

Package ‘couplr’

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Title Optimal Pairing and Matching via Linear Assignment

Version 1.2.1

Description Solves optimal pairing and matching problems using linear assignment algorithms. Provides implementations of the Hungarian method (Kuhn 1955) <doi:10.1002/nav.3800020109>, Jonker-Volgenant shortest path algorithm (Jonker and Volgenant 1987) <doi:10.1007/BF02278710>, Auction algorithm (Bertsekas 1988) <doi:10.1007/BF02186476>, cost-scaling (Goldberg and Kennedy 1995) <doi:10.1007/BF01585996>, scaling algorithms (Gabow and Tarjan 1989) <doi:10.1137/0218069>, push-relabel (Goldberg and Tarjan 1988) <doi:10.1145/48014.61051>, and Sinkhorn entropy-regularized transport (Cuturi 2013) <doi:10.48550/arxiv.1306.0895>. Designed for matching plots, sites, samples, or any pairwise optimization problem. Supports rectangular matrices, forbidden assignments, data frame inputs, batch solving, k-best solutions, and pixel-level image morphing for visualization. Includes automatic preprocessing with variable health checks, multiple scaling methods (standardized, range, robust), greedy matching algorithms, and comprehensive balance diagnostics for assessing match quality using standardized differences and distribution comparisons.

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assignment

*Linear assignment solver***Description**

Solve the linear assignment problem (minimum- or maximum-cost matching) using several algorithms. Forbidden edges can be marked as NA or Inf.

Usage

```
assignment(
  cost,
  maximize = FALSE,
  method = c("auto", "jv", "hungarian", "auction", "auction_gs", "auction_scaled", "sap",
    "ssp", "csflow", "hk01", "bruteforce", "ssap_bucket", "cycle_cancel", "gabow_tarjan",
    "lapmod", "csa", "ramshaw_tarjan", "push_relabel", "orlin", "network_simplex"),
  auction_eps = NULL,
  eps = NULL
)
```

Arguments

cost	Numeric matrix; rows = tasks, columns = agents. NA or Inf entries are treated as forbidden assignments.
maximize	Logical; if TRUE, maximizes the total cost instead of minimizing.
method	Character string indicating the algorithm to use. Options: General-purpose solvers: <ul style="list-style-type: none"> • "auto" — Automatic selection based on problem characteristics (default) • "jv" — 'Jonker-Volgenant', fast general-purpose $O(n^3)$ • "hungarian" — Classic 'Hungarian' algorithm $O(n^3)$ Auction-based solvers: <ul style="list-style-type: none"> • "auction" — 'Bertsekas' auction with adaptive epsilon • "auction_gs" — 'Gauss-Seidel' variant, good for spatial structure • "auction_scaled" — 'Epsilon-scaling', fastest for large dense problems Specialized solvers: <ul style="list-style-type: none"> • "sap" / "ssp" — Shortest augmenting path, handles sparsity well • "lapmod" — Sparse JV variant, faster when $>50\%$ • "hk01" — 'Hopcroft-Karp' for binary (0/1) costs only • "ssap_bucket" — 'Dial' algorithm for integer costs • "line_metric" — $O(n \log n)$ for 1D assignment problems • "bruteforce" — Exact enumeration for tiny problems ($n \leq 8$) Advanced solvers: <ul style="list-style-type: none"> • "csa" — 'Goldberg-Kennedy' cost-scaling, often fastest for medium-large • "gabow_tarjan" — 'Gabow-Tarjan' bit-scaling with complementary slackness $O(n^3 \log C)$ • "cycle_cancel" — Cycle-canceling with 'Karp' algorithm • "csflow" — Cost-scaling network flow • "network_simplex" — 'Network simplex' with spanning tree representation • "orlin" — 'Orlin-Ahuja' scaling $O(\sqrt{n} * m * \log(nC))$ • "push_relabel" — 'Push-relabel' max-flow based solver

	<ul style="list-style-type: none"> • "ramshaw_tarjan" — 'Ramshaw-Tarjan', optimized for rectangular matrices ($n \neq m$)
auction_eps	Optional numeric epsilon for the 'Auction'/'Auction-GS' methods. If NULL, an internal default (e.g., $1e-9$) is used.
eps	Deprecated. Use auction_eps. If provided and auction_eps is NULL, its value is used for auction_eps.

Details

method = "auto" selects an algorithm based on problem size/shape and data characteristics:

- Very small ($n \leq 8$): "bruteforce" — exact enumeration
- Binary/constant costs: "hk01" — specialized for 0/1 costs
- Large sparse ($n > 100, > 50$)
- Sparse or very rectangular: "sap" — handles sparsity well
- Small-medium ($8 < n \leq 50$): "hungarian" — provides exact dual solutions
- Medium ($50 < n \leq 75$): "jv" — fast general-purpose solver
- Large ($n > 75$): "auction_scaled" — fastest for large dense problems

Benchmarks show 'Auction-scaled' and 'JV' are 100-1500x faster than 'Hungarian' at $n=500$.

Value

An object of class lap_solve_result, a list with elements:

- match — integer vector of length $\min(\text{nrow}(\text{cost}), \text{ncol}(\text{cost}))$ giving the assigned column for each row (0 if unassigned).
- total_cost — numeric scalar, the objective value.
- status — character scalar, e.g. "optimal".
- method_used — character scalar, the algorithm actually used.

See Also

- [lap_solve\(\)](#) — Tidy interface returning tibbles
- [lap_solve_kbest\(\)](#) — Find k-best assignments ('Murty' algorithm)
- [assignment_duals\(\)](#) — Extract dual variables for sensitivity analysis
- [bottleneck_assignment\(\)](#) — Minimize maximum edge cost (minimax)
- [sinkhorn\(\)](#) — Entropy-regularized optimal transport

Examples

```
cost <- matrix(c(4,2,5, 3,3,6, 7,5,4), nrow = 3, byrow = TRUE)
res <- assignment(cost)
res$match; res$total_cost
```

assignment_duals *Solve assignment problem and return dual variables*

Description

Solves the linear assignment problem and returns dual potentials (u, v) in addition to the optimal matching. The dual variables provide an optimality certificate and enable sensitivity analysis.

Usage

```
assignment_duals(cost, maximize = FALSE)
```

Arguments

cost	Numeric matrix; rows = tasks, columns = agents. NA or Inf entries are treated as forbidden assignments.
maximize	Logical; if TRUE, maximizes the total cost instead of minimizing.

Details

The dual variables satisfy the complementary slackness conditions:

- For minimization: $u[i] + v[j] \leq \text{cost}[i, j]$ for all (i,j)
- For any assigned pair (i,j): $u[i] + v[j] = \text{cost}[i, j]$

This implies that $\text{sum}(u) + \text{sum}(v) = \text{total_cost}$ (strong duality).

Applications of dual variables:

- **Optimality verification:** Check that duals satisfy constraints
- **Sensitivity analysis:** Reduced cost $c[i, j] - u[i] - v[j]$ shows how much an edge cost must decrease before it enters the solution
- **Pricing in column generation:** Use duals to price new columns
- **Warm starting:** Reuse duals when costs change slightly

Value

A list with class "assignment_duals_result" containing:

- match - integer vector of column assignments (1-based)
- total_cost - optimal objective value
- u - numeric vector of row dual variables (length n)
- v - numeric vector of column dual variables (length m)
- status - character, e.g. "optimal"

See Also

[assignment\(\)](#) for standard assignment without duals

Examples

```

cost <- matrix(c(4, 2, 5, 3, 3, 6, 7, 5, 4), nrow = 3, byrow = TRUE)
result <- assignment_duals(cost)

# Check optimality: u + v should equal cost for assigned pairs
for (i in 1:3) {
  j <- result$match[i]
  cat(sprintf("Row %d -> Col %d: u + v = %.2f, cost = %.2f\n",
            i, j, result$u[i] + result$v[j], cost[i, j]))
}

# Verify strong duality
cat("sum(u) + sum(v) =", sum(result$u) + sum(result$v), "\n")
cat("total_cost =", result$total_cost, "\n")

# Reduced costs (how much must cost decrease to enter solution)
reduced <- outer(result$u, result$v, "+")
reduced_cost <- cost - reduced
print(round(reduced_cost, 2))

```

as_assignment_matrix *Convert assignment result to a binary matrix*

Description

Turns a tidy assignment result back into a 0/1 assignment matrix.

Usage

```
as_assignment_matrix(x, n_sources = NULL, n_targets = NULL)
```

Arguments

x	An assignment result object of class lap_solve_result
n_sources	Number of source nodes, optional
n_targets	Number of target nodes, optional

Value

Integer matrix with 0 and 1 entries

as_matchit

*Convert couplr Result to matchit Object***Description**

Constructs a matchit-class S3 object from a couplr result, enabling use with any function that accepts **MatchIt** objects (e.g., **cobalt**, **marginaleffects**).

Usage

```
as_matchit(
  result,
  left,
  right,
  formula = NULL,
  left_id = "id",
  right_id = "id",
  ...
)
```

Arguments

result	A couplr result object (matching_result, full_matching_result, cem_result, or subclass_result)
left	Data frame of left (treated) units
right	Data frame of right (control) units
formula	Optional formula used for matching. If not provided, a default formula is constructed from result\$info\$vars.
left_id	Name of ID column in left (default: "id")
right_id	Name of ID column in right (default: "id")
...	Additional arguments (ignored)

Value

An S3 object of class "matchit" with fields:

match.matrix Match matrix (treated x controls)

treat Named treatment vector (1/0)

weights Matching weights

X Covariate matrix

call Original call

info Metadata from couplr

Examples

```

left <- data.frame(id = 1:5, age = c(25, 35, 45, 55, 65))
right <- data.frame(id = 6:15, age = runif(10, 20, 70))
result <- match_couples(left, right, vars = "age")
mi <- as_matchit(result, left, right)
if (requireNamespace("cobalt", quietly = TRUE)) {
  cobalt::bal.tab(mi)
}

```

augment

Generic Augment Function

Description

S3 generic for augmenting model results with original data.

Usage

```
augment(x, ...)
```

Arguments

x An object to augment
... Additional arguments passed to methods

Value

Augmented data (depends on method)

augment.matching_result

Augment Matching Results with Original Data (broom-style)

Description

S3 method for augmenting matching results following the broom package conventions. This is a thin wrapper around `join_matched()` with sensible defaults for quick exploration.

Usage

```

## S3 method for class 'matching_result'
augment(x, left, right, ...)

```

Arguments

<code>x</code>	A <code>matching_result</code> object
<code>left</code>	The original left dataset
<code>right</code>	The original right dataset
<code>...</code>	Additional arguments passed to <code>join_matched()</code>

Details

This method follows the `augment()` convention from the `broom` package, making it easy to integrate `couplr` into `tidymodels` workflows. It's equivalent to calling `join_matched()` with default parameters.

If the `broom` package is not loaded, you can use `couplr::augment()` to access this function.

Value

A tibble with matched pairs and original data (see `join_matched()`)

Examples

```
left <- data.frame(
  id = 1:5,
  treatment = 1,
  age = c(25, 30, 35, 40, 45)
)

right <- data.frame(
  id = 6:10,
  treatment = 0,
  age = c(24, 29, 36, 41, 44)
)

result <- match_couples(left, right, vars = "age")
couplr::augment(result, left, right)
```

autoplot.balance_diagnostics
ggplot2 autoplot for balance diagnostics

Description

Produces `ggplot2`-based balance assessment plots. Returns a `ggplot` object.

Usage

```
## S3 method for class 'balance_diagnostics'
autoplot(
  object,
  type = c("love", "histogram", "variance"),
  threshold = 0.1,
  ...
)
```

Arguments

object	A balance_diagnostics object
type	Type of plot: "love" (default), "histogram", or "variance"
threshold	Threshold for standardized differences (default: 0.1)
...	Additional arguments (ignored)

Value

A ggplot object

Examples

```
if (requireNamespace("ggplot2", quietly = TRUE)) {
  set.seed(42)
  left <- data.frame(id = 1:10, age = rnorm(10, 45, 10),
                    income = rnorm(10, 50000, 15000))
  right <- data.frame(id = 11:30, age = rnorm(20, 47, 10),
                    income = rnorm(20, 52000, 15000))
  result <- match_couples(left, right, vars = c("age", "income"))
  bal <- balance_diagnostics(result, left, right, vars = c("age", "income"))
  ggplot2::autoplot(bal)
}
```

autoplot.matching_result

ggplot2 autoplot for matching results

Description

Produces ggplot2-based visualizations of matching distance distributions. Returns a ggplot object that can be further customized.

Usage

```
## S3 method for class 'matching_result'
autoplot(object, type = c("histogram", "density", "ecdf"), ...)
```

Arguments

object	A matching_result object
type	Type of plot: "histogram" (default), "density", or "ecdf"
...	Additional arguments (ignored)

Details

Use plot() for base graphics or autoplot() for ggplot2 output. The ggplot2 package must be installed.

Value

A ggplot object

Examples

```
if (requireNamespace("ggplot2", quietly = TRUE)) {
  left <- data.frame(id = 1:5, x = c(1, 2, 3, 4, 5))
  right <- data.frame(id = 6:10, x = c(1.1, 2.2, 3.1, 4.2, 5.1))
  result <- match_couples(left, right, vars = "x")
  ggplot2::autoplot(result)
  ggplot2::autoplot(result, type = "density")
}
```

autoplot.sensitivity_analysis

ggplot2 autoplot for sensitivity analysis

Description

Plots p-value upper bounds against sensitivity parameter Gamma.

Usage

```
## S3 method for class 'sensitivity_analysis'
autoplot(object, alpha = 0.05, ...)
```

Arguments

object	A sensitivity_analysis object
alpha	Significance level (default: 0.05)
...	Additional arguments (ignored)

Value

A ggplot object

`bal.tab.matching_result`*Balance Table for Matching Results (cobalt integration)*

Description

S3 method enabling `cobalt::bal.tab()` on couplr result objects. Requires the **cobalt** package to be installed.

Usage

```
bal.tab.matching_result(x, left, right, ...)
```

```
bal.tab.full_matching_result(x, left, right, ...)
```

```
bal.tab.cem_result(x, left, right, ...)
```

```
bal.tab.subclass_result(x, data = NULL, ...)
```

Arguments

<code>x</code>	A couplr result object
<code>left</code>	Data frame of left (treated) units
<code>right</code>	Data frame of right (control) units
<code>...</code>	Additional arguments passed to <code>cobalt::bal.tab()</code>
<code>data</code>	Data frame used for subclassification (for <code>subclass_result</code> only)

Details

These methods convert couplr results to the format cobalt expects (a `matchit`-class object) and then delegate to cobalt's own `bal.tab.matchit()` method. The **cobalt** package must be installed but is not required for couplr to function.

Value

A cobalt balance table object

balance_diagnostics *Balance Diagnostics for Matched Pairs*

Description

Computes comprehensive balance statistics comparing the distribution of matching variables between left and right units in the matched sample.

Usage

```
balance_diagnostics(result, ...)  
  
## S3 method for class 'matching_result'  
balance_diagnostics(  
  result,  
  left,  
  right,  
  vars = NULL,  
  left_id = "id",  
  right_id = "id",  
  ...  
)  
  
## S3 method for class 'full_matching_result'  
balance_diagnostics(  
  result,  
  left,  
  right,  
  vars = NULL,  
  left_id = "id",  
  right_id = "id",  
  ...  
)  
  
## S3 method for class 'cem_result'  
balance_diagnostics(  
  result,  
  left,  
  right,  
  vars = NULL,  
  left_id = "id",  
  right_id = "id",  
  ...  
)  
  
## S3 method for class 'subclass_result'  
balance_diagnostics(result, data = NULL, vars = NULL, ...)
```

Arguments

<code>result</code>	A matching result object from <code>match_couples()</code> , <code>greedy_couples()</code> , <code>full_match()</code> , <code>cem_match()</code> , or <code>subclass_match()</code>
<code>...</code>	Additional arguments passed to methods
<code>left</code>	Data frame of left units
<code>right</code>	Data frame of right units
<code>vars</code>	Character vector of variable names to check balance for. Defaults to the variables used in matching (if available in result).
<code>left_id</code>	Character, name of ID column in left data (default: "id")
<code>right_id</code>	Character, name of ID column in right data (default: "id")
<code>data</code>	Data frame used for subclassification (when <code>result</code> is a <code>subclass_result</code>)

Details

This function computes several balance metrics:

Standardized Difference: The difference in means divided by the pooled standard deviation. Values less than 0.1 indicate excellent balance, 0.1-0.25 good balance.

Variance Ratio: The ratio of standard deviations (left/right). Values close to 1 are ideal.

KS Statistic: Kolmogorov-Smirnov test statistic comparing distributions. Lower values indicate more similar distributions.

Overall Metrics include mean absolute standardized difference across all variables, proportion of variables with large imbalance (`lstddiff > 0.25`), and maximum standardized difference.

Value

An S3 object of class `balance_diagnostics` containing:

var_stats Tibble with per-variable balance statistics

overall List with overall balance metrics

pairs Tibble of matched pairs with variables

n_matched Number of matched pairs

n_unmatched_left Number of unmatched left units

n_unmatched_right Number of unmatched right units

method Matching method used

has_blocks Whether blocking was used

block_stats Per-block statistics (if blocking used)

Examples

```
# Create sample data
set.seed(123)
left <- data.frame(
  id = 1:10,
  age = rnorm(10, 45, 10),
  income = rnorm(10, 50000, 15000)
)
right <- data.frame(
  id = 11:30,
  age = rnorm(20, 47, 10),
  income = rnorm(20, 52000, 15000)
)

# Match
result <- match_couples(left, right, vars = c("age", "income"))

# Get balance diagnostics
balance <- balance_diagnostics(result, left, right, vars = c("age", "income"))
print(balance)

# Get balance table
balance_table(balance)
```

balance_table	<i>Create Balance Table</i>
---------------	-----------------------------

Description

Formats balance diagnostics into a clean table for display or export.

Usage

```
balance_table(balance, digits = 3)
```

Arguments

balance	A balance_diagnostics object from balance_diagnostics()
digits	Number of decimal places for rounding (default: 3)

Value

A tibble with formatted balance statistics

bottleneck_assignment *Solve the Bottleneck Assignment Problem*

Description

Finds an assignment that minimizes (or maximizes) the maximum edge cost in a perfect matching. Unlike standard LAP which minimizes the sum of costs, BAP minimizes the maximum (bottleneck) cost.

Usage

```
bottleneck_assignment(cost, maximize = FALSE)
```

Arguments

cost	Numeric matrix; rows = tasks, columns = agents. NA or Inf entries are treated as forbidden assignments.
maximize	Logical; if TRUE, maximizes the minimum edge cost instead of minimizing the maximum (maximin objective). Default is FALSE (minimax).

Details

The Bottleneck Assignment Problem (BAP) is a variant of the Linear Assignment Problem where instead of minimizing the sum of assignment costs, we minimize the maximum cost among all assignments (minimax objective).

Algorithm: Uses binary search on the sorted unique costs combined with 'Hopcroft-Karp' bipartite matching to find the minimum threshold that allows a perfect matching.

Complexity: $O(E * \sqrt{V} * \log(\text{unique costs}))$ where $E = \text{edges}$, $V = \text{vertices}$.

Applications:

- Task scheduling with deadline constraints (minimize latest completion)
- Resource allocation (minimize maximum load/distance)
- Network routing (minimize maximum link utilization)
- Fair division problems (minimize maximum disparity)

Value

A list with class "bottleneck_result" containing:

- match - integer vector of length `nrow(cost)` giving the assigned column for each row (1-based indexing)
- bottleneck - numeric scalar, the bottleneck (max/min edge) value
- status - character scalar, e.g. "optimal"

See Also

[assignment\(\)](#) for standard LAP (sum objective), [lap_solve\(\)](#) for tidy LAP interface

Examples

```
# Simple example: minimize max cost
cost <- matrix(c(1, 5, 3,
                2, 4, 6,
                7, 1, 2), nrow = 3, byrow = TRUE)
result <- bottleneck_assignment(cost)
result$bottleneck # Maximum edge cost in optimal assignment

# Maximize minimum (fair allocation)
profits <- matrix(c(10, 5, 8,
                  6, 12, 4,
                  3, 7, 11), nrow = 3, byrow = TRUE)
result <- bottleneck_assignment(profits, maximize = TRUE)
result$bottleneck # Minimum profit among all assignments

# With forbidden assignments
cost <- matrix(c(1, NA, 3,
                2, 4, Inf,
                5, 1, 2), nrow = 3, byrow = TRUE)
result <- bottleneck_assignment(cost)
```

cardinality_match *Cardinality Matching*

Description

Maximizes the number of matched pairs subject to balance constraints. Uses iterative pruning: starts with a full match, then removes pairs that contribute most to imbalance until all variables satisfy the standardized difference threshold.

Usage

```
cardinality_match(
  left,
  right,
  vars,
  max_std_diff = 0.1,
  distance = "euclidean",
  weights = NULL,
  scale = FALSE,
  auto_scale = FALSE,
  method = "auto",
  max_iter = 100L,
  batch_fraction = 0.1
)
```

Arguments

<code>left</code>	Data frame of "left" units
<code>right</code>	Data frame of "right" units
<code>vars</code>	Character vector of matching variable names
<code>max_std_diff</code>	Maximum allowed absolute standardized difference (default: 0.1, corresponding to "excellent" balance)
<code>distance</code>	Distance metric (default: "euclidean")
<code>weights</code>	Optional named vector of variable weights
<code>scale</code>	Scaling method (default: FALSE)
<code>auto_scale</code>	If TRUE, automatically select scaling (default: FALSE)
<code>method</code>	LAP solver method (default: "auto")
<code>max_iter</code>	Maximum pruning iterations (default: 100)
<code>batch_fraction</code>	Fraction of worst pairs to remove per iteration (default: 0.1). Larger values speed up convergence but may over-prune.

Details

Cardinality matching (Zubizarreta 2014) finds the largest matched sample that satisfies pre-specified balance constraints. This implementation uses an iterative pruning heuristic:

1. Run full optimal matching via `match_couples()`
2. Compute balance diagnostics
3. While any `lstd_diff` exceeds `max_std_diff`:
 - Identify the variable with worst balance
 - Remove the batch of pairs contributing most to that imbalance
 - Recompute balance

Value

A `matching_result` object with additional cardinality info: `result$info$pruning_iterations`, `result$info$pairs_removed`, `result$info$final_balance`.

Examples

```
set.seed(42)
left <- data.frame(id = 1:20, x = rnorm(20, 0, 1), y = rnorm(20, 0, 1))
right <- data.frame(id = 21:50, x = rnorm(30, 0.5, 1), y = rnorm(30, 0.3, 1))
result <- cardinality_match(left, right, vars = c("x", "y"), max_std_diff = 0.2)
print(result)
```

cem_match

*Coarsened Exact Matching***Description**

Coarsens continuous variables into bins, then performs exact matching on the coarsened values. Units in strata containing both left and right units are kept; others are pruned. Matched units receive weights inversely proportional to stratum sizes to maintain balance.

Usage

```
cem_match(
  left,
  right,
  vars,
  cutpoints = NULL,
  n_bins = "sturges",
  grouping = NULL,
  keep = "all",
  left_id = "id",
  right_id = "id"
)
```

Arguments

left	Data frame of left (treated) units
right	Data frame of right (control) units
vars	Character vector of variable names to coarsen and match on
cutpoints	Named list of break vectors per variable. If NULL, automatic binning is used.
n_bins	Binning method when cutpoints is NULL: "sturges" (default), "fd" (Freedman-Diaconis), "scott", or an integer specifying the number of bins for all variables.
grouping	Character vector of variable names to match exactly (without coarsening). These are typically categorical variables.
keep	Which units to return: "all" (default) returns all units with weight 0 for unmatched, "matched" drops unmatched units.
left_id	Name of ID column in left (default: "id")
right_id	Name of ID column in right (default: "id")

Details

CEM algorithm:

1. Coarsen each numeric variable using `cut` with either user-specified breakpoints or automatic binning (Sturges, FD, or Scott rule)
2. Categorical variables in `grouping` are kept as-is

3. Create strata by concatenating all coarsened values
4. Drop strata with 0 left or 0 right units
5. Compute CEM weights: left units get weight 1, right units get weight $n_{\text{left_in_stratum}} / n_{\text{right_in_stratum}}$ so that the total weight of right units in each stratum equals the number of left units

Value

An S3 object of class `c("cem_result", "couplr_result")` containing:

matched Tibble with columns `id`, `side`, `stratum`, `weight`

strata_summary Tibble with per-stratum counts

info List with `n_strata`, `n_matched_left`, `n_matched_right`, `n_pruned_left`, `n_pruned_right`, `method`, `vars`

Examples

```
set.seed(42)
left <- data.frame(
  id = 1:20, age = rnorm(20, 40, 10),
  income = rnorm(20, 50000, 10000)
)
right <- data.frame(
  id = 21:60, age = rnorm(40, 42, 10),
  income = rnorm(40, 52000, 10000)
)
result <- cem_match(left, right, vars = c("age", "income"))
print(result)
```

compute_distances

Compute and Cache Distance Matrix for Reuse

Description

Precomputes a distance matrix between left and right datasets, allowing it to be reused across multiple matching operations with different constraints. This is particularly useful when exploring different matching parameters (`max_distance`, `calipers`, `methods`) without recomputing distances.

Usage

```
compute_distances(
  left,
  right,
  vars,
  distance = "euclidean",
  weights = NULL,
```

```

scale = FALSE,
auto_scale = FALSE,
left_id = "id",
right_id = "id",
block_id = NULL
)

```

Arguments

left	Left dataset (data frame)
right	Right dataset (data frame)
vars	Character vector of variable names to use for distance computation
distance	Distance metric (default: "euclidean")
weights	Optional numeric vector of variable weights
scale	Scaling method: FALSE, "standardize", "range", or "robust"
auto_scale	Apply automatic preprocessing (default: FALSE)
left_id	Name of ID column in left (default: "id")
right_id	Name of ID column in right (default: "id")
block_id	Optional block ID column name for blocked matching

Details

This function computes distances once and stores them in a reusable object. The resulting `distance_object` can be passed to `match_couples()` or `greedy_couples()` instead of providing datasets and variables.

Benefits:

- **Performance:** Avoid recomputing distances when trying different constraints
- **Exploration:** Quickly test `max_distance`, `calipers`, or `methods`
- **Consistency:** Ensures same distances used across comparisons
- **Memory efficient:** Can use sparse matrices when many pairs are forbidden

The `distance_object` stores the original datasets, allowing downstream functions like `join_matched()` to work seamlessly.

Value

An S3 object of class "distance_object" containing:

- `cost_matrix`: Numeric matrix of distances
- `left_ids`: Character vector of left IDs
- `right_ids`: Character vector of right IDs
- `block_id`: Block ID column name (if specified)
- `metadata`: List with computation details (`vars`, `distance`, `scale`, etc.)
- `original_left`: Original left dataset (for later joining)
- `original_right`: Original right dataset (for later joining)

Examples

```
# Compute distances once
left <- data.frame(id = 1:5, age = c(25, 30, 35, 40, 45), income = c(45, 52, 48, 61, 55) * 1000)
right <- data.frame(id = 6:10, age = c(24, 29, 36, 41, 44), income = c(46, 51, 47, 60, 54) * 1000)

dist_obj <- compute_distances(
  left, right,
  vars = c("age", "income"),
  scale = "standardize"
)

# Reuse for different matching strategies
result1 <- match_couples(dist_obj, max_distance = 0.5)
result2 <- match_couples(dist_obj, max_distance = 1.0)
result3 <- greedy_couples(dist_obj, strategy = "sorted")

# All use the same precomputed distances
```

diagnose_distance_matrix

Diagnose distance matrix and suggest fixes

Description

Comprehensive diagnostics for a distance matrix with actionable suggestions.

Usage

```
diagnose_distance_matrix(
  cost_matrix,
  left = NULL,
  right = NULL,
  vars = NULL,
  warn = TRUE
)
```

Arguments

cost_matrix	Numeric matrix of distances
left	Left dataset (for variable checking)
right	Right dataset (for variable checking)
vars	Variables used for matching
warn	If TRUE, issue warnings

Value

List with diagnostic results and suggestions

`example_costs`*Example cost matrices for assignment problems*

Description

Small example datasets for demonstrating couplr functionality across different assignment problem types: square, rectangular, sparse, and binary.

Usage

```
example_costs
```

Format

A list containing four example cost matrices:

simple_3x3 A 3x3 cost matrix with costs ranging from 2-7. Optimal assignment: row 1 -> col 2 (cost 2), row 2 -> col 1 (cost 3), row 3 -> col 3 (cost 4). Total optimal cost: 9.

rectangular_3x5 A 3x5 rectangular cost matrix demonstrating assignment when rows < columns. Each of 3 rows is assigned to one of 5 columns; 2 columns remain unassigned. Costs range 1-6.

sparse_with_na A 3x3 matrix with NA values indicating forbidden assignments. Use this to test algorithms' handling of constraints. Position (1,3), (2,2), and (3,1) are forbidden.

binary_costs A 3x3 matrix with binary (0/1) costs, suitable for testing the HK01 algorithm. Diagonal entries are 0 (preferred), off-diagonal entries are 1 (penalty).

Details

These matrices are designed to test different aspects of LAP solvers:

simple_3x3: Basic functionality test. Any correct solver should find total cost = 9.

rectangular_3x5: Tests handling of non-square problems. The optimal solution assigns all 3 rows with minimum total cost.

sparse_with_na: Tests constraint handling. Algorithms must avoid NA positions while finding an optimal assignment among valid entries.

binary_costs: Tests specialized binary cost algorithms. The optimal assignment uses all diagonal entries (total cost = 0).

See Also

[lap_solve](#), [example_df](#)

Examples

```
# Simple 3x3 assignment
result <- lap_solve(example_costs$simple_3x3)
print(result)
# Optimal: sources 1,2,3 -> targets 2,1,3 with cost 9

# Rectangular problem (3 sources, 5 targets)
result <- lap_solve(example_costs$rectangular_3x5)
print(result)
# All 3 sources assigned; 2 targets unassigned

# Sparse problem with forbidden assignments
result <- lap_solve(example_costs$sparse_with_na)
print(result)
# Avoids NA positions

# Binary costs - test HK01 algorithm
result <- lap_solve(example_costs$binary_costs, method = "hk01")
print(result)
# Finds diagonal assignment (cost = 0)
```

example_df

Example assignment problem data frame

Description

A tidy data frame representation of assignment problems, suitable for use with grouped workflows and batch solving. Contains two independent 3x3 assignment problems in long format.

Usage

```
example_df
```

Format

A tibble with 18 rows and 4 columns:

sim Simulation/problem identifier. Integer with values 1 or 2, distinguishing two independent assignment problems. Use with `group_by(sim)` for grouped solving.

source Source node index. Integer 1-3 representing the row (source) in each 3x3 cost matrix.

target Target node index. Integer 1-3 representing the column (target) in each 3x3 cost matrix.

cost Cost of assigning source to target. Numeric values ranging from 1-7. Each source-target pair has exactly one cost entry.

Details

This dataset demonstrates couplr's data frame interface for LAP solving. The long format (one row per source-target pair) is converted internally to a cost matrix for solving.

Simulation 1: Costs from `example_costs$simple_3x3`

- Optimal assignment: (1->2, 2->1, 3->3)
- Total cost: 9

Simulation 2: Different cost structure

- Optimal assignment: (1->1, 2->3, 3->3) or equivalent
- Total cost: 4

See Also

[lap_solve](#), [lap_solve_batch](#), [example_costs](#)

Examples

```
library(dplyr)

# Solve both problems with grouped workflow
example_df |>
  group_by(sim) |>
  lap_solve(source, target, cost)

# Batch solving for efficiency
example_df |>
  group_by(sim) |>
  lap_solve_batch(source, target, cost)

# Inspect the data structure
example_df |>
  group_by(sim) |>
  summarise(
    n_pairs = n(),
    min_cost = min(cost),
    max_cost = max(cost)
  )
```

full_match

Full Matching

Description

Assigns every unit (left and right) to a matched group with variable ratios (1:k or k:1). Unlike 1:1 matching, full matching does not discard units, producing matched groups where each group contains at least one left and one right unit.

Usage

```

full_match(
  left,
  right,
  vars,
  distance = "euclidean",
  min_controls = 1,
  max_controls = Inf,
  caliper = NULL,
  caliper_sd = NULL,
  weights = NULL,
  scale = FALSE,
  auto_scale = FALSE,
  sigma = NULL,
  left_id = "id",
  right_id = "id",
  method = "optimal"
)

```

Arguments

left	Data frame of left (treated) units
right	Data frame of right (control) units
vars	Character vector of variable names to match on
distance	Distance metric: "euclidean" (default), "mahalanobis", "manhattan", or a custom function
min_controls	Minimum number of right units per group (default: 1)
max_controls	Maximum number of right units per group (default: Inf)
caliper	Maximum allowable distance for a match. Units with no eligible partner within the caliper are left unmatched.
caliper_sd	If not NULL, caliper is expressed in standard deviations of the pooled distance distribution rather than absolute units.
weights	Named numeric vector of variable weights
scale	Scaling method: FALSE (default), "robust", "standardize", or "range"
auto_scale	If TRUE, automatically preprocess and scale variables
sigma	Optional covariance matrix for Mahalanobis distance
left_id	Name of ID column in left (default: "id")
right_id	Name of ID column in right (default: "id")
method	Matching algorithm: "optimal" (default) uses min-cost max-flow to find the globally optimal group assignment minimizing total distance; "greedy" uses a fast two-pass heuristic.

Details

Full matching creates matched groups of variable size. Two algorithms are available:

Optimal (method = "optimal", default): Solves a min-cost max-flow problem that minimizes total distance across all group assignments simultaneously. Each left unit becomes a group center absorbing 1 to max_controls right units, with the globally optimal assignment found via Dijkstra's algorithm with Johnson potentials. When n_left > n_right, roles are transposed automatically.

Greedy (method = "greedy"): A fast two-pass heuristic:

1. Each left unit picks its nearest eligible right unit
2. Remaining right units are assigned to their nearest already-matched left unit, respecting max_controls

This is faster but does not guarantee globally optimal results.

Weights are computed so that within each group, the total weight of right units equals the total weight of left units (which is 1). For a group with 1 left and k right units, each right unit receives weight 1/k.

Value

An S3 object of class c("full_matching_result", "couplr_result") containing:

groups Tibble with columns group_id, id, side ("left"/"right"), and weight

info List with n_groups, n_left, n_right, n_unmatched_left, n_unmatched_right, method, vars

unmatched List of unmatched left and right IDs (if caliper excludes units)

Examples

```
set.seed(42)
left <- data.frame(id = 1:5, age = c(25, 35, 45, 55, 65))
right <- data.frame(id = 6:20, age = runif(15, 20, 70))
result <- full_match(left, right, vars = "age")
print(result)
```

get_method_used

Extract method used from assignment result

Description

Extract method used from assignment result

Usage

```
get_method_used(x)
```

Arguments

x An assignment result object

Value

Character string indicating method used

get_total_cost	<i>Extract total cost from assignment result</i>
----------------	--

Description

Extract total cost from assignment result

Usage

```
get_total_cost(x)
```

Arguments

x An assignment result object

Value

Numeric total cost

greedy_couples	<i>Fast approximate matching using greedy algorithm</i>
----------------	---

Description

Performs fast one-to-one matching using greedy strategies. Does not guarantee optimal total distance but is much faster than `match_couples()` for large datasets. Supports blocking, distance constraints, and various distance metrics.

Usage

```
greedy_couples(
  left,
  right = NULL,
  vars = NULL,
  distance = "euclidean",
  weights = NULL,
  scale = FALSE,
  auto_scale = FALSE,
```

```

max_distance = Inf,
calipers = NULL,
block_id = NULL,
ignore_blocks = FALSE,
require_full_matching = FALSE,
strategy = c("row_best", "sorted", "pq"),
return_unmatched = TRUE,
return_diagnostics = FALSE,
parallel = FALSE,
replace = FALSE,
ratio = 1L,
check_costs = TRUE,
sigma = NULL
)

```

Arguments

<code>left</code>	Data frame of "left" units (e.g., treated, cases)
<code>right</code>	Data frame of "right" units (e.g., control, controls)
<code>vars</code>	Variable names to use for distance computation
<code>distance</code>	Distance metric: "euclidean", "manhattan", "mahalanobis", or a custom function
<code>weights</code>	Optional named vector of variable weights
<code>scale</code>	Scaling method: FALSE (none), "standardize", "range", or "robust"
<code>auto_scale</code>	If TRUE, automatically check variable health and select scaling method (default: FALSE)
<code>max_distance</code>	Maximum allowed distance (pairs exceeding this are forbidden)
<code>calipers</code>	Named list of per-variable maximum absolute differences
<code>block_id</code>	Column name containing block IDs (for stratified matching)
<code>ignore_blocks</code>	If TRUE, ignore <code>block_id</code> even if present
<code>require_full_matching</code>	If TRUE, error if any units remain unmatched
<code>strategy</code>	Greedy strategy: <ul style="list-style-type: none"> • "row_best": For each row, find best available column (default) • "sorted": Sort all pairs by distance, greedily assign • "pq": Use priority queue (good for very large problems)
<code>return_unmatched</code>	Include unmatched units in output
<code>return_diagnostics</code>	Include detailed diagnostics in output
<code>parallel</code>	Enable parallel processing for blocked matching. Requires 'future' and 'future.apply' packages. Can be: <ul style="list-style-type: none"> • FALSE: Sequential processing (default) • TRUE: Auto-configure parallel backend

	<ul style="list-style-type: none"> • Character: Specify future plan (e.g., "multisession", "multicore")
replace	If TRUE, allow matching with replacement (same right unit can be matched to multiple left units). Default: FALSE.
ratio	Integer, number of right units to match per left unit. Default: 1 (one-to-one matching). For k:1 matching, set ratio = k.
check_costs	If TRUE, check distance distribution for potential problems and provide helpful warnings before matching (default: TRUE)
sigma	Optional covariance matrix for Mahalanobis distance. If NULL (default), the pooled sample covariance is used. Only relevant when distance = "mahalanobis".

Details

Greedy strategies do not guarantee optimal total distance but are much faster:

- "row_best": $O(n*m)$ time, simple and often produces good results
- "sorted": $O(nm \log(n*m))$ time, better quality but slower
- "pq": $O(nm \log(n*m))$ time, memory-efficient for large problems

Use greedy_couples when:

- Dataset is very large ($> 10,000 \times 10,000$)
- Approximate solution is acceptable
- Speed is more important than optimality

Value

A list with class "matching_result" (same structure as match_couples)

Examples

```
# Basic greedy matching
left <- data.frame(id = 1:100, x = rnorm(100))
right <- data.frame(id = 101:200, x = rnorm(100))
result <- greedy_couples(left, right, vars = "x")

# Compare to optimal
result_opt <- match_couples(left, right, vars = "x")
result_greedy <- greedy_couples(left, right, vars = "x")
result_greedy$info$total_distance / result_opt$info$total_distance # Quality ratio
```

hospital_staff	<i>Hospital staff scheduling example dataset</i>
----------------	--

Description

A comprehensive example dataset for demonstrating couplr functionality across vignettes. Contains hospital staff scheduling data with nurses, shifts, costs, and preference scores suitable for assignment problems, as well as nurse characteristics for matching workflows.

Usage

```
hospital_staff
```

Format

A list containing eight related datasets:

basic_costs A 10x10 numeric cost matrix for assigning 10 nurses to 10 shifts. Values range from approximately 1-15, where lower values indicate better fit (less overtime, matches skills, respects preferences). Use with `lap_solve()` for basic assignment.

preferences A 10x10 numeric preference matrix on a 0-10 scale, where higher values indicate stronger nurse preference for a shift. Use with `lap_solve(..., maximize = TRUE)` to optimize preferences rather than minimize costs.

schedule_df A tibble with 100 rows (10 nurses x 10 shifts) in long format for data frame workflows:

nurse_id Integer 1-10. Unique identifier for each nurse.

shift_id Integer 1-10. Unique identifier for each shift.

cost Numeric. Assignment cost (same values as `basic_costs`).

preference Numeric 0-10. Nurse preference score.

skill_match Integer 0/1. Binary indicator: 1 if nurse skills match shift requirements, 0 otherwise.

nurses A tibble with 10 rows describing nurse characteristics:

nurse_id Integer 1-10. Links to `schedule_df` and `basic_costs` rows.

experience_years Numeric 1-20. Years of nursing experience.

department Character. Primary department: "ICU", "ER", "General", or "Pediatrics".

shift_preference Character. Preferred shift type: "day", "evening", or "night".

certification_level Integer 1-3. Certification level where 3 is highest (e.g., 1=RN, 2=BSN, 3=MSN).

shifts A tibble with 10 rows describing shift requirements:

shift_id Integer 1-10. Links to `schedule_df` and `basic_costs` cols.

department Character. Department needing coverage.

shift_type Character. Shift type: "day", "evening", or "night".

min_experience Numeric. Minimum years of experience required.

- min_certification** Integer 1-3. Minimum certification level.
- weekly_df** A tibble for batch solving with 500 rows (5 days x 10 nurses x 10 shifts):
- day** Character. Day of week: "Mon", "Tue", "Wed", "Thu", "Fri".
 - nurse_id** Integer 1-10. Nurse identifier.
 - shift_id** Integer 1-10. Shift identifier.
 - cost** Numeric. Daily assignment cost (varies by day).
 - preference** Numeric 0-10. Daily preference score.
- Use with `group_by(day)` for solving each day's schedule.
- nurses_extended** A tibble with 200 nurses for matching examples, representing a treatment group (e.g., full-time nurses):
- nurse_id** Integer 1-200. Unique identifier.
 - age** Numeric 22-65. Nurse age in years.
 - experience_years** Numeric 0-40. Years of nursing experience.
 - hourly_rate** Numeric 25-75. Hourly wage in dollars.
 - department** Character. Primary department assignment.
 - certification_level** Integer 1-3. Certification level.
 - is_fulltime** Logical. TRUE for full-time status.
- controls_extended** A tibble with 300 potential control nurses (e.g., part-time or registry nurses) for matching. Same structure as `nurses_extended`. Designed to have systematic differences from `nurses_extended` (older, less experience on average) to demonstrate matching's ability to create comparable groups.

Details

This dataset is used throughout the `couplr` documentation to provide a consistent, realistic example that evolves in complexity. It supports three use cases: (1) basic LAP solving with cost matrices, (2) batch solving across multiple days, and (3) matching workflows comparing nurse groups.

The dataset is designed to demonstrate progressively complex scenarios:

Basic LAP (`vignette("getting-started")`):

- `basic_costs`: Simple 10x10 assignment
- `preferences`: Maximization problem
- `schedule_df`: Data frame input, grouped workflows
- `weekly_df`: Batch solving across days

Algorithm comparison (`vignette("algorithms")`):

- Use `basic_costs` to compare algorithm behavior
- Modify with NA values for sparse scenarios

Matching workflows (`vignette("matching-workflows")`):

- `nurses_extended`: Treatment group (full-time nurses)
- `controls_extended`: Control pool (part-time/registry nurses)
- Match on age, experience, department for causal analysis

See Also

[lap_solve](#) for basic assignment solving, [lap_solve_batch](#) for batch solving, [match_couples](#) for matching workflows, [vignette\("getting-started"\)](#) for introductory tutorial

Examples

```
# Basic assignment: assign nurses to shifts minimizing cost
lap_solve(hospital_staff$basic_costs)

# Maximize preferences instead
lap_solve(hospital_staff$preferences, maximize = TRUE)

# Data frame workflow
library(dplyr)
hospital_staff$schedule_df |>
  lap_solve(nurse_id, shift_id, cost)

# Batch solve weekly schedule
hospital_staff$weekly_df |>
  group_by(day) |>
  lap_solve(nurse_id, shift_id, cost)

# Matching workflow: match full-time to part-time nurses
match_couples(
  left = hospital_staff$nurses_extended,
  right = hospital_staff$controls_extended,
  vars = c("age", "experience_years", "certification_level"),
  auto_scale = TRUE
)
```

is_distance_object *Check if Object is a Distance Object*

Description

Check if Object is a Distance Object

Usage

```
is_distance_object(x)
```

Arguments

x Object to check

Value

Logical: TRUE if x is a distance_object

Examples

```
left <- data.frame(id = 1:3, x = c(1, 2, 3))
right <- data.frame(id = 4:6, x = c(1.1, 2.1, 3.1))
dist_obj <- compute_distances(left, right, vars = "x")
is_distance_object(dist_obj) # TRUE
is_distance_object(list()) # FALSE
```

is_lap_solve_batch_result

Check if object is a batch assignment result

Description

Check if object is a batch assignment result

Usage

```
is_lap_solve_batch_result(x)
```

Arguments

x Object to test

Value

Logical indicating if x is a batch assignment result

is_lap_solve_kbest_result

Check if object is a k-best assignment result

Description

Check if object is a k-best assignment result

Usage

```
is_lap_solve_kbest_result(x)
```

Arguments

x Object to test

Value

Logical indicating if x is a k-best assignment result

is_lap_solve_result *Check if object is an assignment result*

Description

Check if object is an assignment result

Usage

```
is_lap_solve_result(x)
```

Arguments

x Object to test

Value

Logical indicating if x is an assignment result

join_matched *Join Matched Pairs with Original Data*

Description

Creates an analysis-ready dataset by joining matched pairs with variables from the original left and right datasets. This eliminates the need for manual joins and provides a convenient format for downstream analysis.

Usage

```
join_matched(result, ...)

## S3 method for class 'matching_result'
join_matched(
  result,
  left,
  right,
  left_vars = NULL,
  right_vars = NULL,
  left_id = "id",
  right_id = "id",
  suffix = c("_left", "_right"),
  include_distance = TRUE,
  include_pair_id = TRUE,
  include_block_id = TRUE,
  ...
)
```

```

)

## S3 method for class 'full_matching_result'
join_matched(result, left, right, left_id = "id", right_id = "id", ...)

## S3 method for class 'cem_result'
join_matched(result, left, right, left_id = "id", right_id = "id", ...)

## S3 method for class 'subclass_result'
join_matched(result, data = NULL, ...)

```

Arguments

result	A result object from <code>match_couples()</code> , <code>greedy_couples()</code> , <code>full_match()</code> , or <code>cem_match()</code>
...	Additional arguments passed to methods
left	The original left dataset
right	The original right dataset
left_vars	Character vector of variable names to include from left. If NULL (default), includes all variables except the ID column.
right_vars	Character vector of variable names to include from right. If NULL (default), includes all variables except the ID column.
left_id	Name of the ID column in left dataset (default: "id")
right_id	Name of the ID column in right dataset (default: "id")
suffix	Character vector of length 2 specifying suffixes for left and right variables (default: <code>c("_left", "_right")</code>)
include_distance	Include the matching distance in output (default: TRUE)
include_pair_id	Include <code>pair_id</code> column (default: TRUE)
include_block_id	Include <code>block_id</code> if blocking was used (default: TRUE)
data	Data frame used for subclassification

Details

This function simplifies the common workflow of joining matched pairs with original data. Instead of manually merging `result$pairs` with left and right datasets, `join_matched()` handles the joins automatically and applies consistent naming conventions.

When variables appear in both left and right datasets, suffixes are appended to distinguish them (e.g., "age_left" and "age_right"). This makes it easy to compute differences or use both values in models.

Value

A tibble with one row per matched pair, containing:

- `pair_id`: Sequential pair identifier (if `include_pair_id = TRUE`)
- `left_id`: ID from left dataset
- `right_id`: ID from right dataset
- `distance`: Matching distance (if `include_distance = TRUE`)
- `block_id`: Block identifier (if blocking used and `include_block_id = TRUE`)
- Variables from left dataset (with left suffix)
- Variables from right dataset (with right suffix)

Examples

```
# Basic usage
left <- data.frame(
  id = 1:5,
  treatment = 1,
  age = c(25, 30, 35, 40, 45),
  income = c(45000, 52000, 48000, 61000, 55000)
)

right <- data.frame(
  id = 6:10,
  treatment = 0,
  age = c(24, 29, 36, 41, 44),
  income = c(46000, 51500, 47500, 60000, 54000)
)

result <- match_couples(left, right, vars = c("age", "income"))
matched_data <- join_matched(result, left, right)
head(matched_data)

# Specify which variables to include
matched_data <- join_matched(
  result, left, right,
  left_vars = c("treatment", "age", "income"),
  right_vars = c("age", "income"),
  suffix = c("_treated", "_control")
)

# Without distance or pair_id
matched_data <- join_matched(
  result, left, right,
  include_distance = FALSE,
  include_pair_id = FALSE
)
```

lap_solve	<i>Solve linear assignment problems</i>
-----------	---

Description

Provides a tidy interface for solving the linear assignment problem using 'Hungarian' or 'Jonker-Volgenant' algorithms. Supports rectangular matrices, NA/Inf masking, and data frame inputs.

Usage

```
lap_solve(
  x,
  source = NULL,
  target = NULL,
  cost = NULL,
  maximize = FALSE,
  method = "auto",
  forbidden = NA
)
```

Arguments

<code>x</code>	Cost matrix, data frame, or tibble. If a data frame/tibble, must include columns specified by <code>source</code> , <code>target</code> , and <code>cost</code> .
<code>source</code>	Column name for source/row indices (if <code>x</code> is a data frame)
<code>target</code>	Column name for target/column indices (if <code>x</code> is a data frame)
<code>cost</code>	Column name for costs (if <code>x</code> is a data frame)
<code>maximize</code>	Logical; if TRUE, maximizes total cost instead of minimizing (default: FALSE)
<code>method</code>	Algorithm to use. One of: <ul style="list-style-type: none"> • "auto" (default): automatically selects best algorithm • "jv": 'Jonker-Volgenant' algorithm (general purpose, fast) • "hungarian": Classic 'Hungarian' algorithm • "auction": 'Bertsekas' auction algorithm (good for large dense problems) • "sap": Sparse assignment (good for sparse/rectangular problems) • "hk01": 'Hopcroft-Karp' for binary/uniform costs
<code>forbidden</code>	Value to mark forbidden assignments (default: NA). Can also use Inf.

Value

A tibble with columns:

- `source`: row/source indices
- `target`: column/target indices
- `cost`: cost of each assignment
- `total_cost`: total cost (attribute)

Examples

```
# Matrix input
cost <- matrix(c(4, 2, 5, 3, 3, 6, 7, 5, 4), nrow = 3)
lap_solve(cost)

# Data frame input
library(dplyr)
df <- tibble(
  source = rep(1:3, each = 3),
  target = rep(1:3, times = 3),
  cost = c(4, 2, 5, 3, 3, 6, 7, 5, 4)
)
lap_solve(df, source, target, cost)

# With NA masking (forbidden assignments)
cost[1, 3] <- NA
lap_solve(cost)

# Grouped data frames
df <- tibble(
  sim = rep(1:2, each = 9),
  source = rep(1:3, times = 6),
  target = rep(1:3, each = 3, times = 2),
  cost = runif(18, 1, 10)
)
df |> group_by(sim) |> lap_solve(source, target, cost)
```

lap_solve_batch

Solve multiple assignment problems efficiently

Description

Solve many independent assignment problems at once. Supports lists of matrices, 3D arrays, or grouped data frames. Optional parallel execution via `n_threads`.

Usage

```
lap_solve_batch(
  x,
  source = NULL,
  target = NULL,
  cost = NULL,
  maximize = FALSE,
  method = "auto",
  n_threads = 1,
  forbidden = NA
)
```

Arguments

x	One of: List of cost matrices, 3D array, or grouped data frame
source	Column name for source indices (if x is a grouped data frame)
target	Column name for target indices (if x is a grouped data frame)
cost	Column name for costs (if x is a grouped data frame)
maximize	Logical; if TRUE, maximizes total cost (default: FALSE)
method	Algorithm to use (default: "auto"). See lap_solve for options.
n_threads	Number of threads for parallel execution (default: 1). Set to NULL to use all available cores.
forbidden	Value to mark forbidden assignments (default: NA)

Value

A tibble with columns:

- problem_id: identifier for each problem
- source: source indices for assignments
- target: target indices for assignments
- cost: cost of each assignment
- total_cost: total cost for each problem
- method_used: algorithm used for each problem

Examples

```
# List of matrices
costs <- list(
  matrix(c(1, 2, 3, 4), 2, 2),
  matrix(c(5, 6, 7, 8), 2, 2)
)
lap_solve_batch(costs)

# 3D array
arr <- array(runif(2 * 2 * 10), dim = c(2, 2, 10))
lap_solve_batch(arr)

# Grouped data frame
library(dplyr)
df <- tibble(
  sim = rep(1:5, each = 9),
  source = rep(1:3, times = 15),
  target = rep(1:3, each = 3, times = 5),
  cost = runif(45, 1, 10)
)
df |> group_by(sim) |> lap_solve_batch(source, target, cost)

# Parallel execution (requires n_threads > 1)
lap_solve_batch(costs, n_threads = 2)
```

lap_solve_kbest	<i>Find k-best optimal assignments</i>
-----------------	--

Description

Returns the top k optimal (or near-optimal) assignments using 'Murty' algorithm. Useful for exploring alternative optimal solutions or finding robust assignments.

Usage

```
lap_solve_kbest(
  x,
  k = 3,
  source = NULL,
  target = NULL,
  cost = NULL,
  maximize = FALSE,
  method = "murty",
  single_method = "jv",
  forbidden = NA
)
```

Arguments

x	Cost matrix, data frame, or tibble. If a data frame/tibble, must include columns specified by source, target, and cost.
k	Number of best solutions to return (default: 3)
source	Column name for source/row indices (if x is a data frame)
target	Column name for target/column indices (if x is a data frame)
cost	Column name for costs (if x is a data frame)
maximize	Logical; if TRUE, finds k-best maximizing assignments (default: FALSE)
method	Algorithm for each sub-problem (default: "murty"). Future versions may support additional methods.
single_method	Algorithm used for solving each node in the search tree (default: "jv")
forbidden	Value to mark forbidden assignments (default: NA)

Value

A tibble with columns:

- rank: ranking of solutions (1 = best, 2 = second best, etc.)
- solution_id: unique identifier for each solution
- source: source indices
- target: target indices
- cost: cost of each edge in the assignment
- total_cost: total cost of the complete solution

Examples

```
# Matrix input - find 5 best solutions
cost <- matrix(c(4, 2, 5, 3, 3, 6, 7, 5, 4), nrow = 3)
lap_solve_kbest(cost, k = 5)

# Data frame input
library(dplyr)
df <- tibble(
  source = rep(1:3, each = 3),
  target = rep(1:3, times = 3),
  cost = c(4, 2, 5, 3, 3, 6, 7, 5, 4)
)
lap_solve_kbest(df, k = 3, source, target, cost)

# With maximization
lap_solve_kbest(cost, k = 3, maximize = TRUE)
```

lap_solve_line_metric *Solve 1-D Line Assignment Problem*

Description

Solves the linear assignment problem when both sources and targets are ordered points on a line. Uses efficient $O(n*m)$ dynamic programming for rectangular problems and $O(n)$ sorting for square problems.

Usage

```
lap_solve_line_metric(x, y, cost = "L1", maximize = FALSE)
```

Arguments

x	Numeric vector of source positions (will be sorted internally)
y	Numeric vector of target positions (will be sorted internally)
cost	Cost function for distance. Either: <ul style="list-style-type: none"> • "L1" (default): absolute distance ('Manhattan' distance) • "L2": squared distance (squared 'Euclidean' distance) Can also use aliases: "abs", "manhattan" for L1; "sq", "squared", "quadratic" for L2
maximize	Logical; if TRUE, maximizes total cost instead of minimizing (default: FALSE)

Details

This is a specialized solver that exploits the structure of 1-dimensional assignment problems where costs depend only on the distance between points on a line. It is much faster than general LAP solvers for this special case.

The algorithm works as follows:

Square case ($n == m$): Both vectors are sorted and matched in order: $x[1] \rightarrow y[1]$, $x[2] \rightarrow y[2]$, etc. This is optimal for any metric cost function on a line.

Rectangular case ($n < m$): Uses dynamic programming to find the optimal assignment that matches all n sources to a subset of the m targets, minimizing total distance. The DP recurrence is:

$$dp[i][j] = \min(dp[i][j-1], dp[i-1][j-1] + \text{cost}(x[i], y[j]))$$

This finds the minimum cost to match the first i sources to the first j targets.

Complexity:

- Time: $O(n*m)$ for rectangular, $O(n \log n)$ for square
- Space: $O(n*m)$ for DP table

Value

A list with components:

- `match`: Integer vector of length n with 1-based column indices
- `total_cost`: Total cost of the assignment

Examples

```
# Square case: equal number of sources and targets
x <- c(1.5, 3.2, 5.1)
y <- c(2.0, 3.0, 5.5)
result <- lap_solve_line_metric(x, y, cost = "L1")
print(result)

# Rectangular case: more targets than sources
x <- c(1.0, 3.0, 5.0)
y <- c(0.5, 2.0, 3.5, 4.5, 6.0)
result <- lap_solve_line_metric(x, y, cost = "L2")
print(result)

# With unsorted inputs (will be sorted internally)
x <- c(5.0, 1.0, 3.0)
y <- c(4.5, 0.5, 6.0, 2.0, 3.5)
result <- lap_solve_line_metric(x, y, cost = "L1")
print(result)
```

matchmaker

Create blocks for stratified matching

Description

Constructs blocks (strata) for matching, using either grouping variables or clustering algorithms. Returns the input data frames with block IDs assigned, along with block summary statistics.

Usage

```

matchmaker(
  left,
  right,
  block_type = c("none", "group", "cluster"),
  block_by = NULL,
  block_vars = NULL,
  block_method = "kmeans",
  n_blocks = NULL,
  min_left = 1,
  min_right = 1,
  drop_imbalanced = FALSE,
  imbalance_threshold = Inf,
  return_dropped = TRUE,
  ...
)

```

Arguments

<code>left</code>	Data frame of "left" units (e.g., treated, cases)
<code>right</code>	Data frame of "right" units (e.g., control, controls)
<code>block_type</code>	Type of blocking to use: <ul style="list-style-type: none"> • "none": No blocking (default) • "group": Block by existing categorical variable(s) • "cluster": Block using clustering algorithm
<code>block_by</code>	Variable name(s) for grouping (if <code>block_type = "group"</code>)
<code>block_vars</code>	Variable names for clustering (if <code>block_type = "cluster"</code>)
<code>block_method</code>	Clustering method (if <code>block_type = "cluster"</code>): <ul style="list-style-type: none"> • "kmeans": K-means clustering • "hclust": Hierarchical clustering
<code>n_blocks</code>	Target number of blocks (for clustering)
<code>min_left</code>	Minimum number of left units per block
<code>min_right</code>	Minimum number of right units per block
<code>drop_imbalanced</code>	Drop blocks with extreme imbalance
<code>imbalance_threshold</code>	Maximum allowed $ n_left - n_right / \max(n_left, n_right)$
<code>return_dropped</code>	Include dropped blocks in output
<code>...</code>	Additional arguments passed to clustering function

Details

This function does NOT perform matching - it only creates the block structure. Use [match_couples\(\)](#) or [greedy_couples\(\)](#) to perform matching within blocks.

Value

A list with class "matchmaker_result" containing:

- left: Left data frame with block_id column added
- right: Right data frame with block_id column added
- block_summary: Summary statistics for each block
- dropped: Information about dropped blocks (if any)
- info: Metadata about blocking process

Examples

```
# Group blocking
left <- data.frame(id = 1:10, region = rep(c("A", "B"), each = 5), x = rnorm(10))
right <- data.frame(id = 11:20, region = rep(c("A", "B"), each = 5), x = rnorm(10))
blocks <- matchmaker(left, right, block_type = "group", block_by = "region")
print(blocks$block_summary)

# Clustering
blocks <- matchmaker(left, right, block_type = "cluster",
                     block_vars = "x", n_blocks = 3)
```

match_couples

Optimal matching using linear assignment

Description

Performs optimal one-to-one matching between two datasets using linear assignment problem (LAP) solvers. Supports blocking, distance constraints, and various distance metrics.

Usage

```
match_couples(
  left,
  right = NULL,
  vars = NULL,
  distance = "euclidean",
  weights = NULL,
  scale = FALSE,
  auto_scale = FALSE,
  max_distance = Inf,
  calipers = NULL,
  block_id = NULL,
  ignore_blocks = FALSE,
  require_full_matching = FALSE,
  method = "auto",
```

```

    return_unmatched = TRUE,
    return_diagnostics = FALSE,
    parallel = FALSE,
    replace = FALSE,
    ratio = 1L,
    check_costs = TRUE,
    sigma = NULL
  )

```

Arguments

left	Data frame of "left" units (e.g., treated, cases)
right	Data frame of "right" units (e.g., control, controls)
vars	Variable names to use for distance computation
distance	Distance metric: "euclidean", "manhattan", "mahalanobis", or a custom function
weights	Optional named vector of variable weights
scale	Scaling method: FALSE (none), "standardize", "range", or "robust"
auto_scale	If TRUE, automatically check variable health and select scaling method (default: FALSE)
max_distance	Maximum allowed distance (pairs exceeding this are forbidden)
calipers	Named list of per-variable maximum absolute differences
block_id	Column name containing block IDs (for stratified matching)
ignore_blocks	If TRUE, ignore block_id even if present
require_full_matching	If TRUE, error if any units remain unmatched
method	LAP solver: "auto", "hungarian", "jv", "gabow_tarjan", etc.
return_unmatched	Include unmatched units in output
return_diagnostics	Include detailed diagnostics in output
parallel	Enable parallel processing for blocked matching. Requires 'future' and 'future.apply' packages. Can be: <ul style="list-style-type: none"> • FALSE: Sequential processing (default) • TRUE: Auto-configure parallel backend • Character: Specify future plan (e.g., "multisession", "multicore")
replace	If TRUE, allow matching with replacement (same right unit can be matched to multiple left units). Default: FALSE.
ratio	Integer, number of right units to match per left unit. Default: 1 (one-to-one matching). For k:1 matching, set ratio = k.
check_costs	If TRUE, check distance distribution for potential problems and provide helpful warnings before matching (default: TRUE)
sigma	Optional covariance matrix for Mahalanobis distance. If NULL (default), the pooled sample covariance is used. Only relevant when distance = "mahalanobis".

Details

This function finds the matching that minimizes total distance among all feasible matchings, subject to constraints. Use `greedy_couples()` for faster approximate matching on large datasets.

Value

A list with class "matching_result" containing:

- `pairs`: Tibble of matched pairs with distances
- `unmatched`: List of unmatched left and right IDs
- `info`: Matching diagnostics and metadata

Examples

```
# Basic matching
left <- data.frame(id = 1:5, x = c(1, 2, 3, 4, 5), y = c(2, 4, 6, 8, 10))
right <- data.frame(id = 6:10, x = c(1.1, 2.2, 3.1, 4.2, 5.1), y = c(2.1, 4.1, 6.2, 8.1, 10.1))
result <- match_couples(left, right, vars = c("x", "y"))
print(result$pairs)

# With constraints
result <- match_couples(left, right, vars = c("x", "y"),
                        max_distance = 1,
                        calipers = list(x = 0.5))

# With blocking
left$region <- c("A", "A", "B", "B", "B")
right$region <- c("A", "A", "B", "B", "B")
blocks <- matchmaker(left, right, block_type = "group", block_by = "region")
result <- match_couples(blocks$left, blocks$right, vars = c("x", "y"))
```

match_data

Extract Analysis-Ready Data from Matching Results

Description

A generic function that converts any couplr matching result into a single analysis-ready data frame with weights, subclass, and distance columns. This is the couplr equivalent of MatchIt's `match.data()`.

Usage

```
match_data(result, ...)
```

```
## S3 method for class 'matching_result'
```

```
match_data(result, left, right, left_id = "id", right_id = "id", ...)
```

```
## S3 method for class 'full_matching_result'
match_data(result, left, right, left_id = "id", right_id = "id", ...)

## S3 method for class 'cem_result'
match_data(result, left, right, left_id = "id", right_id = "id", ...)

## S3 method for class 'subclass_result'
match_data(result, data = NULL, ...)
```

Arguments

result	A couplr result object (matching_result, full_matching_result, cem_result, or subclass_result)
...	Additional arguments passed to methods
left	Data frame of left (treated) units
right	Data frame of right (control) units
left_id	Name of ID column in left (default: "id")
right_id	Name of ID column in right (default: "id")
data	Data frame containing all units (for CEM and subclassification, left and right are not always needed separately)

Details

The output format is compatible with downstream packages like **cobalt**, **WeightIt**, and **marginal-effects**. The stacked (long) format with treatment and weights columns is the standard layout expected by these tools.

Value

A tibble with all original variables plus standardized columns:

id Unit identifier
treatment 1 for left/treated, 0 for right/control
weights Matching weights
subclass Matched group/stratum identifier
distance Matching distance (where applicable)

Examples

```
set.seed(42)
left <- data.frame(id = 1:5, age = c(25, 35, 45, 55, 65))
right <- data.frame(id = 6:15, age = runif(10, 20, 70))
result <- match_couples(left, right, vars = "age")
md <- match_data(result, left, right)
head(md)
```

 pixel_morph

Pixel-level image morphing (final frame only)

Description

Computes optimal pixel assignment from A to B and returns the final transported frame (without intermediate animation frames).

Usage

```
pixel_morph(
  imgA,
  imgB,
  n_frames = 16L,
  mode = c("color_walk", "exact", "recursive"),
  lap_method = "jv",
  maximize = FALSE,
  quantize_bits = 5L,
  downscale_steps = 0L,
  alpha = 1,
  beta = 0,
  patch_size = 1L,
  upscale = 1,
  show = interactive()
)
```

Arguments

imgA	Source image (file path or magick image object)
imgB	Target image (file path or magick image object)
n_frames	Internal parameter for rendering (default: 16)
mode	Assignment algorithm: "color_walk" (default), "exact", or "recursive"
lap_method	LAP solver method (default: "jv")
maximize	Logical, maximize instead of minimize cost (default: FALSE)
quantize_bits	Color quantization for "color_walk" mode (default: 5)
downscale_steps	Number of 2x reductions before computing assignment (default: 0)
alpha	Weight for color distance in cost function (default: 1)
beta	Weight for spatial distance in cost function (default: 0)
patch_size	Tile size for tiled modes (default: 1)
upscale	Post-rendering upscaling factor (default: 1)
show	Logical, display result in viewer (default: interactive())

Details

Transport-Only Semantics:

This function returns a SHARP, pixel-perfect transport of A's pixels to positions determined by the assignment to B.

Key Points:

- Assignment computed using: $\text{cost} = \alpha * \text{color_dist} + \beta * \text{spatial_dist}$
- B's COLORS influence assignment but DO NOT appear in output
- Result has A's colors arranged to match B's layout
- No motion blur (unlike intermediate frames in animation)

See [pixel_morph_animate](#) for detailed explanation of assignment vs rendering semantics.

Permutation Warnings:

Assignment is guaranteed to be a bijection (permutation) ONLY when:

- `downscale_steps = 0` (no resolution changes)
- `mode = "exact"` with `patch_size = 1`

With downscaling or tiled modes, assignment may have:

- **Overlaps:** Multiple source pixels map to same destination (last write wins)
- **Holes:** Some destinations never filled (remain transparent)

If assignment is not a bijection (due to downscaling or tiling), a warning will be issued. The result may contain:

- Overlapped pixels (multiple sources -> one destination)
- Transparent holes (some destinations unfilled)

For guaranteed pixel-perfect results, use:

```
pixel_morph(A, B, mode = "exact", downscale_steps = 0)
```

Value

magick image object of the final transported frame

See Also

[pixel_morph_animate](#) for animated version

Examples

```
if (requireNamespace("magick", quietly = TRUE)) {
  imgA <- system.file("extdata/icons/circleA_40.png", package = "couplr")
  imgB <- system.file("extdata/icons/circleB_40.png", package = "couplr")
  if (nzchar(imgA) && nzchar(imgB)) {
    result <- pixel_morph(imgA, imgB, n_frames = 4, show = FALSE)
  }
}
```

pixel_morph_animate *Pixel-level image morphing (animation)*

Description

Creates an animated morph by computing optimal pixel assignment from image A to image B, then rendering intermediate frames showing the transport.

Usage

```
pixel_morph_animate(
  imgA,
  imgB,
  n_frames = 16L,
  fps = 10L,
  format = c("gif", "webp", "mp4"),
  outfile = NULL,
  show = interactive(),
  mode = c("color_walk", "exact", "recursive"),
  lap_method = "jv",
  maximize = FALSE,
  quantize_bits = 5L,
  downscale_steps = 0L,
  alpha = 1,
  beta = 0,
  patch_size = 1L,
  upscale = 1
)
```

Arguments

imgA	Source image (file path or magick image object)
imgB	Target image (file path or magick image object)
n_frames	Integer number of animation frames (default: 16)
fps	Frames per second for playback (default: 10)
format	Output format: "gif", "webp", or "mp4"
outfile	Optional output file path
show	Logical, display animation in viewer (default: interactive())
mode	Assignment algorithm: "color_walk" (default), "exact", or "recursive"
lap_method	LAP solver method (default: "jv")
maximize	Logical, maximize instead of minimize cost (default: FALSE)
quantize_bits	Color quantization for "color_walk" mode (default: 5)
downscale_steps	Number of 2x reductions before computing assignment (default: 0)

alpha	Weight for color distance in cost function (default: 1)
beta	Weight for spatial distance in cost function (default: 0)
patch_size	Tile size for tiled modes (default: 1)
upscale	Post-rendering upscaling factor (default: 1)

Details

Assignment vs Rendering Semantics:

CRITICAL: This function has two separate phases with different semantics:

Phase 1 - Assignment Computation:

The assignment is computed by minimizing:

$$\text{cost}(i, j) = \text{alpha} * \text{color_distance}(A[i], B[j]) + \text{beta} * \text{spatial_distance}(\text{pos}_i, \text{pos}_j)$$

This means B's COLORS influence which pixels from A map to which positions.

Phase 2 - Rendering (Transport-Only):

The renderer uses ONLY A's colors:

- Intermediate frames: A's pixels move along paths with motion blur
- Final frame: A's pixels at their assigned positions (sharp, no blur)
- B's colors NEVER appear in the output

Result: You get A's colors rearranged to match B's geometry/layout.

What This Means:

- B influences WHERE pixels go (via similarity in cost function)
- B does NOT determine WHAT COLORS appear in output
- Final image has A's palette arranged to mimic B's structure

Parameter Guidance:

For pure spatial rearrangement (ignore B's colors in assignment):

```
pixel_morph_animate(A, B, alpha = 0, beta = 1)
```

For color-similarity matching (default):

```
pixel_morph_animate(A, B, alpha = 1, beta = 0)
```

For hybrid (color + spatial):

```
pixel_morph_animate(A, B, alpha = 1, beta = 0.2)
```

Permutation Guarantees:

Assignment is guaranteed to be a bijection (permutation) ONLY when:

- `downscale_steps = 0` (no resolution changes)
- `mode = "exact"` with `patch_size = 1`

With downscaling or tiled modes, assignment may have:

- **Overlaps:** Multiple source pixels map to same destination (last write wins)
- **Holes:** Some destinations never filled (remain transparent)

A warning is issued if overlaps/holes are detected in the final frame.

Value

Invisibly returns a list with animation object and metadata:

animation	magick animation object
width	Image width in pixels
height	Image height in pixels
assignment	Integer vector of 1-based assignment indices (R convention)
n_pixels	Total number of pixels
mode	Mode used for matching
upscale	Upscaling factor applied

Examples

```
if (requireNamespace("magick", quietly = TRUE)) {
  imgA <- system.file("extdata/icons/circleA_40.png", package = "couplr")
  imgB <- system.file("extdata/icons/circleB_40.png", package = "couplr")
  if (nzchar(imgA) && nzchar(imgB)) {
    outfile <- tempfile(fileext = ".gif")
    pixel_morph_animate(imgA, imgB, outfile = outfile, n_frames = 4, show = FALSE)
  }
}
```

plot.balance_diagnostics

Plot method for balance diagnostics

Description

Produces a Love plot (dot plot) of standardized differences.

Usage

```
## S3 method for class 'balance_diagnostics'
plot(x, type = c("love", "histogram", "variance"), threshold = 0.1, ...)
```

Arguments

x	A balance_diagnostics object
type	Type of plot: "love" (default), "histogram", or "variance"
threshold	Threshold line for standardized differences (default: 0.1)
...	Additional arguments passed to plotting functions

Value

The balance_diagnostics object (invisibly)

plot.matching_result *Plot method for matching results*

Description

Produces a histogram of pairwise distances from a matching result.

Usage

```
## S3 method for class 'matching_result'  
plot(x, type = c("histogram", "density", "ecdf"), ...)
```

Arguments

x	A matching_result object
type	Type of plot: "histogram" (default), "density", or "ecdf"
...	Additional arguments passed to plotting functions

Value

The matching_result object (invisibly)

plot.sensitivity_analysis
Plot method for sensitivity analysis (base graphics)

Description

Plot method for sensitivity analysis (base graphics)

Usage

```
## S3 method for class 'sensitivity_analysis'  
plot(x, alpha = 0.05, ...)
```

Arguments

x	A sensitivity_analysis object
alpha	Significance level (default: 0.05)
...	Additional arguments passed to plot

Value

The sensitivity_analysis object (invisibly)

`preprocess_matching_vars`*Preprocess matching variables with automatic checks and scaling*

Description

Main preprocessing function that orchestrates variable health checks, categorical encoding, and automatic scaling selection.

Usage

```
preprocess_matching_vars(  
  left,  
  right,  
  vars,  
  auto_scale = TRUE,  
  scale_method = "auto",  
  check_health = TRUE,  
  remove_problematic = TRUE,  
  verbose = TRUE  
)
```

Arguments

<code>left</code>	Data frame of left units
<code>right</code>	Data frame of right units
<code>vars</code>	Character vector of variable names
<code>auto_scale</code>	Logical, whether to perform automatic preprocessing (default: TRUE)
<code>scale_method</code>	Scaling method: "auto", "standardize", "range", "robust", or FALSE
<code>check_health</code>	Logical, whether to check variable health (default: TRUE)
<code>remove_problematic</code>	Logical, automatically exclude constant/all-NA variables (default: TRUE)
<code>verbose</code>	Logical, whether to print warnings (default: TRUE)

Value

A list with class "preprocessing_result" containing:

- `left`: Preprocessed left data frame
- `right`: Preprocessed right data frame
- `vars`: Final variable names (after exclusions)
- `health`: Variable health diagnostics
- `scaling_method`: Selected scaling method
- `excluded_vars`: Variables that were excluded
- `warnings`: List of warnings issued

```
print.balance_diagnostics
```

Print Method for Balance Diagnostics

Description

Print Method for Balance Diagnostics

Usage

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

Arguments

x	A balance_diagnostics object
...	Additional arguments (ignored)

Value

Invisibly returns the input object x.

```
print.cem_result
```

Print Method for CEM Results

Description

Print Method for CEM Results

Usage

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

Arguments

x	A cem_result object
...	Additional arguments (ignored)

Value

Invisibly returns the input object x.

`print.distance_object` *Print Method for Distance Objects*

Description

Print Method for Distance Objects

Usage

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

Arguments

<code>x</code>	A <code>distance_object</code>
<code>...</code>	Additional arguments (ignored)

Value

Invisibly returns the input object `x`.

`print.full_matching_result`
Print Method for Full Matching Results

Description

Print Method for Full Matching Results

Usage

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

Arguments

<code>x</code>	A <code>full_matching_result</code> object
<code>...</code>	Additional arguments (ignored)

Value

Invisibly returns the input object `x`.

```
print.lap_solve_batch_result
```

Print method for batch assignment results

Description

Prints a summary and the table of results for a batch of assignment problems solved with `lap_solve_batch()`.

Usage

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

Arguments

`x` A `lap_solve_batch_result` object.
`...` Additional arguments passed to `print()`. Currently ignored.

Value

Invisibly returns the input object `x`.

```
print.lap_solve_kbest_result
```

Print method for k-best assignment results

Description

Print method for k-best assignment results

Usage

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

Arguments

`x` A `lap_solve_kbest_result`.
`...` Additional arguments passed to `print()`. Ignored.

Value

Invisibly returns the input object `x`.

```
print.lap_solve_result
```

Print method for assignment results

Description

Nicely prints a lap_solve_result object, including the assignments, total cost, and method used.

Usage

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

Arguments

x	A lap_solve_result object.
...	Additional arguments passed to print(). Currently ignored.

Value

Invisibly returns the input object x.

```
print.matching_result
```

Print method for matching results

Description

Print method for matching results

Usage

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

Arguments

x	A matching_result object
...	Additional arguments (ignored)

Value

Invisibly returns the input object x.

```
print.matchmaker_result
```

Print method for matchmaker results

Description

Print method for matchmaker results

Usage

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

Arguments

x	A matchmaker_result object
...	Additional arguments (ignored)

Value

Invisibly returns the input object x.

```
print.preprocessing_result
```

Print method for preprocessing result

Description

Print method for preprocessing result

Usage

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

Arguments

x	A preprocessing_result object
...	Additional arguments (ignored)

Value

Invisibly returns the input object x.

```
print.sensitivity_analysis  
    Print method for sensitivity analysis
```

Description

Print method for sensitivity analysis

Usage

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

Arguments

x	A sensitivity_analysis object
...	Additional arguments (ignored)

Value

Invisibly returns the input object x.

```
print.subclass_result Print Method for Subclassification Results
```

Description

Print Method for Subclassification Results

Usage

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

Arguments

x	A subclass_result object
...	Additional arguments (ignored)

Value

Invisibly returns the input object x.

print.variable_health *Print method for variable health*

Description

Print method for variable health

Usage

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

Arguments

x	A variable_health object
...	Additional arguments (ignored)

Value

Invisibly returns the input object x.

ps_match *Propensity Score Matching*

Description

Matches treated and control units based on estimated propensity scores. Fits a logistic regression model (or accepts a pre-fitted one), computes logit propensity scores, and calls [match_couples\(\)](#) with a caliper on the logit scale.

Usage

```
ps_match(  
  formula = NULL,  
  data = NULL,  
  treatment = NULL,  
  ps_model = NULL,  
  caliper_sd = 0.2,  
  method = "auto",  
  replace = FALSE,  
  ratio = 1L,  
  ...  
)
```

Arguments

formula	Formula for propensity score model (treatment ~ covariates). Required if ps_model is NULL.
data	Combined dataset containing both treated and control units
treatment	Name of the binary treatment column (0/1 or logical)
ps_model	Pre-fitted glm object (alternative to formula). If provided, formula is ignored.
caliper_sd	Caliper width in standard deviations of logit(PS). Default: 0.2 (Rosenbaum and Rubin recommendation).
method	LAP solver method (default: "auto")
replace	If TRUE, match with replacement (default: FALSE)
ratio	Integer k for k:1 matching (default: 1)
...	Additional arguments passed to match_couples()

Details

The propensity score is the probability of treatment assignment conditional on observed covariates. Matching is performed on the logit of the propensity score (Rosenbaum and Rubin 1985), which provides better distributional properties than matching on the raw probability scale.

The default caliper of 0.2 SD of logit(PS) is recommended by Austin (2011) as removing approximately 98% of bias.

Value

A `matching_result` object with additional propensity score info in `result$info$ps_model` and `result$info$caliper_value`.

Examples

```
set.seed(42)
n <- 100
data <- data.frame(
  id = seq_len(n),
  treated = rbinom(n, 1, 0.4),
  age = rnorm(n, 50, 10),
  income = rnorm(n, 50000, 15000)
)
result <- ps_match(treated ~ age + income, data = data, treatment = "treated")
print(result)
```

Description

Assesses how sensitive a matched comparison is to hidden bias using Rosenbaum bounds on the Wilcoxon signed-rank statistic.

Usage

```
sensitivity_analysis(  
  result,  
  left,  
  right,  
  outcome_var,  
  gamma = seq(1, 3, by = 0.25),  
  alternative = c("greater", "less", "two.sided"),  
  left_id = "id",  
  right_id = "id"  
)
```

Arguments

<code>result</code>	A <code>matching_result</code> object from <code>match_couples()</code> or <code>greedy_couples()</code>
<code>left</code>	Original left (treated) dataset
<code>right</code>	Original right (control) dataset
<code>outcome_var</code>	Name of the outcome column in <code>left</code> and <code>right</code>
<code>gamma</code>	Numeric vector of sensitivity parameters (default: <code>seq(1, 3, by = 0.25)</code>). Gamma = 1 means no hidden bias.
<code>alternative</code>	Direction of the test: "greater" (default), "less", or "two.sided"
<code>left_id</code>	Name of ID column in left (default: "id")
<code>right_id</code>	Name of ID column in right (default: "id")

Details

Rosenbaum (2002, Chapter 4) bounds quantify how much hidden bias (an unobserved confounder) would be needed to explain away the observed treatment effect. The sensitivity parameter Gamma represents the maximum ratio of treatment odds between two matched units:

- Gamma = 1: No hidden bias (standard Wilcoxon test)
- Gamma = 2: One unit could be twice as likely to receive treatment due to an unobserved factor

The function computes upper and lower bounds on the p-value of the Wilcoxon signed-rank test under each level of hidden bias. A finding is "insensitive to bias" if `p_upper` remains below 0.05 even at large Gamma.

Value

An S3 object of class `sensitivity_analysis` containing:

results Tibble with columns: `gamma`, `t_stat`, `p_upper`, `p_lower`

n_pairs Number of matched pairs analyzed

critical_gamma Smallest `gamma` at which `p_upper > 0.05`

alternative Direction of test

References

Rosenbaum, P.R. (2002). *Observational Studies*, 2nd edition. Springer.

Examples

```
set.seed(42)
left <- data.frame(id = 1:20, x = rnorm(20), outcome = rnorm(20, 1, 1))
right <- data.frame(id = 21:40, x = rnorm(20), outcome = rnorm(20, 0, 1))
result <- match_couples(left, right, vars = "x")
sens <- sensitivity_analysis(result, left, right, outcome_var = "outcome")
print(sens)
```

sinkhorn

'Sinkhorn-Knopp' optimal transport solver

Description

Compute an entropy-regularized optimal transport plan using the 'Sinkhorn-Knopp' algorithm. Unlike other LAP solvers that return a hard 1-to-1 assignment, this returns a soft assignment (doubly stochastic matrix).

Usage

```
sinkhorn(
  cost,
  lambda = 10,
  tol = 1e-09,
  max_iter = 1000,
  r_weights = NULL,
  c_weights = NULL
)
```

Arguments

<code>cost</code>	Numeric matrix of transport costs. NA or Inf entries are treated as very high cost (effectively forbidden).
<code>lambda</code>	Regularization parameter (default 10). Higher values produce sharper (more deterministic) transport plans; lower values produce smoother distributions. Typical range: 1-100.
<code>tol</code>	Convergence tolerance (default 1e-9).
<code>max_iter</code>	Maximum iterations (default 1000).
<code>r_weights</code>	Optional numeric vector of row marginals (source distribution). Default is uniform. Will be normalized to sum to 1.
<code>c_weights</code>	Optional numeric vector of column marginals (target distribution). Default is uniform. Will be normalized to sum to 1.

Details

The 'Sinkhorn-Knopp' algorithm solves the entropy-regularized optimal transport problem:

$$P^* = \arg \min_P \langle C, P \rangle - \frac{1}{\lambda} H(P)$$

subject to row sums = `r_weights` and column sums = `c_weights`.

The entropy term $H(P)$ encourages spread in the transport plan. As $\lambda \rightarrow \text{Inf}$, the solution approaches the standard (unregularized) optimal transport.

Key differences from standard LAP solvers:

- Returns a soft assignment (probabilities) not a hard 1-to-1 matching
- Supports unequal marginals (weighted distributions)
- Differentiable, making it useful in ML pipelines
- Very fast: $O(n^2)$ per iteration with typically $O(1/\text{tol}^2)$ iterations

Use `sinkhorn_to_assignment()` to round the soft assignment to a hard matching.

Value

A list with elements:

- `transport_plan` — numeric matrix, the optimal transport plan P . Row sums approximate `r_weights`, column sums approximate `c_weights`.
- `cost` — the transport cost $\langle C, P \rangle$ (without entropy term).
- `u, v` — scaling vectors ($P = \text{diag}(u) * K * \text{diag}(v)$ where $K = \exp(-\lambda * C)$).
- `converged` — logical, whether the algorithm converged.
- `iterations` — number of iterations used.
- `lambda` — the regularization parameter used.

References

Cuturi, M. (2013). 'Sinkhorn Distances': Lightspeed Computation of Optimal Transport. *Advances in Neural Information Processing Systems*, 26.

See Also

[assignment\(\)](#) for hard 1-to-1 matching, [sinkhorn_to_assignment\(\)](#) to round soft assignments.

Examples

```
cost <- matrix(c(1, 2, 3, 4, 5, 6, 7, 8, 9), nrow = 3, byrow = TRUE)

# Soft assignment with default parameters
result <- sinkhorn(cost)
print(round(result$transport_plan, 3))

# Sharper assignment (higher lambda)
result_sharp <- sinkhorn(cost, lambda = 50)
print(round(result_sharp$transport_plan, 3))

# With custom marginals (more mass from row 1)
result_weighted <- sinkhorn(cost, r_weights = c(0.5, 0.25, 0.25))
print(round(result_weighted$transport_plan, 3))

# Round to hard assignment
hard_match <- sinkhorn_to_assignment(result)
print(hard_match)
```

sinkhorn_to_assignment

Round 'Sinkhorn' transport plan to hard assignment

Description

Convert a soft transport plan from [sinkhorn\(\)](#) to a hard 1-to-1 assignment using greedy rounding.

Usage

```
sinkhorn_to_assignment(result)
```

Arguments

result Either a result from [sinkhorn\(\)](#) or a transport plan matrix.

Details

Greedy rounding iteratively assigns each row to its most probable column, ensuring no column is assigned twice. This may not give the globally optimal hard assignment; for that, use the transport plan as a cost matrix with [assignment\(\)](#).

Value

Integer vector of column assignments (1-based), same format as `assignment()`.

See Also

`sinkhorn()`

Examples

```
cost <- matrix(c(1, 2, 3, 4, 5, 6, 7, 8, 9), nrow = 3, byrow = TRUE)
result <- sinkhorn(cost, lambda = 20)
hard_match <- sinkhorn_to_assignment(result)
print(hard_match)
```

 subclass_match

Subclassification on Propensity Score

Description

Divides units into K strata based on quantiles of the propensity score, then computes within-stratum weights for treatment effect estimation. This is a simple, transparent approach to propensity score adjustment that allows visual inspection of balance within each subclass.

Usage

```
subclass_match(
  formula = NULL,
  data = NULL,
  treatment = NULL,
  n_subclasses = 5L,
  ps = NULL,
  ps_model = NULL,
  estimand = "ATT"
)
```

Arguments

formula	Formula for propensity score model (e.g., <code>treatment ~ age + income</code>). Ignored if ps is provided.
data	Data frame containing all variables
treatment	Character, name of the binary treatment column (0/1)
n_subclasses	Integer, number of subclasses to create (default: 5). Cochran (1968) showed that 5 subclasses removes over 90% a single covariate.
ps	Optional pre-computed numeric vector of propensity scores (one per row in data). If NULL, a logistic regression model is fit using formula.
ps_model	Optional pre-fitted glm object for propensity scores
estimand	Target estimand: "ATT" (default), "ATE", or "ATC"

Details

The algorithm:

1. Estimate propensity scores via logistic regression (or use pre-computed scores)
2. Divide the propensity score distribution into K quantile-based strata
3. For each stratum, check overlap (both treated and control units present)
4. Compute within-stratum weights based on the target estimand:
 - **ATT**: Treated units get weight 1; control units get weight $n_{\text{treated_in_stratum}} / n_{\text{control_in_stratum}}$
 - **ATE**: Both groups get weight proportional to stratum size relative to total sample
 - **ATC**: Control units get weight 1; treated units get weight $n_{\text{control_in_stratum}} / n_{\text{treated_in_stratum}}$

Value

An S3 object of class c("subclass_result", "couplr_result") containing:

matched Tibble with columns id, side, subclass, ps, weight

subclass_summary Tibble with per-subclass statistics: counts, mean PS, and overlap status

info List with n_subclasses, estimand, n_left, n_right, method, vars

Examples

```
set.seed(42)
n <- 200
data <- data.frame(
  id = 1:n,
  age = rnorm(n, 40, 10),
  income = rnorm(n, 50000, 15000)
)
data$treatment <- rbinom(n, 1, plogis(-2 + 0.05 * data$age))
result <- subclass_match(treatment ~ age + income, data, treatment = "treatment")
print(result)
```

```
summary.balance_diagnostics
```

Summary method for balance diagnostics

Description

Summary method for balance diagnostics

Usage

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

Arguments

object A balance_diagnostics object
 ... Additional arguments (ignored)

Value

A list containing summary statistics (invisibly)

summary.distance_object

Summary Method for Distance Objects

Description

Summary Method for Distance Objects

Usage

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

Arguments

object A distance_object
 ... Additional arguments (ignored)

Value

Invisibly returns the input object.

summary.lap_solve_kbest_result

Get summary of k-best results

Description

Extract summary information from k-best assignment results.

Usage

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

Arguments

object An object of class lap_solve_kbest_result.
... Additional arguments (unused).

Value

A tibble with one row per solution containing:

- rank: solution rank
- solution_id: solution identifier
- total_cost: total cost of the solution
- n_assignments: number of assignments in the solution

summary.matching_result

Summary method for matching results

Description

Summary method for matching results

Usage

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

Arguments

object A matching_result object
... Additional arguments (ignored)

Value

A list containing summary statistics (invisibly)

```
summary.sensitivity_analysis
      Summary method for sensitivity analysis
```

Description

Summary method for sensitivity analysis

Usage

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

Arguments

object	A sensitivity_analysis object
...	Additional arguments (ignored)

Value

A list containing summary statistics (invisibly)

```
update_constraints      Update Constraints on Distance Object
```

Description

Apply new constraints to a precomputed distance object without recomputing the underlying distances. This is useful for exploring different constraint scenarios quickly.

Usage

```
update_constraints(dist_obj, max_distance = Inf, calipers = NULL)
```

Arguments

dist_obj	A distance_object from compute_distances()
max_distance	Maximum allowed distance (pairs with distance > max_distance become Inf)
calipers	Named list of per-variable calipers

Details

This function creates a new `distance_object` with modified constraints applied to the cost matrix. The original `distance_object` is not modified.

Constraints:

- `max_distance`: Sets cost to Inf for pairs exceeding this threshold
- `calipers`: Per-variable restrictions (e.g., `calipers = list(age = 5)`)

The function returns a new object rather than modifying in place, following R's copy-on-modify semantics.

Value

A new `distance_object` with updated `cost_matrix`

Examples

```
left <- data.frame(id = 1:5, age = c(25, 30, 35, 40, 45))
right <- data.frame(id = 6:10, age = c(24, 29, 36, 41, 44))
dist_obj <- compute_distances(left, right, vars = "age")

# Apply constraints
constrained <- update_constraints(dist_obj, max_distance = 2)
result <- match_couples(constrained)
```

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