use = "groupranks", ranks will be returned with respect to the groups in this column.

out2  If "df", outputs a data frame (tibble). Else if "COIN" attaches to .$Results in an updated COIN.

Details

Although results are available in a COIN in .$Data, the format makes it difficult to quickly present results. This function generates results tables that are suitable for immediate presentation, i.e. sorted by index or other indicators, and only including relevant columns. Scores are also rounded by default, and there is the option to present scores or ranks.

Value

If out2 = "df", the results table is returned as a data frame. If out2 = "COIN", this function returns an updated COIN with the results table attached to .$Results.

See Also

• resultsDash() Interactive results dashboard
• coin2Excel() Export results to Excel

Examples

# build ASEM COIN up to aggregation
ASEM <- build_ASEM()

# results table of scores for index and aggregates (excluding indicator scores)
dfResults <- getResults(ASEM, tab_type = "Aggregates", out2 = "df")
getStats

Get table of indicator statistics for any data set

Description

Takes a COIN or data frame and returns a table of statistics for each column, including max, min, median, mean, standard deviation, kurtosis, etc. Flags indicators with possible outliers, and checks for collinearity with other indicators and denominators (if any). Also checks number of unique values and percentage of zeros. Also returns correlation matrices and a table of outliers, as a list.

Usage

getStats(
  COIN,
  icodes = NULL,
  dset = "Raw",
  out2 = "COIN",
  cortype = "pearson",
  t_skew = 2,
  t_kurt = 3.5,
  t_colin = 0.9,
  t_denom = 0.7,
  t_missing = 65,
  IQR_coef = 1.5
)

Arguments

COIN A COIN object or data frame of indicator data
icodes A character vector of indicator names to analyse. Defaults to all indicators.
dset The data set to analyse.
out2 Where to output the results: if "COIN" (default), appends to the COIN, otherwise if "list", outputs to a separate list.
cortype The type of correlation to calculate, either "pearson", "spearman", or "kendall". See stats::cor.
t_skew Skewness threshold.
t_kurt Kurtosis threshold.
t_colin Collinearity threshold (absolute value of correlation).
t_denom High correlation with denominator threshold.
t_missing Missing data threshold, in percent.
IQR_coef Interquartile range coefficient, used for identifying outliers.
getStrengthNWeak

Value
If out2 = "COIN" (default), results are appended to the COIN in .$Analysis, otherwise if out2 = "list", outputs to a separate list. In both cases, the result is a list containing:

- A data frame of statistics for each indicator column
- A data frame indicating which points may be considered outliers according to the interquartile range
- A data frame of correlations between indicators
- A data frame of correlations between indicators and any denominators present in .$Input$Denominators

Examples

```r
# build ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# get list of stats from raw data set
stat_list <- getStats(ASEM, dset = "Raw", out2 = "list")
```

---

**getStrengthNWeak**

*Generate strengths and weaknesses for a specified unit*

Description

Generates a table of strengths and weaknesses for a selected unit, based on ranks, or ranks within a specified grouping variable.

Usage

```r
getStrengthNWeak(
  COIN,
  dset = NULL,
  usel = NULL,
  topN = 5,
  bottomN = 5,
  withcodes = TRUE,
  use_group = NULL,
  unq_discard = NULL,
  min_discard = TRUE,
  report_level = NULL,
  with_units = TRUE,
  adjust_direction = NULL)
```
Arguments

- **COIN**: A COIN
- **dset**: The data set to extract indicator data from, to use as strengths and weaknesses.
- **use_l**: A selected unit code
- **topN**: The top N indicators to report
- **bottomN**: The bottom N indicators to report
- **withcodes**: If TRUE (default), also includes a column of indicator codes. Setting to FALSE may be more useful in generating reports, where codes are not helpful.
- **use_group**: An optional grouping variable (named column of .Data$Aggregated) to use for reporting in-group ranks. Specifying this will report the ranks of the selected unit within the group of use_group to which it belongs.
- **unq_discard**: Optional parameter for handling discrete indicators. Some indicators may be binary variables of the type "yes = 1", "no = 0". These may be picked up as strengths or weaknesses, when they may not be wanted to be highlighted, since e.g. maybe half of units will have a zero or a one. This argument takes a number between 0 and 1 specifying a unique value threshold for ignoring indicators as strengths. E.g. setting prc_unq_discard = 0.2 will ensure that only indicators with at least 20% unique values will be highlighted as strengths or weaknesses. Set to NULL to disable (default).
- **min_discard**: If TRUE (default), discards any strengths which correspond to the minimum rank for the given indicator. See details.
- **report_level**: Optional aggregation level to report parent codes from. For example, setting report_level = 2 will add a column to the strengths and weaknesses tables which reports the aggregation group from level 2, to which each reported indicator belongs. Set to NULL to disable (default).
- **with_units**: If TRUE (default), includes indicator units in output tables.
- **adjust_direction**: If TRUE, will adjust directions of indicators according to the "Direction" column of IndMeta. By default, this is TRUE if dset = "Raw", and FALSE otherwise.

Details

This currently only works at the indicator level. Indicators with NA values for the selected unit are ignored. Strengths and weaknesses mean the top N-ranked indicators for the selected unit. Effectively, this takes the rank that the selected unit has in each indicator, sorts the ranks, and takes the top N highest and lowest.

This function must be used with a little care: indicators should be adjusted for their directions before use, otherwise a weakness might be counted as a strength, and vice versa. Use the adjust_direction parameter to help here.

A further useful parameter is unq_discard, which also filters out any indicators with a low number of unique values, based on a specified threshold. Also min_discard which filters out any indicators which have the minimum rank.

The best way to use this function is to play around with the settings a little bit. The reason being that in practice, indicators have very different distributions and these can sometimes lead to unexpected
outcomes. An example is if you have an indicator with 50% zero values, and the rest non-zero (but unique). Using the sport ranking system, all units with zero values will receive a rank which is equal to the number of units divided by two. This then might be counted as a "strength" for some units with overall low scores. But a zero value can hardly be called a strength. This is where the min_discard function can help out.

Problems such as these mainly arise when e.g. generating a large number of country profiles.

Value

A list containing a data frame .$Strengths, and a data frame .$Weaknesses. Each data frame has columns with indicator code, name, rank and value (for the selected unit).

See Also

- getUnitReport() Automatic unit report as html, pdf or Word
- getUnitSummary() Summary of scores for a given unit

Examples

```r
# build ASEM COIN up to aggregation
ASEM <- build_ASEM()
# generate top 5 strengths and weaknesses for GBR
getStrengthNWeak(ASEM, usel = "GBR")
```

---

### Description

Generates a scorecard for a given unit using an R Markdown template.

### Usage

```r
getUnitReport(
  COIN,
  usel,
  out_type = ".html",
  outdir = NULL,
  rmd_template = NULL
)
```
**getUnitReport**

**Arguments**

- **COIN**  
  A COIN

- **usel**  
  A selected unit code, or a character vector of unit codes (for multiple reports).

- **out_type**  
  A string specifying the output type. Can be either ".docx" (Word), ".pdf" or ".html". IMPORTANT: if the template includes interactive plots (e.g. the iplot() functions from COINr), writing to .docx or .pdf will not work unless you have installed the webshot package. To do this, run: `install.packages("webshot")`
  `webshot::install_phantomjs()`

- **outdir**  
  Character string specifying the output directory (defaults to current working directory).

- **rmd_template**  
  A character string specifying the full file path to an R Markdown template which is used to generate the report. If this is not specified, defaults to COINr’s inbuilt template example.

**Details**

Most likely you will want to customise the template which can be found in the COINr installed package directory under /UnitReport. Currently, a few examples are given, such as some charts and basic summary statistics.

This function will render the unit report to either pdf, html or word doc. As mentioned below, if you have HTML widgets such as interactive plotly plots, or COINr iplot() functions, you will need to install the webshot package to be able to render to pdf or word formats.

To customise the template, copy the .rmd template found in /UnitReport and alter it, then point the rmd_template argument to your new template.

Note that this function is particularly useful for generating a large number of reports, e.g. we can generate reports for all units at once using a for loop, purrr::map or apply() or similar.

**Value**

Markdown document rendered to HTML, pdf or Word. This function requires Pandoc to be installed. If Pandoc is not found, then it returns a warning and a printed message (string).

**Examples**

```
# build ASEM COIN up to aggregation
ASEM <- build_ASEM()
# Generate a unit report for NZ
# This is written to the temporary directory to avoid polluting other directories
# during automated testing.
# It will be deleted at the end of the R session.
# Normally you would set the directory to somewhere else to save the resulting files
getUnitReport(ASEM, usel = "NZL", out_type = ".html", outdir = tempdir())
# You can find this file in the temporary directory:
print(tempdir())
# We will now delete the file to keep things tidy in testing
unlink(paste0(tempdir(),"\NZL_report.html"))
```
getUnitSummary  Generate unit summary table

Description
Generates a summary table for a single unit. This is mostly useful in unit reports.

Usage
getUnitSummary(COIN, usel, aglevs)

Arguments
- COIN  A COIN
- usel  A selected unit code
- aglevs  The aggregation levels to display results from.

Details
This returns the scores and ranks for each indicator/aggregate as specified in aglevs. It orders the table so that the highest aggregation levels are first. This means that if the index level is included, it will be first.

Value
A summary table as a data frame, containing scores and ranks for specified indicators/aggregates.

See Also
- `getUnitReport()` Automatic unit report as html, pdf or Word
- `getStrengthNWeak()` Top N-ranking indicators for a given unit

Examples
```r
# build ASEM COIN up to aggregation
ASEM <- build_ASEM()
# generate unit summary for NZ - index and sub-indexes only
getUnitSummary(ASEM, usel ="NZL", aglevs = c(4,3))
```
**Description**

Weighted harmonic mean of a vector. NA are skipped by default. This function is used inside `aggregate()`.

**Usage**

```
harMean(x, w = NULL)
```

**Arguments**

- `x` A numeric vector of positive values.
- `w` A vector of weights, which should have length equal to `length(x)`. Weights are relative and will be re-scaled to sum to 1. If `w` is not specified, defaults to equal weights.

**Value**

Weighted harmonic mean, as a numeric value.

**Examples**

```r
# a vector of values
x <- 1:10
# a vector of weights
w <- runif(10)
# weighted harmonic mean
harMean(x, w)
```

**Description**

This returns a data frame of any highly correlated indicators within the same aggregation group. The level of the aggregation group can be controlled by the `grouplev` argument.
Usage

hicorrSP(
  COIN,
  dset = "Normalised",
  hicorval = 0.9,
  cortype = "pearson",
  grouplev = NULL
)

Arguments

COIN                  Data frame with first col indicator codes, second is weights, third is correlations
dset                  The data set to use for correlations.
hicorval              A threshold to flag high correlation. Default 0.9.
cortype               The type of correlation, either "pearson" (default), "spearman" or "kendall". See stats::cor.
grouplev              The level to group indicators in. E.g. if grouplev = 2 it will look for high correlations between indicators that belong to the same group in Level 2.

Value

A data frame with one entry for every indicator pair that is highly correlated within the same group, at the specified level. Pairs are only reported once, i.e. only uses the upper triangle of the correlation matrix.

See Also

• rew8r() Interactive app for adjusting weights and seeing effects on correlations
  • getCorr() Get correlations between indicators/levels

Examples

# Assemble ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# check for any within-pillar correlations of > 0.7
hicorrSP(ASEM, dset = "Raw", hicorval = 0.7, , grouplev = 2)

Description

Imputation of missing data data sets using a variety of methods (see imtype). This also includes the possibility of imputing by grouping variables, i.e. columns of IndData that are prefaced by "Group_".
Usage

```r
impute(
  COIN,
  imtype = NULL,
  dset = NULL,
  groupvar = NULL,
  EMaglev = NULL,
  out2 = "COIN"
)
```

Arguments

- **COIN**: A COIN or a data frame
- **imtype**: The type of imputation method. Either:
  - "agg_mean" (the mean of normalised indicators inside the aggregation group),
  - "agg_median" (the median of normalised indicators inside the aggregation group),
  - "ind_mean" (the mean of all the other units in the indicator),
  - "ind_median" (the median of all the other units in the indicator),
  - "indgroup_mean" (the mean of all the other units in the indicator, in the same group),
  - "indgroup_median" (the median of all the other units in the indicator, in the same group),
  - "EM" (expectation maximisation algorithm via AMELIA package, currently without bootstrapping)
  - "none" (no imputation, returns original data set)
- **dset**: The data set in .$Data to impute
- **groupvar**: The name of the column to use for by-group imputation. Only applies when imtype is set to a group option.
- **EMaglev**: The aggregation level to use if imtype = "EM".
- **out2**: Where to output the imputed data frame. If "COIN" (default for COIN input), creates a new data set .$Data$Imputed. Otherwise if "df" outputs directly to a data frame.

Details

See online documentation for further details and examples.

Value

If out2 = "COIN" (default for COIN input), creates a new data set .$Data$Imputed. Otherwise if out2 = "df" outputs directly to a data frame.
indChange

Add and remove indicators

Description

A shortcut function to add and remove indicators. This will make the relevant changes and recalculate the index if asked. Adding and removing is done relative to the current set of indicators used in calculating the index results. Any indicators that are added must of course be present in .$Input$Original (in both IndData and IndMeta).

Usage

indChange(COIN, add = NULL, drop = NULL, regen = FALSE)

Arguments

<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>COIN</td>
<td>COIN object</td>
</tr>
<tr>
<td>add</td>
<td>A character vector of indicator codes to add (must be present in the original input data)</td>
</tr>
<tr>
<td>drop</td>
<td>A character vector of indicator codes to remove (must be present in the original input data)</td>
</tr>
<tr>
<td>regen</td>
<td>Logical (default): if TRUE, automatically regenerates the results based on the new specs Otherwise, just updates the .$Method$assemble parameters. This latter might be useful if you want to Make other changes before re-running using the regen() function.</td>
</tr>
</tbody>
</table>

Value

An updated COIN, with regenerated results if regen = TRUE.
See Also

- `regen()` regenerate a COIN
- `compTable()` compare two different COINs
- `compTableMulti()` compare multiple COINs

Examples

```r
# build ASEM example
ASEM <- build_ASEM()
# remove one indicator and regenerate results
ASEM2 <- indChange(ASEM, drop = "UNVote", regen = TRUE)
# compare the differences
CT <- compTable(ASEM, ASEM2, dset = "Aggregated", isel = "Index")
```

Description

Generates an interactive visualisation of one or two indicators at a time. Requires Shiny and an active R session. This dashboard is useful for quickly exploring indicator data, and seeing e.g. an untreated indicator distribution against its treated equivalent.

Usage

```r
indDash(COIN)
```

Arguments

- **COIN** The COIN object

Details

This function requires an interactive R session. Otherwise it will simply generate a message to that effect.

Value

Interactive app (if running an interactive R session). This app is purely exploratory and does not return anything back to R.
Examples

```r
# Only run in interactive mode
if(interactive()){
  # build ASEM COIN
  ASEM <- build_ASEM()
  # view dashboard
  indDash(ASEM)
}
```

**iplotBar**

*Bar chart*

**Description**

Generates an interactive bar chart. This function is simply a wrapper for the `plotly` bar chart function, but accesses the COIN object to get the relevant indicator. Also has click event data for Shiny. Allows construction of stacked bar charts which show underlying components (for aggregated data only), and plots of only specified groups.

**Usage**

```r
iplotBar(
  COIN,
  dset = "Raw",
  isel = NULL,
  usel = NULL,
  aglev = NULL,
  stack_children = FALSE,
  from_group = NULL
)
```

**Arguments**

- **COIN**
  The COIN object, or a data frame of indicator data.
- **dset**
  The data set to plot.
- **isel**
  The selected indicator code or aggregate (does not support multiple indicators)
- **usel**
  A character vector of unit codes to highlight on the bar chart (optional)
- **aglev**
  The aggregation level to collect the indicator data from (this needs to be specified)
- **stack_children**
  If TRUE, produces a stacked bar chart with any children of isel. In this case, usel is ignored. This only works if dset = "Aggregated" and aglev > 1.
- **from_group**
  Filters the bar chart to a specified group using a group column that is present in the specified data set. Specified as list(group_variable = selected_group).
**iplotCorr**

### Value
Interactive bar chart generated by plotly.

### See Also
- `iplotMap()` bar chart of indicator or aggregate
- `resultsDash()` interactive dashboard of indicator data

### Examples
```r
# assemble ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# plot Flights indicator
iplotBar(ASEM, dset = "Raw", isel = "Flights", aglev = 1)
```

---

### iplotCorr

#### Correlation heatmap

### Description
Plots an interactive heatmap of a correlation matrix. Currently this only works with the aggregated data set, i.e. you need to have aggregated the data first before using this.

### Usage
```
iplotCorr(
  COIN,  # The COIN object
  aglevs = NULL,  # A two length vector specifying which level to plot against which level. E.g. c(2,4) for COIN plots sub-pillars against sub-indexes. If NULL, plots everything against everything.
  insig = FALSE,  # Logical: if TRUE, all correlation values are plotted; if FALSE (default), does not plot insignificant correlations.
  lev = TRUE,  
  grouprects = TRUE,  
  flagcolours = TRUE,  
  corthresh = NULL,  
  showvals = TRUE,  
  cortype = "pearson",  
  useweights = NULL
)
```

### Arguments
- **COIN**
  - The COIN object
- **aglevs**
  - A two length vector specifying which level to plot against which level. E.g. c(2,4) for COIN plots sub-pillars against sub-indexes. If NULL, plots everything against everything.
- **insig**
  - Logical: if TRUE, all correlation values are plotted; if FALSE (default), does not plot insignificant correlations.
<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>levs</td>
<td>Logical: if TRUE, plots lines showing the division between different levels. Only works if aglevs = NULL.</td>
</tr>
<tr>
<td>grouprects</td>
<td>Logical: if TRUE, plots rectangles showing aggregation groups.</td>
</tr>
<tr>
<td>flagcolours</td>
<td>If TRUE uses a discrete colour scale specified by corthresh. Otherwise uses a continuous colour map.</td>
</tr>
<tr>
<td>corthresh</td>
<td>A named list specifying the colour thresholds to use if flagcolours = TRUE. Entries should specify correlation thresholds and can specify any of clow, cmid and chi. Anything below clow will be coloured red. Anything between clow and cmid will be grey. Anything between cmid and chigh will be blue. Anything above chigh will be green. Default is list(clow = -0.4, cmid = 0.4, chigh = 0.85). You can specify a subset of these and the others will revert to defaults.</td>
</tr>
<tr>
<td>showvals</td>
<td>Logical: if TRUE, overlays correlation values on each square.</td>
</tr>
<tr>
<td>cortype</td>
<td>The type of correlation: either &quot;pearson&quot; (default), &quot;spearman&quot; or &quot;kendall&quot;. See stats::cor.</td>
</tr>
<tr>
<td>useweights</td>
<td>An optional list of weights to use (this is used mainly in the rew8r() app).</td>
</tr>
</tbody>
</table>

**Details**

This is a slightly involved wrapper for plotly. It allows plotting any level against any other, and outputs correlation heat maps as HTML widgets. It has some flexibility regarding grouping of indicators, colouring, treatment of insignificant correlations, and the correlation type. Explore the arguments and see.

**Value**

A plotly correlation map (figure).

**See Also**

- plotCorr() Static correlation heat maps
- rew8r() Interactive app for adjusting weights and seeing effects on correlations
- getCorr() Get correlations between indicators/levels

**Examples**

```r
# build ASEM COIN up to aggregation
ASEM <- build_ASEM()

# correlation heatmap of pillars against sub-indexes
iplotCorr(ASEM, aglevs = c(2,3))
```
iplotIndDist  

Interactive indicator distribution plots

Description

Generates a JavaScript distribution plot of a single indicator, using plotly. Plot can be embedded e.g. in HTML documents, websites, etc, or used for more interactive data exploration. This only plots one indicator at a time - for multiple plots you can use plotIndDist().

Usage

iplotIndDist(
  COIN,
  dset = "Raw",
  icodes = NULL,
  ptype = "Violin",
  aglev = 1,
  axlims = NULL
)

Arguments

COIN The COIN object, or a data frame of indicator data  
dset The source data set to use for indicator data (if input is COIN object)  
icodes A character vector of a single indicator name or aggregate name to plot.  
ptype The type of plot to produce. Currently supports "Violin" and "Histogram".  
aglev The aggregation level to extract the indicator data from. Defaults to indicator level (1)  
axlims Optional parameter specifying axis limits. Useful mainly for matching with another plot.

Value

Plots generated with plotly. These can be edited further with plotly commands.

Examples

# build ASEM COIN  
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)  
# plot renewable energy indicator  
iplotIndDist(ASEM, "Raw", "Renew", ptype = "Violin")
iplotIndDist2  

Interactive indicator distribution plots for two indicators simultaneously

Description

Generates a JavaScript distribution plot of two indicators, using Plotly. Plot can be embedded e.g. in HTML documents, websites, etc, or used for more interactive data exploration.

Usage

```r
iplotIndDist2(
  COIN,
  dsets = "Raw",
  icodes = NULL,
  ptype = "Scatter",
  aglevs = 1
)
```

Arguments

- **COIN**: The COIN, or a data frame of indicator data
- **dsets**: The source data sets to use for indicator data (if input is COIN object). If the source data sets are the same, this can be a single character string, otherwise, a character vector, e.g. `c("Raw","Treated")`.
- **icodes**: A character vector of two indicator codes to plot (corresponding to the two dsets specified)
- **ptype**: The type of plot to produce. Currently supports "Histogram" and "Scatter".
- **aglevs**: The aggregation level to extract the indicator data from. Defaults to indicator level (1). This also can be specified as a vector if the two indicators are from different levels.

Value

Plots generated with `plotly`. These can be edited further with `plotly` commands.

Examples

```r
# build ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# plot CO2 against renewable energy indicator
iplotIndDist2(ASEM, dsets = "Raw", icodes = c("Renew", "CO2"))
```
**Description**

Generates an interactive choropleth map of specified indicator data. Only works on national level data (i.e. one point per country), with countries labelled by ISO alpha-3 codes. This function is simply a quick wrapper for the `plotly` choropleth map function.

**Usage**

```r
iplotMap(COIN, dset = "Raw", isel)
```

**Arguments**

- **COIN**: The COIN object, or a data frame of indicator data.
- **dset**: The data set to plot.
- **isel**: The selected indicator code or aggregate

**Value**

Interactive map generated by plotly.

**See Also**

- `iplotBar()` bar chart of indicator or aggregate
- `resultsDash()` interactive dashboard of indicator data

**Examples**

```r
# assemble ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# map CO2 indicator
iplotMap(ASEM, dset = "Raw", isel = "CO2")
```
iplotRadar  

Radar chart

Description

Generates an interactive radar chart for a specified unit or set of units.

Usage

iplotRadar(
  COIN,
  dset = "Raw",
  usel = NULL,
  aglev = NULL,
  isel = NULL,
  addstat = "none",
  statgroup = NULL,
  statgroup_name = NULL
)

Arguments

COIN  
The COIN object, or a data frame of indicator data.

Dset  
The data set to use in the table

usel  
Character vector of unit code(s) to plot data from

aglev  
The selected aggregation level to take indicator data from, where 1 is the base indicator level, and 2, 3 etc are higher aggregation levels

isel  
The indicator or aggregation code(s) to plot

addstat  
Adds the statistic of the scores in each dimension as a separate trace. If "mean" adds the overall mean for each dimension/indicator. If "median" adds the overall median. If "groupmean" or "groupmedian", adds the group mean or median respectively of the first unit specified in usel, using the group specified by statgroup. Default "none", i.e. no extra trace. Using a group mean or median won’t make sense unless all of selected units are from the same group.

statgroup  
A grouping variable (must be present in dset) if addstat = "groupmean" or "groupmedian"

statgroup_name  
An optional name to display for statgroup. In the legend this will appear as e.g. "statgroup_name group mean". Defaults to statgroup.

Details

This function uses plotly to generate a radar chart for showing one or more units, compared using a specified set of indicators. Optionally, you can add mean/median or group mean/median as an extra trace. The point being to show how a particular unit compares to its peers.
Value

Interactive radar chart generated using plotly.

See Also

- resultsDash() Interactive results dashboard.

Examples

# build ASEM COIN up to aggregation
ASEM <- build_ASEM()
# radar chart of Austria vs. China in Political indicators
iplotRadar(ASEM, dset = "Aggregated", usel = c("AUT", "CHN"), isel = "Political", aglev = 1)

iplotTable

Results table

Description

Generates an interactive table of data. For use in e.g. Shiny or HTML documents.

Usage

iplotTable(
  COIN,
  dset = "Raw",
  isel = NULL,
  aglev = NULL,
  nround = 1,
  extra_cols = FALSE
)

Arguments

- **COIN**: The COIN object, or a data frame of indicator data.
- **dset**: The data set to use in the table
- **isel**: The selected indicator codes (default all in dset)
- **aglev**: The aggregation level to select data from
- **nround**: The number of decimal places to round scores to (default 1).
- **extra_cols**: If FALSE (default), excludes group columns and similar. If TRUE, includes them.

Details

This function is a wrapper for the reactable package and offers a fast way to make interactive tables. It also applies conditional formatting (colouring by cell value), and sorts by the first column by default. Like other COINr functions, it can target subsets of indicators.
is.COIN

Value

An interactive table generated by reactable.

See Also

- `colourTable()` Conditionally-formatted table for any data frame
- `resultsDash()` interactive dashboard of indicator data
- `getResults()` results summary tables

Examples

# assemble ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# table of indicators in "Political" pillar
iplotTable(ASEM, dset = "Raw", isel = "Political", aglev = 1)

is.COIN

Check if an object is a COIN

Description

Returns TRUE if an input object is a COIN, otherwise FALSE if not.

Usage

is.COIN(obj)

Arguments

obj

An input object to test

Value

Logical: TRUE if input is a COIN, otherwise FALSE

See Also

- `getIn()` Get subset of indicator data from either a COIN or data frame.
- `assemble()` Assemble a COIN from indicator data and metadata

Examples

# build the ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# check class
stopifnot(is.COIN(ASEM))
Description

This applies various simple transformations, to be used by the treat() function. This function is probably not very useful on its own because it requires params, a list of parameters which are used to output the type and status of treatment applied.

Usage

loggish(x, ltype, params = NULL)

Arguments

x A vector or column of data to transform
ltype The type of log transformation - see deflog in treat().
params Some extra parameters to pass. These parameters mostly concern internal messages for treat() and this can be left unspecified unless ltype == "boxcox", in which case there should be a parameter params$boxlam specified (see BoxCox()). However, if you wish to use a Box Cox transformation, it is better to use BoxCox() directly.

Value

A list with

- .$x is the transformed vector of data
- .$Flag is a flag of the type of treatment specified (used inside treat())
- .$Treatment the treatment applied (used inside treat())
- .$TreatSpec the treatment specified (used inside treat())

See Also

- treat() Outlier treatment

Examples

# get a column of data with outliers
x <- ASEMIndData$Tariff
# apply a GII transformation
xdash <- loggish(x, ltype = "GIIlog")
# plot one against the other
plot(x, xdash$x)
names2Codes

Generate short codes from long names

Description

Given a character vector of long names (probably with spaces), generates short codes. Intended for use when importing from the COIN Tool.

Usage

names2Codes(cvec, maxword = 2, maxlet = 4)

Arguments

cvec
A character vector of names
maxword
The maximum number of words to use in building a short name (default 2)
maxlet
The number of letters to take from each word (default 4)

Value

A corresponding character vector, but with short codes, and no duplicates.

See Also

• COINToolIn() Import data from the COIN Tool (Excel).

Examples

# generate codes for indicators in the ASEM data set (first five only)
names2Codes(ASEMIndMeta$IndName[1:5], maxlet = 3)

noisyWeights

Noisy replications of weights

Description

Given a set of weights, this function returns multiple replicates of the weights, with added noise. This is intended for use in uncertainty and sensitivity analysis.

Usage

noisyWeights(w, noise_specs, Nrep)
normalise

Arguments

- **w**: A data frame of weights, in the format found in `.Parameters$Weights`.
- **noise_specs**: a data frame with columns:
  - AgLevel: The aggregation level to apply noise to
  - NoiseFactor: The size of the perturbation: setting e.g. 0.2 perturbs by +/- 20% of nominal values.
- **Nrep**: The number of weight replications to generate.

Details

Weights are expected to be in a long data frame format with columns Aglevel, Code and Weight, as used inside COINs.

Noise is added using the `noise_specs` argument, which is specified by a data frame with columns AgLevel and NoiseFactor. The aggregation level refers to number of the aggregation level to target while the NoiseFactor refers to the size of the perturbation. If e.g. a row is AgLevel = 1 and NoiseFactor = 0.2, this will allow the weights in aggregation level 1 to deviate by +/- 20% of their nominal values (the values in `w`).

Value

A list of `Nrep` sets of weights (data frames).

See Also

- `sensitivity()` Perform global sensitivity or uncertainty analysis on a COIN

Examples

```r
# build ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)

# generate 2 sets of weights based on original ASEM weights, # perturbed by +/-20% only at indicator level
wlist <- noisyWeights(ASEM$Parameters$Weights$Original, noise_specs = data.frame(AgLevel = 1, NoiseFactor = 0.2), Nrep = 2)
```

---

**Normalise indicator data sets**

**Description**

A dataset of indicators is normalised using one of several methods. This function also supports custom normalisation.
Usage

normalise(
  COIN,
  ntype = "minmax",
  npara = NULL,
  dset = NULL,
  directions = NULL,
  individual = NULL,
  indiv_only = NULL,
  out2 = NULL
)

Arguments

COIN Either the COIN object, or a data frame of indicator data
ntype The type of normalisation method. Either "minmax", "zscore", "scaled", "goalposts", "rank", "borda", "prank", "fracmax", "dist2targ", "dist2ref", "dist2max", "custom" or "none". See the online documentation.
npara Supporting object for ntype. This is a list of the form list(ntype = parameters_for_ntype). So, if ntype = "minmax", npara could be list(minmax = c(0,100)) to scale into the 0 to 100 interval. If ntype = "zscore", npara could be list(zscore = c(0,1)) to scale to mean zero and standard deviation 1. This means you can store parameters for more than one normalisation type side by side, which helps in comparisons, adjustments, and sensitivity analyses.
dset The data set to normalise.
directions A vector specifying the direction assigned to each indicator. Needs to be the same length as the number of indicators, or the number of indicators in icodes, if specified.
individual A list of named lists specifying individual normalisation to apply to specific indicators. Should be structured as follows: The name of each sub-list should be the indicator code. The the list elements are:
  • .$ntype is the type of normalisation to apply
  • .$npara is a corresponding object or parameters that are used by ntype, in the same format as npara above.
indiv_only Logical: if FALSE (default), indicators not specified in individual are subjected to default normalisation. Otherwise if TRUE they are not normalised.
out2 Where to output the results. If "COIN" (default for COIN input), appends to updated COIN, otherwise if "df" outputs to data frame.

details

Normalisation refers to the operation of bringing variables (indicators) onto a common scale. This is typically done by matching one or more indicator statistics. For example, the min-max method operates a linear transformation to make the minimum and maximum values of each indicator to be equal. The z-score method makes the standard deviation and variance equal. And so on.
This function supports a range of normalisation methods - see ntype. Some of these require supporting parameters or similar - to see full details check the online documentation.

Indicators can also be each normalised by a different method. See individual.

Value

If out2 = "COIN" (default for COIN input), returns an updated COIN object with a data frame $Data$Normalised added. Else if out2 = "df" outputs a data frame of normalised data.

Examples

```r
# build ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# directly normalise raw data using min-max, onto 0-10 interval
ASEM <- normalise(ASEM, dset = "Raw", ntype = "minmax", npara = list(minmax = c(0,10)))
# Check: get indicator data first
NormData <- getIn(ASEM, dset = "Normalised")$ind_data_only
# ensure that min is 0 and max is 10 for all columns
stopifnot(
  all(apply(NormData, MARGIN = 2, min, na.rm = TRUE) == 0),
  all(apply(NormData, MARGIN = 2, max, na.rm = TRUE) == 10)
)
```

outrankMatrix

```
outrankMatrix  Outranking matrix

Description

Constructs an outranking matrix based on a data frame of indicator data and corresponding weights. This function is used inside aggregate().

Usage

```
outrankMatrix(ind_data, w = NULL)
```

Arguments

```
ind_data  A data frame or matrix of indicator data, with observations as rows and indicators as columns. No other columns should be present (e.g. label columns).

w         A vector of weights, which should have length equal to ncol(ind_data). Weights are relative and will be re-scaled to sum to 1. If w is not specified, defaults to equal weights.
```
Value

A list with:

- `.OutRankMatrix` the outranking matrix with `nrow(ind_data)` rows and columns (matrix class).
- `.nDominant` the number of dominance/robust pairs
- `.fracDominant` the percentage of dominance/robust pairs

Examples

```r
# get a sample of a few indicators
ind_data <- COINr::ASEMIndData[12:16]
# calculate outranking matrix
outlist <- outrankMatrix(ind_data)
# see fraction of dominant pairs (robustness)
outlist$fracDominant
```

plotCorr

*Static heatmaps of correlation matrices*

Description

Generates heatmaps of correlation matrices using `ggplot2`. This enables correlating any set of indicators against any other, and supports calling named aggregation groups of indicators. The `withparent` argument generates tables of correlations only with parents of each indicator. Also supports discrete colour maps using `flagcolours`, different types of correlation, and groups plots by higher aggregation levels.

Usage

```r
plotCorr(
  COIN,
  dset = "Raw",
  icodes = NULL,
  aglevs = 1,
  cortype = "pearson",
  withparent = "parent",
  grouplev = NULL,
  box_level = NULL,
  showvals = TRUE,
  flagcolours = FALSE,
  flagthresh = c(-0.4, 0.3, 0.9),
  pval = 0.05,
  insig_colour = "#F0F0F0",
  text_colour = NULL,
  discrete_colours = NULL,
)"
```r
box_colour = NULL,
out2 = "fig"
)

Arguments

COIN The COIN object
dset The target data set.
icodes An optional list of character vectors where the first entry specifies the indicator/aggregate codes to correlate against the second entry (also a specification of indicator/aggregate codes)
aglevs The aggregation levels to take the two groups of indicators from. See getIn() for details.
cortype The type of correlation to calculate, either "pearson", "spearman", or "kendall" (see stats::cor()).
withparent If aglev[1] != aglev[2], and equal "parent" will only plot correlations of each row with its parent (default). If "family", plots the lowest aggregation level in aglevs against all its parent levels. If "none" plots the full correlation matrix.
grouplevel The aggregation level to group correlations by if aglev[1] == aglev[2]. By default, groups correlations into the aggregation level above. Set to 0 to disable grouping and plot the full matrix.
box_level The aggregation level to draw boxes around if aglev[1] == aglev[2].
showvals If TRUE, shows correlation values. If FALSE, no values shown.
flagcolours If TRUE, uses discrete colour map with thresholds defined by flagthresh. If FALSE uses continuous colour map.
pval The significance level for plotting correlations. Correlations with \( p < pval \) will be shown, otherwise they will be plotted as the colour specified by insig_colour. Set to 0 to disable this.
insig_colour The colour to plot insignificant correlations. Defaults to a light grey.
text_colour The colour of the correlation value text (default white).
discrete_colours An optional 4-length character vector of colour codes or names to define the discrete colour map if flagcolours = TRUE (from high to low correlation categories). Defaults to a green/blue/grey/purple.
box_colour The line colour of grouping boxes, default black.
out2 If "fig" returns a plot, if "dflong" returns the correlation matrix in long form, if "dfwide", returns the correlation matrix in wide form. The last option here is probably useful if you want to present a table of the data in a report.
```
Details

This function calls `getCorr()`. Note that this function can only call correlations within the same data set (i.e. only one data set in `.Data`).

Value

If `out2` = "fig" returns a plot generated with `ggplot2`. These can be edited further with `ggplot2` commands. If `out2` = "dflong" returns the correlation matrix as a data frame in long form, if `out2` = "dfwide", returns the correlation matrix in wide form. The last option here is probably useful if you want to present a table of the data in a report.

See Also

- `getCorr()` Getting correlation matrices of indicator subsets

Examples

```r
# build ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# correlation data frame of indicators in connectivity sub-index, grouped by pillar
corrs <- plotCorr(ASEM, dset = "Raw", icodes = "Conn", aglevs = 1,
showvals = F, out2 = "dflong")
# NOTE to create a plot instead set out2 = "fig"
```

Description

Plots the structure of the index using a sunburst plot using `plotly`. Output can be used as an interactive plot in html documents, e.g. via R Markdown.

Usage

`plotframework(COIN)`

Arguments

- `COIN` COIN object, or list with first entry is the indicator metadata, second entry is the aggregation metadata

Details

Note that this plot is sensitive to the order of the elements. If you use `assemble()` and input a COIN, this plot should work automatically. If you input a list, you should make sure that the indicator metadata is ordered by descending order of the hierarchy (i.e. highest level, working downwards).
plotIndDist

Value

Interactive sunburst plot built using plotly. This can be edited further with plotly commands.

Examples

# build ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# plot framework
plotframework(ASEM)

plotIndDist

Static indicator distribution plots

Description

Plots indicator distributions using box plots, dot plots, violin plots, violin-dot plots, and histograms. Supports plotting multiple indicators by calling aggregation groups.

Usage

plotIndDist(
  COINobj,
  dset = "Raw",
  icodes = NULL,
  aglev = 1,
  type = "Box",
  ntype = NULL,
  npara = NULL
)

Arguments

<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>COINobj</td>
<td>The COIN object, or a data frame of indicator data</td>
</tr>
<tr>
<td>dset</td>
<td>The source data set to use for indicator data (if input is COIN object)</td>
</tr>
<tr>
<td>icodes</td>
<td>A character vector of indicator names to plot. Defaults to all indicators.</td>
</tr>
<tr>
<td>aglev</td>
<td>The aggregation level to extract the indicator data from. Defaults to indicator level (1).</td>
</tr>
<tr>
<td>type</td>
<td>The type of plot. Currently supported &quot;Box&quot;, &quot;Dot&quot;, &quot;Violin&quot;, &quot;Violindot&quot;, &quot;Histogram&quot;.</td>
</tr>
<tr>
<td>ntype</td>
<td>The type of normalisation to apply. If NULL, no normalisation applied, otherwise specify using ntype options in normalise().</td>
</tr>
<tr>
<td>npara</td>
<td>Optional parameters to pass to normalise() if normalisation required.</td>
</tr>
</tbody>
</table>
Details

This function also optionally normalises indicators so they can be compared more easily side by side. For this purpose it calls `normalise()` - see `ntype` and `npara` arguments.

See COINr online documentation and `getIn()` for more information on accessing/plotting groups.

Value

Plots generated with `ggplot2`. These can be edited further with `ggplot2` commands.

Examples

```r
# build ASEM COIN
ASEM <- assemble(IndData = ASEMIndData, IndMeta = ASEMIndMeta, AggMeta = ASEMAggMeta)
# plot indicators in Physical pillar
plotIndDist(ASEM, type = "Box", dset = "Raw", icodes = "Physical")
```

Description

Plots a single indicator as a line of dots, and optionally highlights a selected unit.

Usage

```r
plotIndDot(
  COIN,
  dset = NULL,
  icode = NULL,
  usel = NULL,
  use_group = NULL,
  marker_type = "circle",
  add_stat = NULL,
  stat_label = NULL,
  show_ticks = TRUE,
  plabel = NULL,
  usel_label = TRUE,
  vert_adjust = 0.5
)
```

Arguments

- `COIN` The COIN
- `dset` The source data set to use for indicator data
- `icode` An indicator code to plot.
use_l | A unit or set of units (as a string or character vector) to highlight.

use_group | The name of a grouping column which must be present in the specified data set. If this is specified, the plot will be restricted to only include units within the group(s) to which use_l belongs. This argument can also be specified as a list - see function documentation for getIn().

marker_type | The type of marker, either "circle" (default) or "cross", or a marker number to pass to ggplot2 (0-25).

add_stat | A statistic to overlay, either "mean", "median" or else a specified value.

stat_label | An optional string to use as label at the point specified by add_stat.

show_ticks | Set FALSE to remove axis ticks.

plabel | Controls the labelling of the indicator. If not specified, returns the indicator name, plus units if found. Otherwise if "indname", returns only indicator name, if "indname+unit", returns indicator name plus unit (if found), if "unit" returns only unit (if found), otherwise if "none", displays no text. Finally, any other string can be passed, so e.g. "My indicator" will display this on the axis.

use_l_label | If TRUE (default) also labels selected units with their unit codes. FALSE to disable.

vert_adjust | Adjusts the vertical height of text labels and stat lines, which matters depending on plot size. Takes a value between 0 to 2 (higher will probably remove the label from the axis space).

Value

Plots generated with ggplot2. These can be edited further with ggplot2 commands.

Examples

# assemble ASEM COIN
ASEM <- assemble(ASEMIndData, ASEMIndMeta, ASEMAggMeta)
# plot CO2 indicator with highlighted countries plus median
plotIndDot(ASEM, dset = "Raw", icode = "CO2",
           use_l = c("GBR", "ESP", "AUS"), add_stat = "median")

plotSA

Plot sensitivity indices

Description

Plots sensitivity indices as bar or pie charts.

Usage

plotSA(SAresults, ptype = "bar")
Arguments

SAresults A list of sensitivity/uncertainty analysis results from sensitivity().

ptype Type of plot to generate - either "bar", "pie" or "box".

Details

To use this function you first need to run sensitivity(). Then enter the resulting list as the SAresults argument here. See COINr online documentation for more details.

Value

A plot of sensitivity indices generated by ggplot2.

See Also

• sensitivity() Perform global sensitivity or uncertainty analysis on a COIN
• plotSARanks() Plot confidence intervals on ranks following a sensitivity analysis

Examples

# build ASEM COIN up to aggregation
ASEM <- build_ASEM()

# define noise to be applied to weights
nspecs <- data.frame(AgLevel = c(2,3), NoiseFactor = c(0.25,0.25))

# create list specifying assumptions to vary and alternatives
SAspecs <- list(
  normalise = list(ntype = c("minmax", "rank", "dist2max")),
  weights = list(NoiseSpecs = nspecs, Nominal = "Original")
)

# run uncertainty analysis
# here we set N deliberately much lower than normal to enable quick testing
# Would recommend in a practical case to increase to perhaps 500 (more is always better)
SAresults <- sensitivity(ASEM, v_targ = "Index",
  SA_specs = SAspecs,
  N = 5,
  SA_type = "SA")

# Plot results as a bar chart
plotSA(SAresults, ptype = "bar")
**plotSARanks**  
*Plot ranks from an uncertainty/sensitivity analysis*

**Description**

Plots the ranks resulting from an uncertainty and sensitivity analysis, in particular plots the median, and 5th/95th percentiles of ranks.

**Usage**

```r
plotSARanks(
  SAresults,
  plot_units = NULL,
  order_by = "nominal",
  dot_colour = NULL,
  line_colour = NULL
)
```

**Arguments**

- `SAresults` A list of sensitivity/uncertainty analysis results from `sensitivity()`.
- `plot_units` A character vector of units to plot. Defaults to all units. You can also set to "top10" to only plot top 10 units, and "bottom10" for bottom ten.
- `order_by` If set to "nominal", orders the rank plot by nominal ranks (i.e. the original ranks prior to the sensitivity analysis). Otherwise if "median", orders by median ranks.
- `dot_colour` Colour of dots representing median ranks.
- `line_colour` Colour of lines connecting 5th and 95th percentiles.

**Details**

To use this function you first need to run `sensitivity()`. Then enter the resulting list as the `SAresults` argument here. See [COINr online documentation](https://coinr.readthedocs.io) for more details.

**Value**

A plot of rank confidence intervals, generated by 'ggplot2'.

**See Also**

- `sensitivity()` Perform global sensitivity or uncertainty analysis on a COIN
- `plotSA()` Plot sensitivity indices following a sensitivity analysis.
Examples

# build ASEM COIN up to aggregation
ASEM <- build_ASEM()

# define noise to be applied to weights
nspecs <- data.frame(AgLevel = c(2,3), NoiseFactor = c(0.25,0.25))

# create list specifying assumptions to vary and alternatives
SAspecs <- list(
    impute = list(imtype = c("indgroup_mean", "ind_mean", "none")),
    normalise = list(ntype = c("minmax", "rank", "dist2max")),
    weights = list(NoiseSpecs = nspecs, Nominal = "Original")
)

# run uncertainty analysis
# here we set N deliberately much lower than normal to enable quick testing
# Would recommend in a practical case to increase to perhaps 500 (more is always better)
SAresults <- sensitivity(ASEM, v_targ = "Index",
    SA_specs = SAspecs,
    N = 20,
    SA_type = "UA")

# plot rank intervals
plotSARanks(SAresults)

---

**print.COIN**

**Print COIN**

**Description**

Some details about the COIN

**Usage**

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

**Arguments**

- `x` A COIN
- `...` Arguments to be passed to or from other methods.

**Value**

Text output
Examples

```
ASEM <- build_ASEM()
print(ASEM)
```

### rankDF

**Convert a data frame to ranks**

Replaces all numerical columns of a data frame with their ranks. Uses sport ranking, i.e. ties share the highest rank place. Ignores non-numerical columns. See `rank()`. Optionally, returns in-group ranks using a specified grouping column.

#### Usage

```
rankDF(df, use_group = NULL)
```

#### Arguments

- **df**: A data frame
- **use_group**: An optional column of df (specified as a string) to use as a grouping variable. If specified, returns ranks inside each group present in this column.

#### Value

A data frame equal to the data frame that was input, but with any numerical columns replaced with ranks.

#### See Also

- `roundDF()`: Round a data frame to a specified number of decimals.

#### Examples

```
# some random data, with a column of characters
df <- data.frame(RName = c("A", "B", "C"),
                   Score1 = runif(3), Score2 = runif(3))
# convert to ranks
rankDF(df)
# grouped ranking - use some example data
df1 <- ASEMIndData[c("UnitCode", "Group_GDP", "Goods", "LPI")]
rankDF(df1, use_group = "Group_GDP")
```
Regenerate COIN object

Description

Function to regenerate the results of the COIN, using the methodological parameters stored in .Method. This function calls the construction functions of COINr in the order that they are found in .Method, along with any custom code found in .Method$Custom.

Usage

regen(COINold, quietly = FALSE)

Arguments

COINold       COIN object containing specifications on how to regenerate.
quietly       Logical: if TRUE suppresses all messages from COINr functions (warnings may still occur though).

Details

Note that while sets of weights will be passed to the regenerated COIN, anything in .Analysis will be removed and will have to be recalculated.

For more details on regeneration of COINs, comparisons and adjustments, see the online documentation.

Value

An updated COIN object, with all data sets recalculated according to specifications in .Method. Weight sets will be passed through.

See Also

- compTable() compare two different COINs
- compTableMulti() compare multiple COINs

Examples

ASEM <- build_ASEM()
# Make a copy
ASEMAltNorm <- ASEM
# Edit .Method
ASEMAltNorm$Method$normalise$ntype <- "borda"
# Regenerate
ASEMAltNorm <- regen(ASEMAltNorm, quietly = TRUE)
removeElements

Check the effect of removing indicators or aggregates

Description

This is an analysis function for seeing what happens when elements of the composite indicator are removed. This can help with "what if" experiments and acts as different measure of the influence of each indicator or aggregate.

Usage

removeElements(COIN, aglev, isel, quietly = FALSE)

Arguments

- **COIN**
  - The COIN, which must be constructed up to and including the aggregation step.
- **aglev**
  - The level at which to remove elements. For example, aglev = 1 would check the effect of removing each indicator, one at a time. aglev = 2 would check the effect of removing each of the aggregation groups above the indicator level, one at a time.
- **isel**
  - A character string indicating the indicator or aggregate code to extract from each iteration. I.e. normally this would be set to the index code to compare the ranks of the index upon removing each indicator or aggregate. But it can be any code that is present in .$Data$Aggregated.
- **quietly**
  - Logical: if FALSE (default) will output to the console an indication of progress. Might be useful when iterating over many indicators. Otherwise set to TRUE to shut this up.

Details

One way of looking at indicator "importance" in a composite indicator is via correlations. A different way is to see what happens if we remove the indicator completely from the framework. If removing an indicator or a whole aggregation of indicators results in very little rank change, it is one indication that perhaps it is not necessary to include it. Emphasis on one: there may be many other things to take into account.

This function works by successively setting the weight of each indicator or aggregate to zero. If the analysis is performed at the indicator level, it creates a copy of the COIN, sets the weight of the first indicator to zero, regenerates the results, and compares to the nominal results (results when no weights are set to zero). It repeats this for each indicator in turn, such that each time one indicator is set to zero weights, and the others retain their original weights. The output is a series of tables comparing scores and ranks (see Value).

Note that "removing the indicator" here means more precisely "setting its weight to zero". In most cases the first implies the second, but check that the aggregation method that you are using satisfies this relationship. For example, if the aggregation method does not use any weights, then setting the weight to zero will have no effect.
Value

A list with elements as follows:

- .$Scores: a data frame where each column is the scores for each unit, with indicator/aggregate corresponding to the column name removed. E.g. .$Scores$Ind1 gives the scores resulting from removing "Ind1".
- .$Ranks: as above but ranks
- .$RankDiffs: as above but difference between nominal rank and rank on removing each indicator/aggregate
- .$RankAbsDiffs: as above but absolute rank differences
- .$MeanAbsDiffs: as above, but the mean of each column. So it is the mean (over units) absolute rank change resulting from removing each indicator or aggregate.

See Also

- compTable() Comparison table between two COINs
- compTableMulti() Comparison table between multiple COINs

Examples

```r
# check the effect of removing ASEM sub-pillars, one at a time
# First build ASEM index
ASEM <- build_ASEM()
# now run check at sub-index level (level 3), on index scores/ranks
CheckPillars <- removeElements(ASEM, 3, "Index")
# summary by pillar
CheckPillars$MeanAbsDiff
# have a look at the rest of the output list to see more details.
```

---

replaceDF

**Replace multiple values in a data frame**

**Description**

Given a data frame (or vector), this function replaces values according to a look up table or dictionary. In COINr this may be useful for exchanging categorical data with numeric scores, prior to assembly. Or for changing codes.

**Usage**

`replaceDF(df, lookup)`

**Arguments**

- `df`: A data frame or a vector
- `lookup`: A data frame with columns `old` (the values to be replaced) and `new` the values to replace with. See details.
Details

The lookup data frame must not have any duplicated values in the old column. This function looks for exact matches of elements of the old column and replaces them with the corresponding value in the new column. For each row of lookup, the class of the old value must match the class of the new value. This is to keep classes of data frames columns consistent. If you wish to replace with a different class, you should convert classes in your data frame before using this function.

Value

A data frame with replaced values

See Also

• assemble() Assemble a COIN - this function optionally calls extractYear().
• rankDF() Replace numeric columns of a data frame with ranks.
• roundDF() Replace numeric columns of a data frame with rounded values.
• compareDF() Detailed comparison of two similar data frames.

Examples

# replace sub-pillar codes in ASEM indicator metadata
codeswap <- data.frame(old = c("Conn", "Sust"), new = c("SI1", "SI2"))
# swap codes in both indmeta and aggmeta
replaceDF(ASEMIndMeta, codeswap)
replaceDF(ASEMAggMeta, codeswap)
Details

This function provides a fast way to present and explore results in a COIN. It plots interactive bar charts of any indicator or aggregate, and maps if the unit codes are ISO alpha-3 country codes. It also includes an interactive results table, and the possibility to quickly compare units on a radar chart.

Value

Interactive app for exploring results (if running an interactive R session). Otherwise will simply generate a message.

See Also

- `indDash()` shiny dashboard for exploring indicator distributions
- `rew8r()` shiny dashboard for altering weights and visualising correlations

Examples

```r
# To be run in an interactive R session...
if(interactive()){
  # build ASEM up to results
  ASEM <- build_ASEM()
  # launch results dashboard
  resultsDash(ASEM)
}
```

rew8r

Re-weight indicators

Description

Interactive gadget which lets you adjust weights and see the effects. Weights can be saved with new names to the COIN object.

Usage

```r
rew8r(COIN)
```

Arguments

- `COIN` COIN object
Details

Correlations between indicators, and between indicators and index, can inform about how well the composite indicator is conveying information in the underlying indicators. For example, a correlation of zero between an indicator and the index shows that knowing the index ranks, you have no indication of what the ranks of the underlying indicator are. Since one of the objectives of a composite indicator is to summarise the information in its underlying indicators, this can be a problem.

The correlation between indicators and index (and other levels) can be adjusted by changing weights. But the effect of changing weights can be hard to gauge. The rew8r() app allows you to play around with the weights at any level, and to see what happens to the resulting correlations of interest. It also demonstrates what happens to the results. Rather than changing weights and manually regenerating the results, the rew8r() app does all this for you. If you find a set or sets of weights that you like, you can also save it/them back to the COIN as a new weight-set(s). To do this, click "Save", and then at the end of the session, "Close app".

Consider that changing weights to "fix" correlations may result in unusual sets of weights that are hard to justify. This tool may be more useful as a curiosity, as part of a "what-if" analysis, or simply to better understand what is going on inside your index.

NOTE that you need to have aggregated your data first before using this. This app also requires an interactive R session to run.

Value

An updated COIN object with additional sets of weights in .$Parameters$Weights, if specified.

See Also

- weightOpt() Correlation-optimised weights
- plotCorr() Correlation plots

Examples

```r
## Only run this example in interactive R sessions
if (interactive()) {
  # build ASEM COIN up to aggregation
  ASEM <- build_ASEM()
  # launch app
  ASEM <- rew8r(ASEM)
}
```

roundDF

Round down a data frame

Description

Tiny function just to round down a data frame for display in a table.
Usage

roundDF(df, decimals = 2)

Arguments

df A data frame to input
decimals The number of decimal places to round to (default 2)

Value

A data frame, with any numeric columns rounded to the specified amount.

See Also

- rankDF() Replace data frame numbers with ranks.

Examples

roundDF( as.data.frame(matrix(runif(20),10,2)), decimals = 3)

---

SA_estimate Estimate sensitivity indices

Description

Post process a sample to obtain sensitivity indices. This function takes a univariate output which is generated as a result of running a Monte Carlo sample from SA_sample() through a system. Then it estimates sensitivity indices using this sample.

Usage

SA_estimate(yy, N, d, Nboot = NULL)

Arguments

yy A vector of model output values, as a result of a \(N(d + 2)\) Monte Carlo design.
N The number of sample points per dimension.
d The dimensionality of the sample
Nboot Number of bootstrap draws for estimates of confidence intervals on sensitivity indices. If this is not specified, bootstrapping is not applied.

Details

This function is built to be used inside sensitivity(). See COInr online documentation for more details.
Generates an input sample for a Monte Carlo estimation of global sensitivity indices. Used in the `sensitivity()` function. The total sample size will be $N(d + 2)$.

**Usage**

```r
SA_sample(N, d)
```

**Arguments**

- `N`: The number of sample points per dimension.
- `d`: The dimensionality of the sample
Details

This function generates a Monte Carlo sample as described e.g. in the Global Sensitivity Analysis: The Primer book. See also COINr online documentation.

Value

A matrix with $N(d + 2)$ rows and $d$ columns.

See Also

- `sensitivity()` Perform global sensitivity or uncertainty analysis on a COIN.
- `SA_estimate()` Estimate sensitivity indices from system output, as a result of input design from SA_sample().

Examples

```r
# sensitivity analysis sample for 3 dimensions with 100 points per dimension
X <- SA_sample(100, 3)
```

Description

Performs global uncertainty and sensitivity analysis on a COIN.

Usage

```r
sensitivity(
  COIN,
  v_targ,
  SA_specs,
  N,
  SA_type = "UA",
  NrepWeights = 1000,
  store_results = "results+params",
  Nboot = NULL,
  quietly = FALSE
)
```

Arguments

- **COIN** A COIN (this function does not support data frame input)
- **v_targ** The target variable to perform SA or UA on. Currently just supports one variable, which should be an indicator/aggregate code present in .$Data$Aggregated.
SA_specs A list which specifies which variables to perturb, and which alternatives/distributions to use.

N The number of Monte Carlo replications.

SA_type The type of analysis to run. "UA" runs an uncertainty analysis. "SA" runs a sensitivity analysis (which anyway includes an uncertainty analysis).

NrepWeights The number of weight-replications to generate. Default 1000.

store_results Which results to store:
- "onlyresults" only stores scores, ranks and rank statistics (e.g. mean, median, quantiles)
- "results+params" (default) stores all of the above, plus a record of the parameter values used for each replication
- "results+method" stores all of the above, plus the full .$Method list of each replication
- "results+COIN" stores all results and the complete COIN of each replication (this could result in a very large list).

Nboot Number of bootstrap draws for estimates of confidence intervals on sensitivity indices. If this is not specified, bootstrapping is not applied. Ignored if SA_type = "UA".

quietly If FALSE (default), gives progress messages. Set TRUE to suppress these.

Details

To perform a sensitivity or uncertainty analysis, you must specify which parameters/assumptions to vary and what their alternative values are. This is the SA_specs argument below. To understand how this works, please see the COINr online documentation.

The output of this function can be visualised with the functions plotSARanks() and plotSA().

Value

Sensitivity analysis results as a list, containing:
- .$Scores a data frame with a row for each unit, and columns are the scores for each replication.
- .$Parameters a record of the parameters used for each iteration
- .$Ranks as .$Scores but for unit ranks
- .$RankStats summary statistics for ranks of each unit
- .$Nominal the nominal scores and ranks of each unit (i.e. from the original COIN)
- Some information on the time elapsed, average time, and the parameters perturbed.
- Depending on the setting of store_results, may also contain a list of Methods or a list of COINs for each replication.

See Also

- plotSARanks() Plot confidence intervals of ranks following UA or SA
- plotSA() Plot sensitivity indices following a sensitivity analysis
Examples

```r
# build ASEM COIN up to aggregation
ASEM <- build_ASEM()

# define noise to be applied to weights
nspecs <- data.frame(AgLevel = c(2,3), NoiseFactor = c(0.25,0.25))

# create list specifying assumptions to vary and alternatives
SAspecs <- list(
  impute = list(imtype = c("indgroup_mean", "ind_mean", "none")),
  normalise = list(ntype = c("minmax", "rank", "dist2max")),
  weights = list(NoiseSpecs = nspecs, Nominal = "Original")
)

# run uncertainty analysis
# here we set N deliberately much lower than normal to enable quick testing
# Would recommend in a practical case to increase to perhaps 500 (more is always better)
SAresults <- sensitivity(ASEM, v_targ = "Index",
                        SA_specs = SAspecs,
                        N = 15,
                        SA_type = "UA")

# to run a sensitivity analysis set SA_type = "SA" (takes longer)
```

treat  Treatment of outliers

Description

Takes the COIN object and Winsorises indicators where necessary or specified, or reverts to log transform or similar. This is done one indicator at a time.

Usage

```r
treat(
  COIN,
  dset = NULL,
  winmax = NULL,
  winchange = NULL,
  deflog = NULL,
  boxlam = NULL,
  t_skew = NULL,
  t_kurt = NULL,
  individual = NULL,
  indiv_only = NULL,
  bypass_all = NULL
)
```
treat

Arguments

**COIN**
The COIN object

**dset**
The data set to treat

**winmax**
The maximum number of points to Winsorise for each indicator. If NA, will keep Winsorising until skewness and kurtosis thresholds achieved (but it is likely this will cause errors).

**winchange**
Logical: if TRUE (default), Winsorisation can change direction from one iteration to the next. Otherwise if FALSE, no change.

**deflog**
The type of transformation to apply if Winsorisation fails. If "log", use simple \( \log(x) \) as log transform (note: indicators containing negative values will be skipped). If "CTlog", will do \( \log(x - \min(x) + a) \), where \( a = 0.01\times(\max(x) - \min(x)) \), similar to that used in the COIN Tool. If "CTlog_orig", this is exactly the COIN Tool log transformation, which is \( \log(x - \min(x) + 1) \). If "GIIlog", use GII log transformation. If "boxcox", performs a Box-Cox transformation. In this latter case, you should also specify boxlam. Finally, if "none", will return the indicator untreated.

**boxlam**
The lambda parameter of the Box-Cox transform.

**t_skew**
Absolute skew threshold (default 2)

**t_kurt**
Kurtosis threshold (default 3.5)

**individual**
A data frame specifying individual treatment for each indicator, with each row corresponding to one indicator to be treated. Columns are:

- **IndCode** The code of the indicator to be treated.
- **Treat** The type of treatment to apply, one of "win" (Winsorise), "log" (log), "GIIlog" (GII log), "CTlog" (COIN Tool log), "boxcox" (Box Cox), or "None" (no treatment).
- **Winmax** The maximum number of points to Winsorise. Ignored if the corresponding entry in "Treat" is not "win".
- **Thresh** Either NA, which means that Winsorisation will continue up to \( \text{winmax} \) with no checks on skew and kurtosis, or "thresh", which uses the skew and kurtosis thresholds specified in \( t_{skew} \) and \( t_{kurt} \).
- **boxlam** Lambda parameter for the Box Cox transformation

**indiv_only**
Logical: if TRUE, only the indicators specified in "individual" are treated. If FALSE, all indicators are treated: any outside of individual will get default treatment.

**bypass_all**
Logical: if TRUE, bypasses all data treatment and returns the original data. This is useful for sensitivity analysis and comparing the effects of turning data treatment on and off.

Details

Outliers are identified according to skewness and kurtosis thresholds. The algorithm attempts to reduce the absolute skew and kurtosis by successively Winsorising points up to a specified limit. If this limit is reached, it applies a nonlinear transformation.

The process is detailed in the COINr online documentation.
Weight optimisation

This function provides optimised weights to agree with a pre-specified vector of "target importances".

Usage

weightOpt(
  COIN,
  itarg,
  aglev,
  cortype = "pearson",
  optype = "balance",
  toler = NULL,
  maxiter = NULL,
  out2 = NULL
)
weightOpt

Arguments

- **COIN**: COIN object
- **itarg**: a vector of (relative) target importances. For example, c(1, 2, 1) would specify that the second indicator should be twice as "important" as the other two.
- **aglev**: The aggregation level to apply the weight adjustment to.
- **cortype**: The type of correlation to use - can be either "pearson", "spearman" or "kendall". See `stats::cor`.
- **optype**: The optimisation type. Either "balance", which aims to balance correlations according to a vector of "importances" specified by `itarg` (default), or "infomax" which aims to maximise overall correlations. *This latter option is experimental and may not yet work very well.*
- **toler**: Tolerance for convergence. Defaults to 0.001 (decrease for more accuracy, increase if convergence problems).
- **maxiter**: Maximum number of iterations. Default 500.
- **out2**: Where to output the results. If "COIN" (default for COIN input), appends to updated COIN, creating a new list of weights in `.Parameters$Weights`. Otherwise if "list" outputs to a list (default).

Details

This is a linear version of the weight optimisation proposed in this paper: doi: 10.1016/j.ecolind.2017.03.056. Weights are optimised to agree with a pre-specified vector of "importances". The optimised weights are returned back to the COIN.

See the chapter in the COINr online documentation for more details.

Value

If out2 = "COIN" returns an updated COIN object with a new set of weights in `.Parameters$Weights`, plus details of the optimisation in `.Analysis`. Else if out2 = "list" the same outputs (new weights plus details of optimisation) are wrapped in a list.

See Also

- `rew8r()` Interactive app for adjusting weights and seeing effects on correlations
- `getPCA()` PCA, including weights from PCA

Examples

```r
# build ASEM COIN up to aggregation
ASEM <- build_ASEM()

# optimise sub-pillar weights to give equal correlations with index
ASEM <- weightOpt(ASEM, itarg = "equal", aglev = 3, out2 = "COIN")
```
weights2corr (Recalculate correlations and ranks based on new weights)

Description
This is a short cut function which takes a new set of indicator weights, and recalculates the COIN results based on these weights. It returns a summary of rankings and the correlations between indicators and index.

Usage
weights2corr(
  COIN, 
  w, 
  aglevs = NULL, 
  icodes = NULL, 
  cortype = "pearson", 
  withparent = TRUE 
)

Arguments
<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>COIN</td>
<td>COIN object</td>
</tr>
<tr>
<td>w</td>
<td>Full data frame of weights for each level</td>
</tr>
<tr>
<td>aglevs</td>
<td>A 2-length vector with two aggregation levels to correlate against each other</td>
</tr>
<tr>
<td>icodes</td>
<td>List of two character vectors of indicator codes, corresponding to the two aggregation levels</td>
</tr>
<tr>
<td>cortype</td>
<td>Correlation type. Either &quot;pearson&quot; (default), &quot;kendall&quot; or &quot;spearman&quot;. See stats::cor.</td>
</tr>
<tr>
<td>withparent</td>
<td>Logical: if TRUE, only correlates with the parent, e.g. sub-pillars are only correlated with their parent pillars and not others.</td>
</tr>
</tbody>
</table>

Details
This function is principally used inside rew8r(). The w argument should be a data frame of weights, of the same format as the data frames found in .$Parameters$Weights.

Value
A list where .$cr is a vector of correlations between each indicator and the index, and .$dat is a data frame of rankings, with unit code, and index, input and output scores

See Also
- rew8r() Interactive app for adjusting weights and seeing effects on correlations
- getCorr() Get correlations between indicators/levels
Examples

# build ASEM COIN up to aggregation
ASEM <- build_ASEM()
# get correlations between pillars (level 2) and index (level 4)
# original weights used just for demonstration, normally you would alter first.
l <- weights2corr(ASEM, ASEM$Parameters$Weights$Original, aglevs = c(2,4))

WorldDenoms

World denomination data

Description

A small selection of common denominator indicators, which includes GDP, Population, Area, GDP per capita and income group. All data sourced from the World Bank as of Feb 2021 (data is typically from 2019). Note that this is intended as example data, and it would be a good idea to use updated data from the World Bank when needed. In this data set, country names have been altered slightly so as to include no accents - this is simply to make it more portable between distributions.

Usage

WorldDenoms

Format

A data frame with 249 rows and 7 variables.

Source

https://data.worldbank.org/
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