

Package ‘softwareRisk’

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Type Package

Title Computation of Node and Path-Level Risk Scores in Scientific Models

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Description It leverages the network-like architecture of scientific models together with software quality metrics to identify chains of function calls that are more prone to generating and propagating errors. It operates on `tbl_graph` objects representing call dependencies between functions (callers and callees) and computes risk scores for individual functions and for paths (sequences of function calls) based on cyclomatic complexity, in-degree and betweenness centrality. The package supports variance-based uncertainty and sensitivity analyses after Puy et al. (2022) <[doi:10.18637/jss.v102.i05](https://doi.org/10.18637/jss.v102.i05)> to assess how risk scores change under alternative risk definitions.

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all_paths_fun	<i>Enumerate entry-to-sink call paths and compute risk metrics at the node and path level</i>
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Description

Given a directed call graph (`tidygraph::tbl_graph`) with a node attribute for cyclomatic complexity, this function:

- computes node-level metrics (in-degree, out-degree, betweenness),
- calculates a node risk score as a weighted combination of rescaled metrics,
- enumerates all simple paths from entry nodes (in-degree = 0) to sink nodes (out-degree = 0),
- computes path-level summaries and a path-level risk score.
- calculates a gini index and the slope of risk at the path-level.

Usage

```
all_paths_fun(
  graph,
  alpha = 0.6,
  beta = 0.3,
  gamma = 0.1,
  p = 1,
  eps = 1e-12,
  complexity_col = "cyclo",
  weight_tol = 1e-08
)
```

Arguments

graph	A directed tidygraph::tbl_graph. Graph nodes must have a name attribute (i.e., igraph::V(as.igraph(graph))\$name) and a numeric node attribute specified by complexity_col.
alpha, beta, gamma	Numeric non-negative weights for the risk score, constrained such that alpha + beta + gamma == 1 (within weight_tol).
p	Numeric scalar. Power parameter for the weighted power mean. Must be finite and lie in the interval $[-1, 2]$. When $p = 1$ the formula reduces to a weighted sum. Default 1.
eps	Numeric. Small positive constant ϵ used for numerical stability in the $p \rightarrow 0$ (geometric mean) case and when $p < 0$, where zero-valued normalized metrics are replaced by ϵ to avoid non-finite intermediate values. Default 1e-12.
complexity_col	Character scalar. Name of the node attribute containing cyclomatic complexity. Default "cyclo".
weight_tol	Numeric tolerance for enforcing the weight-sum constraint. Default 1e-8.

Details

The normalized node metrics are computed using scales::rescale() and denoted by a tilde $\tilde{\cdot}$.

The risk score for node v_i is computed as the weighted power mean of normalized metrics:

$$r_{(v_i)} = \left(\alpha \tilde{C}_{(v_i)}^p + \beta \tilde{d}_{(v_i)}^{np} + \gamma \tilde{b}_{(v_i)}^p \right)^{1/p},$$

where p is the power-mean parameter. When $p = 1$ this reduces to a weighted sum (additive). In the limit $p \rightarrow 0$, this reduces to a weighted geometric mean, implemented with a small constant ϵ to ensure numerical stability:

$$r_{(v_i)} = \exp \left(\alpha \log(\max(\tilde{C}_{(v_i)}, \epsilon)) + \beta \log(\max(\tilde{d}_{(v_i)}^n, \epsilon)) + \gamma \log(\max(\tilde{b}_{(v_i)}, \epsilon)) \right).$$

The path-level risk score is calculated as

$$P_k = 1 - \prod_{i=1}^{n_k} (1 - r_{k(v_i)}),$$

where $r_{k(v_i)}$ is the risk of the i -th function in path k and n_k is the number of functions in that path. The equation behaves like a saturating OR-operator: P_k is at least as large as the maximum individual function risk and monotonically increases as more functions on the path become risky, approaching 1 when several functions have high risk scores.

The Gini index of path k is computed as

$$G_k = \frac{\sum_i \sum_j |r_{k(v_i)} - r_{k(v_j)}|}{2n_k^2 \bar{r}_k},$$

where $\bar{r}_k$ is the mean node risk in path k .

Finally, the trend of risk is defined by the slope of the regression

$$r_{k(v_i)} = \theta_{0k} + \theta_{1k} i + \epsilon_i,$$

where $r_{k(v_i)}$ is the risk score of the function at position i along path k (ordered from upstream to downstream execution) and ϵ_i is a residual term.

The returned paths tibble includes path_cc, a list-column where each element is the vector of per-node cyclomatic complexity values along the path.

Value

A named list with two tibbles:

nodes Node-level metrics with columns name, cyclomatic_complexity, indeg (in-degree), outdeg (out-degree), btw (betweenness), risk_score.

paths Path-level metrics with columns path_id, path_nodes, path_str, hops, path_risk_score, path_cc, gini_node_risk, risk_slope, risk_mean, risk_sum

Examples

```
# synthetic_graph is a tidygraph::tbl_graph with node attribute "cyclo"
data(synthetic_graph)
```

```
# additive risk (p = 1, default)
out1 <- all_paths_fun(
  graph = synthetic_graph,
  alpha = 0.6, beta = 0.3, gamma = 0.1,
  p = 1,
  complexity_col = "cyclo"
)
```

```
# power-mean risk (p = 0 ~ weighted geometric mean)
out2 <- all_paths_fun(
  graph = synthetic_graph,
  alpha = 0.6, beta = 0.3, gamma = 0.1,
  p = 0,
  eps = 1e-12,
  complexity_col = "cyclo"
)
```

```
out1$nodes
out1$paths
```

call_graph_fun	<i>Build a call graph from an R package or a directory of R scripts</i>
----------------	---

Description

Constructs the directed call graph required by [all_paths_fun\(\)](#) automatically, so that no manual preparation of edge lists or complexity spreadsheets is needed for R code. Each node is a function, each edge from `->` to is a call from function from (caller) to function to (callee), and each node carries a `cyclo` attribute with its cyclomatic complexity.

Usage

```
call_graph_fun(pkg = NULL, dir = NULL, exclude = NULL, keep_self_loops = FALSE)
```

Arguments

<code>pkg</code>	Character scalar. Name of an installed package whose namespace functions (exported and internal) are analyzed. Exactly one of <code>pkg</code> or <code>dir</code> must be supplied.
<code>dir</code>	Character scalar. Path to a directory containing <code>.R</code> files. Files are parsed (searched recursively) and all top-level name <code><- function(...)</code> definitions (also with <code>=</code> or <code><<-</code>) are collected. The code is never executed: function literals are only parsed and turned into closures.
<code>exclude</code>	Optional character vector of function names to drop from the graph (e.g., generated or vendored code).
<code>keep_self_loops</code>	Logical. Keep self-recursion edges (a function calling itself)? These edges do not affect the simple-path enumeration of all_paths_fun() but change the in-degree. Default <code>FALSE</code> .

Details

Call detection relies on `codetools::findGlobals()`, i.e., on static analysis of each function body. Two limitations follow: (i) calls built at run time (e.g., `do.call(paste0("f", i), ...)`, `get()`, method dispatch) cannot be detected; (ii) a function of the set that is merely *referenced* (e.g., passed to `lapply()`) is treated as called, since in a dependency sense the caller relies on it.

Cyclomatic complexity is computed internally as 1 plus one unit per decision point (`if`, `for`, `while`, `repeat`, `&&`, `||`, and each `switch()` alternative beyond the first), following McCabe (1976).

Value

A directed `tidygraph::tbl_graph` with node attributes `name` and `cyclo`, ready to be passed to [all_paths_fun\(\)](#).

References

McCabe, T. J. (1976). *A Complexity Measure*. IEEE Transactions on Software Engineering, SE-2(4), 308–320. doi:10.1109/TSE.1976.233837

See Also

`read_call_graph()` to import a call graph prepared outside R (e.g., for Fortran, C or Python models).

Examples

```
# build the call graph of a directory of R scripts
td <- file.path(tempdir(), "cg_example")
dir.create(td, showWarnings = FALSE)
writeLines(c(
  "load_data <- function(x) x",
  "clean_data <- function(x) load_data(x)",
  "calc_scores <- function(x) if (length(x) > 0) mean(x) else 0",
  "compute_risk <- function(x) calc_scores(clean_data(x))"
), file.path(td, "model.R"))

g <- call_graph_fun(dir = td)
g

# the result feeds directly into the pipeline
out <- all_paths_fun(g, alpha = 0.6, beta = 0.3, gamma = 0.1)
out$paths

# build the call graph of an installed package

g_pkg <- call_graph_fun(pkg = "softwareRisk")
g_pkg
```

fix_portfolio_fun

Greedy selection of the most risk-reducing node fixes

Description

Selects, under a budget of budget refactoring interventions, the set of nodes whose fixing (risk set to 0) most reduces the path-level risk of the system. Whereas `path_fix_heatmap()` shows the effect of fixing single nodes on single paths, this function answers the budgeted question directly: *given resources to refactor k functions, which k ?*

Usage

```
fix_portfolio_fun(
  all_paths_out,
  budget = 5,
  objective = c("total", "max"),
  tol = 1e-12
)
```

Arguments

all_paths_out	A list produced by <code>all_paths_fun()</code> with elements nodes and paths. nodes must contain name and risk_score; paths must contain path_nodes and path_risk_score.
budget	Integer. Maximum number of nodes to fix. Default 5.
objective	Character scalar, "total" (default) or "max". See Details.
tol	Numeric. Minimum improvement in the objective required to continue selecting nodes. Default 1e-12.

Details

At each step the algorithm evaluates, for every remaining candidate node, the objective obtained by adding that node to the already-selected set, fixes the best node, and repeats. Path risk is recomputed under the independence assumption used throughout the package:

$$P_k = 1 - \prod_{i=1}^{n_k} (1 - r_{k(v_i)}),$$

with $r = 0$ for fixed nodes. Two objectives are available:

- "total": minimize $\sum_k P_k$, the total path risk. This objective is monotone submodular in the set of fixed nodes, so the greedy solution is guaranteed to achieve at least $1 - 1/e \approx 63\%$ of the reduction of the optimal set of the same size (Nemhauser et al. 1978).
- "max": minimize $\max_k P_k$, the risk of the riskiest path.

The greedy search stops early if no remaining node yields an improvement larger than tol.

Value

A list with two elements:

- portfolio: a tibble with one row per selected node, in selection order, with columns step, node, objective_after (value of the objective once the node is fixed), delta (improvement contributed by the node) and cum_reduction (cumulative fraction of the initial objective removed).
- plot: a **ggplot2** object showing the objective as a function of the number of fixed nodes, with the selected nodes labelled.

References

Nemhauser, G. L., Wolsey, L. A., and Fisher, M. L. (1978). *An analysis of approximations for maximizing submodular set functions—I*. Mathematical Programming, 14(1), 265–294. doi:10.1007/BF01588971

See Also

`path_fix_heatmap()` for the node-by-path improvement matrix and `node_exposure_fun()` for path-aware node criticality.

Examples

```
data(synthetic_graph)
out <- all_paths_fun(graph = synthetic_graph, alpha = 0.6, beta = 0.3,
                    gamma = 0.1, complexity_col = "cyclo")
res <- fix_portfolio_fun(out, budget = 5)
res$portfolio
res$plot
```

gini_index_fun	<i>Compute the Gini index of a numeric vector</i>
----------------	---

Description

Computes the Gini index (a measure of inequality) for a numeric vector. Non-finite (NA, NaN, Inf) values are removed prior to computation. If fewer than two finite values remain, or if all finite values are zero (no inequality is measurable), the function returns 0.

Usage

```
gini_index_fun(x)
```

Arguments

x Numeric vector.

Details

The Gini index ranges from 0 (perfect equality) to 1 (maximal inequality).

Value

A numeric scalar giving the Gini index of x.

Examples

```
gini_index_fun(c(1, 1, 1, 1))
gini_index_fun(c(1, 2, 3, 4))
gini_index_fun(c(NA, 1, 2, Inf, 3))
```

node_exposure_fun	<i>Path-aware node criticality (risk load)</i>
-------------------	--

Description

Summarizes, for each node, its participation in the entry-to-sink paths enumerated by `all_paths_fun()`. Whereas the node risk score captures how error-prone a function is in isolation, the *risk load* captures how many risky execution chains depend on it: a function of moderate complexity can still be critical if most high-risk paths route through it.

Usage

```
node_exposure_fun(all_paths_out)
```

Arguments

`all_paths_out` A list produced by `all_paths_fun()` with elements nodes and paths. nodes must contain name and risk_score; paths must contain path_nodes and path_risk_score.

Details

For node v the function computes:

- `n_paths`: the number of paths that contain v ,
- `path_share`: `n_paths` divided by the total number of paths,
- `risk_load`: $\sum_{k:v \in k} P_k$, the sum of `path_risk_score` over the paths containing v ,
- `risk_load_share`: `risk_load` divided by $\sum_k P_k$,
- `mean_path_risk`: the mean `path_risk_score` of the paths containing v .

The output also reports the rank of each node by `risk_load` and by its standalone `risk_score`, so that structurally exposed functions can be told apart from merely complex ones: a node ranking much higher on `risk_load` than on `risk_score` is a chokepoint rather than a hotspot.

Value

A tibble with one row per node appearing on at least one path, sorted by decreasing `risk_load`, with columns `name`, `risk_score`, `n_paths`, `path_share`, `risk_load`, `risk_load_share`, `mean_path_risk`, `rank_risk_load` and `rank_risk_score`. Nodes on no path (e.g., disconnected helpers) are omitted. If `all_paths_out$paths` is empty, an empty tibble with these columns is returned.

See Also

`path_fix_heatmap()` for the per-path effect of fixing single nodes, and `fix_portfolio_fun()` for a budgeted selection of fixes.

Examples

```
data(synthetic_graph)
out <- all_paths_fun(graph = synthetic_graph, alpha = 0.6, beta = 0.3,
                    gamma = 0.1, complexity_col = "cyclo")
node_exposure_fun(out)
```

path_fix_heatmap

Path-level improvement from fixing high-risk nodes

Description

Compute how much the risk score of the riskiest paths would decrease if selected high-risk nodes were made perfectly reliable (risk fixed to 0), and visualise the result as a heatmap.

Usage

```
path_fix_heatmap(all_paths_out, n_nodes = 20, k_paths = 20)
```

Arguments

all_paths_out	A list returned by <code>all_paths_fun()</code> , with elements nodes and paths. nodes must contain at least the columns name and risk_score. paths must contain at least the columns path_id, path_nodes (list-column of node names) and path_risk_score.
n_nodes	Integer, number of top-risk nodes (by risk_score) to include as rows in the heatmap. Defaults to 20.
k_paths	Integer, number of top-risk paths (by path_risk_score) to include as columns in the heatmap. Defaults to 20.

Details

For each of the top `n_nodes` nodes ranked by `risk_score` and the top `k_paths` paths ranked by `path_risk_score`, the function sets the risk of that node to 0 along the path (for all its occurrences) and recomputes the path risk score under the independence assumption, using

$$P_k = 1 - \prod_{i=1}^{n_k} (1 - r_{k(v_i)})$$

The improvement

$$\Delta P_k = R_{\text{orig}} - R_{\text{fix}}$$

is used as the fill value in the heatmap cells.

Bright cells indicate nodes that act as chokepoints for a given path. Rows with many bright cells correspond to nodes whose refactoring would improve many risky paths (global chokepoints), while columns with a few very bright cells correspond to paths dominated by a single risky node.

Value

A list with two elements:

- `delta_tbl`: a tibble with columns `node`, `path_id` and `deltaR`, containing the reduction in path risk score when fixing the node in that path.
- `plot`: a **ggplot2** object containing the heatmap.

Examples

```
data(synthetic_graph)
out <- all_paths_fun(graph = synthetic_graph, alpha = 0.6, beta = 0.3,
  gamma = 0.1, complexity_col = "cyclo")
res <- path_fix_heatmap(all_paths_out = out, n_nodes = 20, k_paths = 20)
res
```

`path_uncertainty_plot` *Plot path-level uncertainty for the top-risk paths*

Description

Plot the top `n_paths` paths ranked by their mean risk score, with horizontal error bars representing the uncertainty range (minimum and maximum risk) computed from the Monte Carlo samples stored in `uncertainty_analysis`.

Usage

```
path_uncertainty_plot(ua_sa_out, n_paths = 20)
```

Arguments

<code>ua_sa_out</code>	A list returned by <code>uncertainty_fun()</code> containing at least an element <code>\$paths</code> , which must be a data frame with columns <code>path_id</code> and <code>uncertainty_analysis</code> . The column <code>uncertainty_analysis</code> is expected to be a list-column where each element is a numeric vector of path risk values obtained from Monte Carlo sampling.
<code>n_paths</code>	Integer, number of top paths (by mean risk) to include in the plot. Defaults to 20.

Details

This function is designed to work with the `paths` component of the output of `uncertainty_fun()`. For each path, it summarises the vector of path risk values by computing the mean, minimum and maximum values, and then displays these summaries for the `n_paths` most risky paths.

Value

A **ggplot2** object.

Examples

```
data(synthetic_graph)
out <- all_paths_fun(graph = synthetic_graph, alpha = 0.6, beta = 0.3,
  gamma = 0.1, complexity_col = "cyclo")
results <- uncertainty_fun(all_paths_out = out, N = 2^10, order = "first")
path_uncertainty_plot(ua_sa_out = results, n_paths = 20)
```

plot_top_paths_fun	<i>Plot the top risky call paths on a Sugiyama layout</i>
--------------------	---

Description

Visualizes the most risky entry-to-sink paths (by decreasing `path_risk_score`) computed by `all_paths_fun()`. Edges that occur on the top paths are highlighted, with edge colour mapped to the mean path risk and edge width mapped to the number of top paths using that edge. Nodes on the top paths are emphasized, with node size mapped to in-degree and node fill mapped to binned cyclomatic complexity.

Usage

```
plot_top_paths_fun(
  graph,
  all_paths_out,
  model.name = "",
  language = "",
  top_n = 10,
  alpha_non_top = 0.05
)
```

Arguments

<code>graph</code>	A directed <code>tidygraph::tbl_graph</code> representing the call graph to plot (typically the same graph used as input to <code>all_paths_fun()</code>).
<code>all_paths_out</code>	Output from <code>all_paths_fun()</code> , i.e. a list with elements <code>nodes</code> and <code>paths</code> .
<code>model.name</code>	Character scalar used in the plot title (e.g., model name).
<code>language</code>	Character scalar used in the plot title (e.g., language name).
<code>top_n</code>	Integer. Number of highest-risk paths to display (default 10).
<code>alpha_non_top</code>	Numeric between 0 and 1. Alpha (transparency) for edges that are not on the top-risk paths. Smaller values fade background edges more.

Details

The function selects the top_n paths by sorting paths_tbl on path_risk_score (descending). For those paths, it:

- builds an edge list from path_nodes,
- marks graph edges that appear on at least one top path,
- computes path_freq (how many top paths include each edge),
- computes risk_mean_path (mean of risk_sum across top paths that include each edge),
- highlights nodes that appear on any top path.

Node fills are based on cyclomatic_complexity using breaks $(-\text{Inf}, 10]$, $(10, 20]$, $(20, 50]$, $(50, \text{Inf}]$ as per Watson & McCabe (1996).

Value

A ggplot object (invisibly). The plot is also printed as a side effect.

References

Watson, A. H. and McCabe, T. J. (1996). *Structured Testing: A Testing Methodology Using the Cyclomatic Complexity Metric*. NIST Special Publication 500-235, National Institute of Standards and Technology, Gaithersburg, MD. doi:10.6028/NIST.SP.500-235

Examples

```
data(synthetic_graph)
out <- all_paths_fun(graph = synthetic_graph, alpha = 0.6, beta = 0.3,
  gamma = 0.1, complexity_col = "cyclo")
p <- plot_top_paths_fun(synthetic_graph, out, model.name = "MyModel", language = "R", top_n = 10)
p
```

rank_robustness_fun	<i>Robustness of the risk ranking under uncertainty</i>
---------------------	---

Description

Quantifies how stable the identification of the top- k riskiest paths (or nodes) is across the uncertainty draws produced by `uncertainty_fun()`. Every draw corresponds to one plausible definition of risk (one sampled combination of the weights α, β, γ and the exponent p); ranking the paths within each draw therefore shows which paths are flagged as risky *regardless* of how risk is weighted and which ones enter the top- k only under specific assumptions.

Usage

```
rank_robustness_fun(ua_sa_out, top_k = 10, what = c("paths", "nodes"))
```

Arguments

ua_sa_out	A list returned by <code>uncertainty_fun()</code> with elements nodes and paths, each containing an <code>uncertainty_analysis</code> list-column of numeric draws.
top_k	Integer. Size of the top set whose membership is tracked. Clamped to the number of items with a message if larger. Default 10.
what	Character scalar, "paths" (default) or "nodes".

Details

For each item (path or node) the function reports the probability of belonging to the top- k across draws and the median and the 5th and 95th percentiles of its rank (rank 1 = riskiest; ties share the minimum rank). As a global summary it also reports the mean Spearman correlation between the risk values of each draw and the consensus (mean across draws) risk values: values close to 1 indicate that the ranking barely responds to the risk definition, values well below 1 that conclusions depend on it.

Draws that are NA for an item are treated as never placing that item in the top- k .

Value

A list with:

- `summary`: a tibble with one row per item, sorted by decreasing `top_k_prob`, with columns `id` (path `path_id` or node name), `path_str` (paths only), `mean_risk`, `top_k_prob`, `rank_median`, `rank_q05` and `rank_q95`.
- `consensus_correlation`: mean Spearman correlation between each draw and the consensus risk values.
- `top_k`, `what`: the settings used.

See Also

`rank_robustness_plot()` to visualize the result and `uncertainty_fun()` to generate the draws.

Examples

```
data(synthetic_graph)
out <- all_paths_fun(graph = synthetic_graph, alpha = 0.6, beta = 0.3,
                    gamma = 0.1, complexity_col = "cyclo")
results <- uncertainty_fun(all_paths_out = out, N = 2^7, order = "first")

rr <- rank_robustness_fun(results, top_k = 10, what = "paths")
rr$summary
rr$consensus_correlation
```

rank_robustness_plot *Plot the robustness of the risk ranking*

Description

Displays, for the items most often ranked in the top- k , the probability of top- k membership across the uncertainty draws computed by `rank_robustness_fun()`. Items with probability close to 1 are flagged as risky under essentially any risk definition; intermediate probabilities identify items whose criticality depends on how the risk score is weighted.

Usage

```
rank_robustness_plot(rank_out, top_n = 20)
```

Arguments

`rank_out` A list returned by `rank_robustness_fun()`.
`top_n` Integer. Number of items (by decreasing `top_k_prob`) to display. Default 20.

Value

A **ggplot2** object.

Examples

```
data(synthetic_graph)
out <- all_paths_fun(graph = synthetic_graph, alpha = 0.6, beta = 0.3,
                    gamma = 0.1, complexity_col = "cyclo")
results <- uncertainty_fun(all_paths_out = out, N = 2^7, order = "first")
rr <- rank_robustness_fun(results, top_k = 10, what = "paths")
rank_robustness_plot(rr, top_n = 20)
```

read_call_graph *Import a call graph from edge-list and complexity tables*

Description

Builds the directed call graph required by `all_paths_fun()` from the two datasets described in the package vignette: an edge list of function calls and a table of per-function cyclomatic complexity. The input is validated so that common preparation errors (misspelled columns, functions missing from the complexity table, duplicated or non-numeric entries) fail early with an informative message. This is the recommended entry point for models written in languages other than R (e.g., Fortran, C, Python), whose edge lists and complexity values are extracted with external tools.

Usage

```
read_call_graph(
  edges,
  metrics,
  from_col = "from",
  to_col = "to",
  name_col = "name",
  complexity_col = "cyclo"
)
```

Arguments

<code>edges</code>	A data.frame (or path to a .csv file) with one row per function call. Must contain the columns given by <code>from_col</code> (caller) and <code>to_col</code> (callee).
<code>metrics</code>	A data.frame (or path to a .csv file) with one row per function. Must contain the columns given by <code>name_col</code> (function name) and <code>complexity_col</code> (cyclomatic complexity).
<code>from_col, to_col</code>	Character scalars. Names of the caller and callee columns in edges. Defaults "from" and "to".
<code>name_col</code>	Character scalar. Name of the function-name column in metrics. Default "name".
<code>complexity_col</code>	Character scalar. Name of the complexity column in metrics. Default "cyclo". The column keeps this name in the returned graph, matching the default of all_paths_fun() .

Value

A directed `tidygraph::tbl_graph` with node attributes `name` and `complexity_col`, ready to be passed to [all_paths_fun\(\)](#).

See Also

[call_graph_fun\(\)](#) to build the graph automatically from R source code.

Examples

```
calls_df <- data.frame(
  from = c("clean_data", "compute_risk", "compute_risk", "calc_scores"),
  to   = c("load_data", "clean_data", "calc_scores", "trim_mean")
)
cyclo_df <- data.frame(
  name = c("clean_data", "load_data", "compute_risk", "calc_scores",
           "trim_mean"),
  cyclo = c(6, 3, 12, 5, 2)
)

g <- read_call_graph(edges = calls_df, metrics = cyclo_df)
g
```

sensitivity_plot_fun *Plot Sobol' sensitivity indices of the node risk scores*

Description

Visualizes the Sobol' indices stored in the `sensitivity_analysis` list-column returned by `uncertainty_fun()`, showing how much of the variance in the node risk scores is driven by each parameter of the risk definition (the weights α , β , γ and the power-mean exponent p).

Usage

```
sensitivity_plot_fun(ua_sa_out, nodes = NULL, index = c("both", "Si", "Ti"))
```

Arguments

<code>ua_sa_out</code>	A list returned by <code>uncertainty_fun()</code> with element nodes containing a <code>sensitivity_analysis</code> list-column.
<code>nodes</code>	Optional character vector of node names to display individually. Default <code>NULL</code> (aggregate view across all nodes).
<code>index</code>	Character scalar: "both" (default), "Si" (first-order) or "Ti" (total-order). Which indices to display.

Details

Two displays are available:

- With `nodes = NULL` (default), the distribution of the indices across *all* nodes is shown as one boxplot per parameter, answering the system-level question of which assumption dominates the risk ranking overall.
- With `nodes` set to a character vector of node names, the indices of those nodes are shown as dodged bars with their confidence intervals, one panel per node.

Nodes whose risk score does not vary across draws (e.g., all metrics zero) yield non-finite indices; these are dropped with a message.

Value

A **ggplot2** object.

References

Puy, A., Lo Piano, S., Saltelli, A., and Levin, S. A. (2022). *sensobol: An R Package to Compute Variance-Based Sensitivity Indices*. *Journal of Statistical Software*, 102(5), 1–37. doi:10.18637/jss.v102.i05

See Also

`uncertainty_fun()` to generate the indices and the **sensobol** package (Puy et al. 2022) for their estimation.

Examples

```
data(synthetic_graph)
out <- all_paths_fun(graph = synthetic_graph, alpha = 0.6, beta = 0.3,
                    gamma = 0.1, complexity_col = "cyclo")
results <- uncertainty_fun(all_paths_out = out, N = 2^7, order = "first")

# system-level view: which assumption drives the risk scores?
sensitivity_plot_fun(results)

# node-level view
sensitivity_plot_fun(results, nodes = results$nodes$name[1:4])
```

slope_fun

*Compute the linear slope of a numeric sequence***Description**

Computes the slope of a simple linear regression of a numeric vector against its index (`seq_along(x)`). Non-finite (NA, NaN, Inf) values are removed prior to computation, but the remaining values keep their original positions as the regressor, so removing values does not distort the trend. If fewer than two finite values remain, the function returns 0.

Usage

```
slope_fun(x)
```

Arguments

`x` Numeric vector.

Details

The slope is estimated from the model $x_i = \beta_0 + \beta_1 i + \varepsilon_i$, where $i = 1, \dots, n$. The function returns the estimated slope β_1 .

This summary is useful for characterizing monotonic trends in ordered risk values along a path.

Value

A numeric scalar giving the slope of the fitted linear trend.

Examples

```
slope_fun(c(1, 2, 3, 4))  
slope_fun(c(4, 3, 2, 1))  
slope_fun(c(NA, 1, 2, Inf, 3))
```

synthetic_graph

Synthetic citation graph for software risk examples

Description

A synthetic directed graph with cyclomatic complexity values to illustrate the functions of the package.

Usage

```
data(synthetic_graph)
```

Format

A `tbl_graph` with:

nodes 55 nodes

edges 122 directed edges

Details

The graph is stored as a `tbl_graph` object with:

- Node attributes: `name`, `cyclo`
- Directed edges defined by `from` → `to`

theme_AP

A clean, publication-oriented ggplot2 theme

Description

`theme_AP()` provides a minimalist, publication-ready theme based on `ggplot2::theme_bw()`, with grid lines removed, compact legends, and harmonized text sizes. It is designed for dense network and path-visualization plots (e.g. call graphs, risk paths).

Usage

```
theme_AP()
```

Details

The theme:

- removes major and minor grid lines,
- uses transparent legend backgrounds and keys,
- standardizes text sizes for axes, legends, strips, and titles,
- reduces legend spacing for compact layouts.

This theme is intended to be composable: it should be added to a ggplot object using `+ theme_AP()`.

Value

A `ggplot2::theme` object.

Examples

```
ggplot2::ggplot(mtcars, ggplot2::aes(mpg, wt)) +
  ggplot2::geom_point() +
  theme_AP()
```

uncertainty_fun

Uncertainty and sensitivity analysis of node and path risk

Description

Runs a full variance-based uncertainty and sensitivity analysis (UA/SA) for node risk scores using the results returned by [all_paths_fun\(\)](#) and the functions provided by the **sensobol** package (Puy et al. 2022).

Usage

```
uncertainty_fun(all_paths_out, N, order, eps = 1e-12)
```

Arguments

- | | |
|----------------------------|--|
| <code>all_paths_out</code> | A list produced by all_paths_fun() with elements <code>nodes</code> and <code>paths</code> . <code>nodes</code> must contain columns <code>name</code> , <code>cyclomatic_complexity</code> , <code>indeg</code> , <code>btw</code> ; <code>paths</code> must contain <code>path_id</code> , <code>path_nodes</code> , <code>path_str</code> , and <code>hops</code> . |
| <code>N</code> | Integer. Base sample size used for Sobol' matrices. |
| <code>order</code> | Passed to <code>sensobol::sobol_matrices()</code> and <code>sensobol::sobol_indices()</code> to control which Sobol indices are computed (e.g., first/total/second order), depending on your implementation. |
| <code>eps</code> | Numeric. Small positive constant ϵ used for numerical stability in the $p \rightarrow 0$ evaluation and when $p < 0$, where zero-valued rescaled metrics are replaced by ϵ to avoid non-finite intermediate values. Default <code>1e-12</code> . |

Details

Uncertainty is induced by jointly sampling the weights (α, β, γ) (renormalized to sum to 1) and the power parameter $p \in [-1, 2]$ used in the node-risk definition:

$$r = \left(\alpha \tilde{C}^p + \beta (\tilde{d}^{\text{in}})^p + \gamma \tilde{b}^p \right)^{1/p}.$$

For each node, risk scores are repeatedly recalculated using the sampled parameter combinations, producing a distribution of possible outcomes. Sobol' first-order and/or total-order sensitivity indices are then computed for all four parameters (α , β , γ , and p), quantifying how much of the variance in the node risk score is attributable to each parameter.

Parameter labels and the Sobol' design. Internally the design samples four independent $U(0, 1)$ values (a_raw, b_raw, c_raw, p_raw) because the Sobol' quasi-random sequence requires independent uniform inputs. Before evaluating the risk model, the raw draws are transformed: the three weight draws are normalised to sum to one, yielding α , β , γ ; and p_raw is mapped linearly to $p \in [-1, 2]$. The sensitivity indices are then attributed to the *transformed* parameters and the output labels them as alpha, beta, gamma, and p rather than the internal raw names, so the results are directly interpretable in terms of the model parameters.

Path-level uncertainty is obtained by propagating node-level uncertainty draws through the path aggregation function:

$$P_k = 1 - \prod_{i=1}^{n_k} (1 - r_{k(v_i)}),$$

where $r_{k(v_i)}$ are node risks along path k .

All uncertainty metrics are computed from the first N Sobol draws (matrix A), while sensitivity indices use the full Sobol' design.

For more information about the uncertainty and sensitivity analysis and the output of this function, see the **sensobol** package (Puy et al. 2022).

The returned node table includes the following columns:

- name: name of the node.
- uncertainty_analysis: numeric vector of length N giving the uncertainty draws in the node risk score (from Sobol matrix A).
- sensitivity_analysis: object returned by `sensobol::sobol_indices()` for that node, containing Sobol' indices labelled alpha, beta, gamma, and p. These correspond to the normalised weights and the power-mean exponent respectively. The indices are computed on the transformed parameters (see Details).

The returned paths table includes:

- path_id: path identifier.
- path_str: sequence of function calls for each path.
- hops: number of edges.
- uncertainty_analysis: numeric vector giving the uncertainty draws in the path risk score.
- gini_index: numeric vector giving the uncertainty draws in the gini index.
- risk_trend: numeric vector giving the uncertainty draws in the risk trend.

Value

A named list with:

nodes A tibble of node results.

paths A tibble of path results. If `all_paths_out$paths` is empty (e.g., the graph has no entry-to-sink paths), an empty tibble with the same columns is returned.

References

Puy, A., Lo Piano, S., Saltelli, A., and Levin, S. A. (2022). *sensobol: An R Package to Compute Variance-Based Sensitivity Indices*. Journal of Statistical Software, 102(5), 1–37. doi:10.18637/jss.v102.i05

Examples

```
data(synthetic_graph)
out <- all_paths_fun(graph = synthetic_graph, alpha = 0.6, beta = 0.3,
                    gamma = 0.1, complexity_col = "cyclo")

# Power-mean risk (increase N to at least 2^10 for a proper UA/SA)
results <- uncertainty_fun(all_paths_out = out, N = 2^2, order = "first")

results$nodes
results$paths
```

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