

Package ‘seqcomp’

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Type Package

Title Sequential Comparison of Probabilistic Forecasts

Version 0.3.0

Description Implements tools for the anytime-valid sequential comparison of two or more probabilistic forecasters. Provides binary, categorical, and quantile scoring rules, together with finite-sample confidence sequences and e-processes following Choe and Ramdas (2024) <[doi:10.1287/opre.2021.0792](https://doi.org/10.1287/opre.2021.0792)>. Extends to multi-model evaluation via Sequential Model Confidence Sets, following Arnold, Gavriloopoulos, Schulz, and Ziegel (2026) <[doi:10.1093/jrsssb/qkag066](https://doi.org/10.1093/jrsssb/qkag066)>, using closure principles, joint confidence sequences, and accelerated closed-testing. Adaptive betting fractions for the strong null (aGRAPA and ONS-m) are adapted from Waudby-Smith and Ramdas (2024) <[doi:10.1093/jrsssb/qkad009](https://doi.org/10.1093/jrsssb/qkad009)>. Also includes Winkler-score comparisons, lag handling, and predictable-bound betting e-processes.

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URL <https://github.com/alasgarliakbar/seqcomp>,
<https://alasgarliakbar.github.io/seqcomp/>

BugReports <https://github.com/alasgarliakbar/seqcomp/issues>

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Description

seqcomp provides tools for comparing probabilistic forecasters sequentially, following the anytime-valid framework of Choe and Ramdas (2024). For three or more forecasters, the package additionally implements Sequential Model Confidence Sets following Arnold, Gavrilopoulos, Schulz and Ziegel (2026).

Details

The package is built around the score difference

$$\hat{\delta}_t = S(p_t, y_t) - S(q_t, y_t),$$

where scores are positively oriented, so larger values are better. Positive score differences favor forecaster p ; negative score differences favor forecaster q .

Main workflow

For most applications, start with `compare_forecasts()`. It computes pointwise scores, running mean score differences, confidence sequences, and e-processes in one call.

Scoring rules

The package includes positively oriented scoring rules such as `brier_score()`, `log_score()`, `spherical_score()`, `tick_loss()`, `qlike_score()`, `winkler_score()`, `crps_normal()`, `crps_empirical()`, and `crps_std()`.

Confidence sequences

Use `cs_hoeffding()` for Hoeffding-style confidence sequences, `cs_bernstein()` for empirical Bernstein confidence sequences, and `cs_asymptotic()` for asymptotic confidence sequences when finite-sample boundedness is not available.

E-processes

Use `eprocess()` for the main sub-exponential mixture e-process and `eprocess_rejections()` to extract first rejection times. For multi-step forecasts, see `eprocess_lag()`. For predictable time-varying bounds, see `eprocess_predictable()`.

Multiple forecasters

For three or more candidate forecasters, use `smcs_compare()` as the main entry point. It extends the pairwise workflow above to a Sequential Model Confidence Set (SMCS): the running set of models not yet shown to underperform some competitor. Lower-level access is available via `smcs_strong()` (strong or uniformly weak null, via closure testing) and `smcs_weak()` (time-varying weak null, via joint confidence sequences), with `vovk_wang_merge()` and `eprocess_betting()` as supporting building blocks. To construct optimal adaptive betting fractions for the strong null, use `build_agrapa_betting_array()` or `build_ons_betting_array()` for constant-bound scores, and `build_quantile_betting_arrays()` for conditionally bounded rules such as tick loss.

Winkler scores

For binary probability forecasts with unbounded base scores, use `winkler_score()`, `winkler_cs()`, `winkler_etest()`, or `winkler_compare()`.

References

- Arnold, S., Gavrilopoulos, G., Schulz, B. and Ziegel, J. (2026). Sequential model confidence sets. *Journal of the Royal Statistical Society Series B: Statistical Methodology*, qkag066.
- Choe, Y. J. and Ramdas, A. (2024). Comparing Sequential Forecasters. *Operations Research*, 72(4), 1368-1387.
- Howard, S. R., Ramdas, A., McAuliffe, J. and Sekhon, J. (2021). Time-uniform, nonparametric, nonasymptotic confidence sequences. *The Annals of Statistics*, 49(2), 1055-1080.
- Waudby-Smith, I. and Ramdas, A. (2024). Estimating means of bounded random variables by betting. *Journal of the Royal Statistical Society Series B: Statistical Methodology*, 86(1), 1-27.

brier_score	<i>Brier score for binary and categorical forecasts</i>
-------------	---

Description

Computes the positively oriented Brier/quadratic score. Vector probability input is treated as binary; matrix probability input is treated as categorical.

Usage

```
brier_score(p, y)
```

Arguments

- | | |
|---|--|
| p | Numeric vector in $[0, 1]$ for binary forecasts, or a numeric matrix whose rows are probability vectors for categorical forecasts. |
| y | For binary vector input, numeric vector in $\{0, 1\}$. For categorical matrix input, integer vector in $\{1, \dots, K\}$, where $K = \text{ncol}(p)$. |

Details

For binary forecasts, this computes

$$S(p, y) = -(p - y)^2.$$

For categorical forecasts, this computes

$$S(\mathbf{p}, y) = -\frac{1}{2} \|\mathbf{p} - e_y\|_2^2,$$

where e_y is the one-hot vector of the realised category.

With the convention that category 2 corresponds to the binary event $y = 1$, the categorical formula recovers the binary formula exactly when $K = 2$.

Value

Numeric vector of scores in $[-1, 0]$. Higher is better.

Bounds

Score differences lie in $[-1, 1]$, so use $c = 1$ for Theorem 1 and $c = 2$ for Theorems 2 and 3.

Examples

```
p <- c(0.2, 0.7, 0.9)
y <- c(0, 1, 1)
brier_score(p, y)
```

```
build_agrapa_betting_array
```

Build 3D Parameter Array for aGRAPA Betting

Description

Pre-computes the dynamic, pair-specific betting fractions required to evaluate the Sequential Model Confidence Set under the strong null hypothesis, using the aGRAPA algorithm.

Usage

```
build_agrapa_betting_array(scores, c_mat, period = 1, ...)
```

Arguments

scores	A $T \times m$ matrix of positively-oriented scores.
c_mat	A numeric scalar, an $m \times m$ matrix, or a $T \times m \times m$ array of predictable bounds (matching the <code>c_param</code> argument of <code>smcs_strong()</code>).
period	Integer. If greater than 1, uses the periodic betting variant of aGRAPA, which updates the betting fraction only every period time steps.
...	Additional arguments passed to <code>lambda_betting_agrapa()</code> (e.g., <code>kappa</code> , <code>prior_mean</code> , <code>prior_variance</code> , <code>fake_obs</code>).

Value

A $T \times m \times m$ numeric array of predictable betting fractions, suitable for passing to the `lambda_param` argument of `smcs_strong()`.

See Also

`lambda_betting_agrapa()`, `eprocess_betting()`, `smcs_strong()`

Examples

```
set.seed(42)
y <- rbinom(50, 1, 0.5)
forecasts <- matrix(runif(150), nrow = 50, ncol = 3)
scores <- matrix(0, nrow = 50, ncol = 3)
for (i in 1:3) scores[, i] <- brier_score(forecasts[, i], y)

# For Brier scores, differences are bounded in [-1, 1], so c = 2
lam_array <- build_agrapa_betting_array(scores, c_mat = 2)
dim(lam_array) # 50 x 3 x 3
```

`build_ons_betting_array`

Build 3D Parameter Array for ONS-m Betting

Description

Pre-computes the dynamic, pair-specific betting fractions required to evaluate the Sequential Model Confidence Set under the strong null hypothesis, using the Online Newton Step (ONS-m) algorithm.

Usage

```
build_ons_betting_array(scores, c_mat, period = 1, ...)
```

Arguments

<code>scores</code>	A $T \times m$ matrix of positively-oriented scores.
<code>c_mat</code>	A numeric scalar, an $m \times m$ matrix, or a $T \times m \times m$ array of predictable bounds (matching the <code>c_param</code> argument of <code>smcs_strong()</code>).
<code>period</code>	Integer. If greater than 1, uses the periodic betting variant of ONS-m, which updates the betting fraction only every period time steps.
<code>...</code>	Additional arguments passed to <code>lambda_betting_ons()</code> (e.g., <code>eta</code>).

Value

A $T \times m \times m$ numeric array of predictable betting fractions, suitable for passing to the `lambda_param` argument of `smcs_strong()`.

See Also

[lambda_betting_ons\(\)](#), [eprocess_betting\(\)](#), [smcs_strong\(\)](#)

Examples

```
set.seed(43)
y <- rbinom(50, 1, 0.5)
forecasts <- matrix(runif(150), nrow = 50, ncol = 3)
scores <- matrix(0, nrow = 50, ncol = 3)
for (i in 1:3) scores[, i] <- brier_score(forecasts[, i], y)

# For Brier scores, differences are bounded in [-1, 1], so c = 2
lam_array <- build_ons_betting_array(scores, c_mat = 2)
dim(lam_array) # 50 x 3 x 3
```

build_quantile_betting_arrays

Build 3D Parameter Arrays for Quantile Betting

Description

Pre-computes the dynamic, pair-specific bounding arrays (`c_array`) and betting fractions (`lambda_array`) required to evaluate the Sequential Model Confidence Set for quantile forecasts under the strong null hypothesis.

Usage

```
build_quantile_betting_arrays(forecasts, scores, tau, eps = 1e-08)
```

Arguments

<code>forecasts</code>	A $T \times m$ matrix of raw quantile forecasts.
<code>scores</code>	A $T \times m$ matrix of positively-oriented tick-loss scores.
<code>tau</code>	Numeric scalar in $(0, 1)$. The quantile level.
<code>eps</code>	Positive numeric scalar. Used both as the denominator safeguard in lambda_betting_quantile() and as the minimum predictable bound returned in <code>c_array</code> . Default: $1e-8$.

Value

A list containing two $T \times m \times m$ numeric arrays: `c_array` and `lambda_array`.

Note

Scale Translation: This function assumes forecasts and scores are evaluated on the raw, linear scale. To replicate the exact log-scale bounds used in the Arnold et al. (2026) Covid-19 case study, the forecast matrix and outcomes must be log-transformed prior to passing them to this pipeline.

Examples

```

set.seed(3)
fcsts <- matrix(runif(150), nrow = 50, ncol = 3)
y <- rbinom(50, 1, 0.5)
scores <- matrix(0, nrow = 50, ncol = 3)
for(i in 1:3) scores[, i] <- tick_loss(fcsts[, i], y, tau = 0.5)

arrays <- build_quantile_betting_arrays(fcsts, scores, tau = 0.5)
dim(arrays$c_array) # 50 x 3 x 3

```

calibrate_p_to_e	<i>P-to-e calibrator</i>
------------------	--------------------------

Description

Converts anytime-valid p-values to e-values using the mixture or simple calibrator, as used in CR24 Section 4.4.

Usage

```
calibrate_p_to_e(p, strategy = "mixture", eps = 1e-16)
```

Arguments

p	Numeric vector of p-values in (0, 1].
strategy	Character. "mixture" (default, from Vovk & Wang 2021) or "simple". See Details for the formulas.
eps	Numeric. Numerical guard for log(0). Default: 1e-16.

Details

Mixture calibrator (default, matches Python comparecast behaviour):

$$f(p) = \frac{1 - p + p \log(p)}{p (\log p)^2}$$

Simple calibrator (strategy = "simple"):

$$f(p) = \frac{1}{2\sqrt{p}}$$

Value

Numeric vector of e-values ≥ 0 .

Examples

```
p <- c(0.5, 0.1, 0.01)
calibrate_p_to_e(p)
calibrate_p_to_e(p, strategy = "simple")
```

cm_boundary	<i>Normal mixture (CM) boundary</i>
-------------	-------------------------------------

Description

Normal mixture (CM) boundary

Usage

```
cm_boundary(v, alpha, rho)
```

Arguments

v	Numeric vector ≥ 0 . Intrinsic time values (V_t or t).
alpha	Numeric in (0,1). ONE-SIDED significance level.
rho	Numeric > 0 . Tuning parameter. Obtain via <code>rho_from_vopt()</code> .

Value

Numeric vector of boundary values (cumulative-sum scale, before $/t$).

See Also

[rho_from_vopt\(\)](#) to compute rho from a target `v_opt`.

Examples

```
rho <- rho_from_vopt(v_opt = 10, alpha = 0.025)
u <- cm_boundary(v = 1:500, alpha = 0.025, rho = rho)
```

compare_forecasts	<i>Compare Two Sequential Forecasters</i>
-------------------	---

Description

Computes pointwise scores for two probabilistic forecasters and compares them sequentially using confidence sequences and, when valid finite-sample bounds are available, e-processes.

Usage

```
compare_forecasts(
  p,
  q,
  y,
  scoring_rule = c("brier", "spherical", "log"),
  alpha = 0.05,
  cs_type = NULL,
  compute_cs = TRUE,
  compute_e = TRUE,
  v_opt = 10,
  boundary = "mixture",
  lcb_only = FALSE,
  ucb_only = FALSE,
  eps = 1e-15,
  clip_max = 1e+07
)
```

Arguments

p	Forecasts from forecaster 1. For binary outcomes, a numeric vector of probabilities for event $y = 1$. For categorical outcomes, a numeric matrix whose rows are probability vectors.
q	Forecasts from forecaster 2, in the same format as p.
y	Outcomes. For binary vector forecasts, a numeric vector in $\{0, 1\}$. For categorical matrix forecasts, integer class labels in $\{1, \dots, K\}$.
scoring_rule	Character. Scoring rule used to compare forecasts. Currently supports "brier", "spherical", and "log".
alpha	Numeric in $(0, 1)$. Significance level. Default is 0.05.
cs_type	Character or NULL. Confidence sequence type: "bernstein", "hoeffding", "asymptotic", or "none". If NULL, the wrapper uses "bernstein" for bounded scoring rules ("brier" and "spherical") and "asymptotic" for "log".
compute_cs	Logical. If TRUE, compute a confidence sequence. Default is TRUE.
compute_e	Logical. If TRUE, compute two one-sided e-processes. Default is TRUE. This is only allowed for bounded score differences under the current wrapper, namely "brier" and "spherical".

v_opt	Numeric > 0. Intrinsic time at which the mixture boundary or e-process is tuned to be tightest. Default is 10.
boundary	Character. Boundary type passed to <code>cs_hoeffding()</code> or <code>cs_bernstein()</code> . Default is "mixture".
lcb_only	Logical. If TRUE, compute a lower one-sided empirical Bernstein CS. Only used when <code>cs_type = "bernstein"</code> .
ucb_only	Logical. If TRUE, compute an upper one-sided empirical Bernstein CS. Only used when <code>cs_type = "bernstein"</code> .
eps	Numeric. Probability floor passed to <code>log_score()</code> when <code>scoring_rule = "log"</code> . Default is 1e-15.
clip_max	Numeric. Maximum e-process value before clipping. Passed to <code>eprocess()</code> . Default is 1e7.

Details

This is a convenience wrapper around `brier_score()`, `spherical_score()`, `log_score()`, `cs_hoeffding()`, `cs_bernstein()`, `cs_asymptotic()`, and `eprocess()`. It is designed for the common workflow where the user has two forecast streams p and q , an outcome stream y , and wants a single tidy output object.

All scoring rules in `seqcomp` are positively oriented: higher scores are better. Therefore

$$\hat{\delta}_t = S(p_t, y_t) - S(q_t, y_t)$$

is positive when forecaster p performs better than forecaster q at time t .

For "brier" and "spherical", score differences are bounded in $[-1, 1]$. The wrapper therefore uses $c = 1$ for Hoeffding-style confidence sequences and $c = 2$ for empirical Bernstein confidence sequences and e-processes.

For "log", score differences are unbounded. The wrapper therefore defaults to `cs_asymptotic()` and refuses to compute finite-sample e-processes. For binary log-score comparisons where the Winkler construction is appropriate, use `winkler_compare()` instead.

Value

A data.frame with one row per time point and columns:

t Time index.

score_p Pointwise score of forecaster p .

score_q Pointwise score of forecaster q .

delta Pointwise score difference, $\text{score}_p - \text{score}_q$.

estimate Running mean score difference. Positive values favour forecaster p ; negative values favour forecaster q .

lower, upper Confidence sequence bounds. These are NA if `compute_cs = FALSE` or `cs_type = "none"`.

e_pq, e_qp One-sided e-processes. e_{pq} tests whether forecaster p outperforms q ; e_{qp} tests the reverse direction. These are NA if `compute_e = FALSE`.

Interpretation

The confidence sequence estimates the running average score advantage of p over q . If the whole interval lies above zero, the data favour p ; if the whole interval lies below zero, the data favour q .

The e-processes are evidence processes for one-sided null hypotheses. At level α , the two-sided rejection threshold used by `eprocess()` is $2 / \alpha$.

Examples

```
set.seed(1)
y <- rbinom(200, 1, 0.5)
p <- rep(0.5, 200)
q <- runif(200)

out <- compare_forecasts(p, q, y, scoring_rule = "brier")
tail(out)
```

crps_empirical	<i>Negated CRPS for empirical predictive distributions</i>
----------------	--

Description

Wrapper over `scoringRules::crps_sample` using `method = "edf"` (empirical distribution function, $O(n \log n)$ via quantile decomposition of Laio & Tamea, 2007). Positively oriented: higher = better.

Usage

```
crps_empirical(ensemble, y)
```

Arguments

ensemble	Matrix. $T \times n$ matrix of forecast draws. Each row corresponds to one observation in y and comprises n simulation draws from the predictive distribution. For Historical Simulation: each row is the past WINDOW returns.
y	Numeric vector of length T . Realised observations.

Details

Requires `nrow(ensemble) == length(y)`. Passes `dat = ensemble` directly to `crps_sample` which handles vectorisation over rows natively. `show_messages` is suppressed as the "edf" method requires no bandwidth selection messages.

Value

Numeric vector of length T of CRPS values in $(-\text{Inf}, 0]$ (negated loss).

Examples

```
if (requireNamespace("scoringRules", quietly = TRUE)) {
  ensemble <- matrix(c(0.1, 0.2, 0.3, 1.0, 1.1, 1.2), nrow = 2, byrow = TRUE)
  crps_ensemble(ensemble, y = c(0.25, 1.05))
}
```

crps_normal

*Negated CRPS for normal predictive distributions***Description**

Computes the Continuous Ranked Probability Score for a normal predictive distribution using `scoringRules::crps_norm()` and negates it so that higher values are better.

Usage

```
crps_normal(mu, sigma, x)
```

Arguments

<code>mu</code>	Numeric vector. Location parameters (conditional means).
<code>sigma</code>	Numeric vector. Scale parameters (conditional SDs, > 0).
<code>x</code>	Numeric vector. Realised observations.

Details

Calls `scoringRules::crps_norm(y = x, mean = mu, sd = sigma)` and negates. Use for GARCH(1,1)-norm forecasts where `mu` is the conditional mean and `sigma` is the conditional standard deviation.

Value

Numeric vector of CRPS values in $(-\infty, 0]$ (negated loss).

Examples

```
if (requireNamespace("scoringRules", quietly = TRUE)) {
  crps_normal(mu = c(0, 1), sigma = c(1, 2), x = c(0.2, 1.3))
}
```

crps_std

*Negated CRPS for Student-t predictive distributions***Description**

Wrapper over `scoringRules::crps_t`. Positively oriented: higher = better. The `dof > 2` constraint ensures finite variance, which is required for the CRPS to be well-defined for the t-distribution.

Usage

```
crps_std(mu, sigma, dof, x)
```

Arguments

<code>mu</code>	Numeric vector. Location parameters (conditional means).
<code>sigma</code>	Numeric vector. Scale parameters (conditional SDs, > 0).
<code>dof</code>	Numeric vector or scalar. Degrees of freedom (> 2). May be scalar if constant across all observations (e.g. estimated once per rolling window).
<code>x</code>	Numeric vector. Realised observations.

Details

Calls `scoringRules::crps_t(y = x, df = dof, location = mu, scale = sigma)` and negates. Use for GARCH(1,1)-std forecasts where `dof` is the estimated degrees-of-freedom parameter from `ugarchroll`.

Value

Numeric vector of CRPS values in $(-\text{Inf}, 0]$ (negated loss).

Examples

```
if (requireNamespace("scoringRules", quietly = TRUE)) {
  crps_std(mu = c(0, 1), sigma = c(1, 2), dof = 5, x = c(0.2, 1.3))
}
```

cs_asymptotic	<i>Asymptotic confidence sequence (EC.3, Eq. EC.29, Choe & Ramdas 2024)</i>
---------------	---

Description

Asymptotic, not finite-sample: coverage $\geq 1 - \alpha$ holds only as $t \rightarrow \infty$. Valid without requiring bounded score differences — requires only that $\hat{\Delta}_t$ has finite variance — so it's appropriate for tick loss and other scoring rules where hard bounds depend on unbounded realised values. Suitable for large evaluation windows.

Usage

```
cs_asymptotic(scores1, scores2, alpha = 0.05, t_star = NULL)
```

Arguments

scores1	Numeric vector. Scores for forecaster 1.
scores2	Numeric vector. Scores for forecaster 2.
alpha	Numeric in (0,1). Significance level. Default: 0.05.
t_star	Numeric > 0. Sample size at which CS is tightest. Default: length(scores1) (tightest at end of sample).

Details

$$C_t^A = \hat{\Delta}_t \pm \sqrt{\frac{2(t\sigma_t^2\rho^2 + 1)}{t^2\rho^2} \log \frac{\sqrt{t\sigma_t^2\rho^2 + 1}}{\alpha}}$$

where $\sigma_t^2 = \frac{1}{t} \sum_{i=1}^t (\hat{\delta}_i - \hat{\Delta}_{i-1})^2$ and ρ is tuned to be tightest at t_{star} :

$$\rho(t_{\text{star}}) = \sqrt{\frac{2 \log(1/\alpha) + \log(1 + 2 \log(1/\alpha))}{t_{\text{star}}}}$$

The running variance estimator uses the predictable mean $\hat{\Delta}_{t-1}$ (not the current mean $\hat{\Delta}_t$) to maintain predictability, with $\hat{\Delta}_0 := 0$.

Value

data.frame with columns t, estimate, lower, upper.

Examples

```
scores1 <- c(-0.4, -0.2, -0.3, -0.1, -0.2)
scores2 <- c(-0.5, -0.3, -0.4, -0.2, -0.3)
cs_asymptotic(scores1, scores2, alpha = 0.05)
```

cs_bernstein	<i>Empirical Bernstein confidence sequence (Theorem 2, Choe & Ramdas 2024)</i>
--------------	--

Description

Constructs a variance-adaptive time-uniform CS using empirical intrinsic time $\hat{V}_t = \sum_{i=1}^t (\hat{\delta}_i - \gamma_i)^2$. Tighter than the Hoeffding CS when score differences have low variance.

Usage

```
cs_bernstein(
  scores1,
  scores2,
  alpha = 0.05,
  c = 2,
  v_opt = 10,
  boundary = "mixture",
  gammas = NULL,
  lcb_only = FALSE,
  ucb_only = FALSE
)
```

Arguments

scores1	Numeric vector. Scores for forecaster 1.
scores2	Numeric vector. Scores for forecaster 2.
alpha	Numeric in (0,1). Significance level. Default: 0.05.
c	Numeric > 0. Sub-exponential scale. The process must satisfy $\text{lhat_delta_il} \leq c/2$. For score differences in $[a-b, b-a]$, $c = b - a$ (e.g. $c = 2$ for Brier score differences in $[-1, 1]$). Default: 2.
v_opt	Numeric > 0. Optimal intrinsic time. Default: 10.
boundary	Character. "mixture" (default, GE mixture) or "stitching" (polynomial stitched) or "hardcoded" (CR24 example formula, only valid for $\alpha=0.05, c=1$).
gammas	Numeric vector or NULL. Predictable centering sequence. If NULL, constructed as lagged running mean (default).
lcb_only	Logical. If TRUE, return lower CS only: $[\text{lower}, +\text{Inf}]$. Requires finite lower bound on hat_delta_i ; provide c .
ucb_only	Logical. If TRUE, return upper CS only: $(-\text{Inf}, \text{upper}]$.

Details

The CS is:

$$C_t^{EB} = \hat{\Delta}_t \pm u_{\alpha/2}^{GE}(\hat{V}_t; \rho, c) / t$$

Value

data.frame with columns t, estimate, lower, upper. lower = -Inf if ucb_only = TRUE; upper = Inf if lcb_only = TRUE.

Examples

```
scores1 <- c(-0.04, -0.09, -0.01, -0.16)
scores2 <- c(-0.09, -0.16, -0.04, -0.25)
cs_bernstein(scores1, scores2, alpha = 0.05)
```

cs_hoeffding	<i>Hoeffding-style confidence sequence (Theorem 1, Choe & Ramdas 2024)</i>
--------------	--

Description

Constructs a time-uniform confidence sequence for the mean score difference $\Delta_t = \frac{1}{t} \sum_{i=1}^t E[\hat{\delta}_i \mid \mathcal{F}_{i-1}]$.

Usage

```
cs_hoeffding(
  scores1,
  scores2,
  alpha = 0.05,
  c = 1,
  v_opt = 10,
  boundary = "mixture"
)
```

Arguments

scores1	Numeric vector. Scores $S(p_t, y_t)$ for forecaster 1.
scores2	Numeric vector. Scores $S(q_t, y_t)$ for forecaster 2.
alpha	Numeric in (0,1). Significance level. The CS has coverage $1 - \alpha$ uniformly over all t . Default: 0.05.
c	Numeric > 0 . Sub-Gaussian scale. The process must satisfy $ \text{hat_delta}_i \leq c$ for all i . For scores in $[a, b]$, the difference is in $[a-b, b-a]$, so $c = b - a$. Default: 1 (appropriate for Brier score differences in $[-1, 1]$).
v_opt	Numeric > 0 . Intrinsic time at which the CS is tightest. Default: 10 (recommended by CR24).
boundary	Character. "mixture" (default, recommended) or "stitching".

Value

data.frame with columns t, estimate, lower, upper.

Assumption

Requires $\hat{\Delta}_i$ to be c -sub-Gaussian given $\mathcal{F}_{i-1}$, i.e. $|\hat{\Delta}_i| \leq c$ for all i .

Boundary

$$C_t^H = \hat{\Delta}_t \pm u_{\alpha/2}^{CM}(c^2 t; \rho)/t$$

where u^{CM} is the normal mixture boundary and $c^2 t$ is the intrinsic time for a c -sub-Gaussian process with deterministic variance proxy. The intrinsic time for Theorem 1 is $v_t = c^2 * t$, not $v_t = t$: the CM boundary implicitly assumes 1-sub-Gaussian inputs, so the c^2 scaling must be applied explicitly. This matches the H21 convention, where the boundary absorbs the sub-Gaussian parameter via the variance process definition.

Relation to Python comparecast: Python uses $v_t = \text{sigma} * t$ where $\text{sigma} = (\text{hi} - \text{lo})/2 = c$. This is equivalent to our $c^2 * t$ only when $c = 1$. For $c \neq 1$ the parametrisations differ; we follow the paper.

Output

Returns a data.frame with one row per t and columns t , estimate (the running mean $\hat{\Delta}_t$), lower , and upper , with coverage guarantee $P(\forall t \geq 1 : \Delta_t \in [\text{lower}_t, \text{upper}_t]) \geq 1 - \alpha$.

Examples

```
scores1 <- c(-0.04, -0.09, -0.01, -0.16)
scores2 <- c(-0.09, -0.16, -0.04, -0.25)
cs_hoeffding(scores1, scores2, alpha = 0.05)
```

eprocess

Sub-exponential mixture e-process (Theorem 3, Choe & Ramdas 2024)

Description

Constructs two simultaneous one-sided e-processes for sequentially testing whether forecaster 1 (p) outperforms forecaster 2 (q) or vice versa.

Usage

```
eprocess(
  scores1,
  scores2,
  alpha = 0.05,
  c = 2,
  v_opt = 10,
  alpha_opt = NULL,
  gammas = NULL,
  clip_max = 1e+07
)
```

Arguments

scores1	Numeric vector. Scores $S(p_t, y_t)$ for forecaster 1.
scores2	Numeric vector. Scores $S(q_t, y_t)$ for forecaster 2.
alpha	Numeric in (0,1). Significance level. Rejection threshold is $2/\alpha$ for the two-sided test. Default: 0.05.
c	Numeric > 0 . Sub-exponential scale. Must satisfy $ \hat{\Delta}_i \leq c/2$ for all i . For score differences in $[-(b-a), b-a]$: $c = 2*(b-a)$. Default: 2 (for Brier score differences in $[-1, 1]$).
v_opt	Numeric > 0 . Intrinsic time at which e-process grows fastest. Default: 10 (recommended by CR24).
alpha_opt	Numeric in (0,1). One-sided alpha used to compute rho. Default: $\alpha/2$ (matches comparecast two-sided convention).
gammas	Numeric vector or NULL. Predictable centering sequence. If NULL, constructed as lagged running mean.
clip_max	Numeric. Maximum e-process value before clipping. Default: $1e7$ (matches Python comparecast).

Details

The mixture e-process at time t is:

$$E_t^{\text{mix}} = m(S_t, \hat{V}_t)$$

where $S_t = \sum_{i=1}^t \hat{\delta}_i$, $\hat{V}_t = \sum_{i=1}^t (\hat{\delta}_i - \gamma_i)^2$, and $m(s, v)$ is the Gamma-Exponential mixture function (Proposition EC.3, CR24).

VARIANCE PROCESS: The intrinsic time $V_{\hat{t}}$ uses NO floor. The GE mixture $m(s, v)$ is well-defined at $v=0$ (returns 1 when $s=0$), so no floor is needed. Adding a floor would yield less power in the e-process.

SCALE CONVENTION: c is the sub-exponential scale parameter such that $|\hat{\Delta}_i| \leq c/2$. This is the Theorems 2 & 3 convention from CR24. For Brier score differences in $[-1, 1]$: $c = 2$. For Winkler scores (bounded above by 1): $c = 2$.

LOG-SPACE: E-process values are computed in log-space and clipped before exponentiating to avoid numerical overflow.

Value

data.frame with the following columns:

t Time index.

e_pq, e_qp One-sided e-processes. e_pq tests $H_0^w(p, q)$: whether forecaster p outperforms q ; e_qp tests $H_0^w(q, p)$: whether forecaster q outperforms p .

log_e_pq, log_e_qp Log-scale values of the e-processes, clipped at $\log(\text{clip_max})$.

Rejection rule

At level α : reject $H_0^w(p, q)$ (conclude p outperforms q) when $e_{pq} \geq 2/\alpha$; reject $H_0^w(q, p)$ (conclude q outperforms p) when $e_{qp} \geq 2/\alpha$. Use `eprocess_rejections()` to extract the first crossing time for each.

Examples

```
scores1 <- c(-0.04, -0.09, -0.01, -0.16)
scores2 <- c(-0.09, -0.16, -0.04, -0.25)
ep <- eprocess(scores1, scores2, alpha = 0.05)
head(ep)
```

eprocess_betting

Betting-style e-process for the strong null hypothesis

Description

Implements the product-form test martingale

$$E_t = \prod_{r=1}^t (1 + \lambda_r \hat{\delta}_r)$$

for testing the strong null $H_0^s(p, q) : \delta_t \leq 0$ for all t .

Usage

```
eprocess_betting(scores1, scores2, c_t, lambda_t = NULL, clip_max = 1e+07)
```

Arguments

scores1	Numeric vector. Scores $S(p_t, y_t)$ for forecaster 1.
scores2	Numeric vector. Scores $S(q_t, y_t)$ for forecaster 2.
c_t	Numeric scalar or vector (same length as scores) of predictable bounds such that $ scores1 - scores2 \leq c_t / 2$ almost surely at every step.
lambda_t	Optional numeric vector of predictable betting fractions in $[0, 1/c_t]$. If NULL (default), uses the fixed fraction $\lambda_t = 1 / (2 * c_t)$.
clip_max	Numeric. Maximum e-process value before clipping. Default: $1e7$.

Value

data.frame with columns t, e_pq, e_qp, log_e_pq, log_e_qp.

Examples

```
set.seed(123)
T_sim <- 100
y <- rbinom(T_sim, size = 1, prob = 0.7)

# Forecaster 1 has a genuine edge (closer to true probability 0.7)
p1 <- runif(T_sim, 0.5, 0.9)
# Forecaster 2 is just guessing uniformly
p2 <- runif(T_sim, 0.1, 0.9)
```

```
# 1. Compute positively oriented Brier scores
s1 <- brier_score(p1, y)
s2 <- brier_score(p2, y)

# 2. Establish the bound.
# Brier scores lie in [-1, 0], so the maximum absolute difference is 1.
# eprocess_betting() requires |s1 - s2| <= c_t / 2, so c_t = 2 is globally valid.
res <- eprocess_betting(scores1 = s1, scores2 = s2, c_t = 2)

# The e-process e_pq tests the null that Forecaster 1 is NOT better than Forecaster 2.
# Because Forecaster 1 is genuinely better, e_pq accumulates massive evidence.
head(res)
tail(res, 3)
```

<code>eprocess_lag</code>	<i>Lag-h e-process for sequential forecast comparison (Propositions 5 & 6)</i>
---------------------------	--

Description

For h -step-ahead forecasts, constructs an anytime-valid e-process by stream splitting and combining h individual e-processes.

Usage

```
eprocess_lag(
  scores1,
  scores2,
  h = 1,
  alpha = 0.05,
  c = 2,
  v_opt = 10,
  null = "pw",
  calibrate = TRUE,
  cal_strategy = "mixture"
)
```

Arguments

<code>scores1</code>	Numeric vector. Scores for forecaster 1.
<code>scores2</code>	Numeric vector. Scores for forecaster 2.
<code>h</code>	Integer ≥ 1 . Forecast lag. For $h=1$, reduces to standard <code>eprocess()</code> — no splitting.
<code>alpha</code>	Numeric in $(0,1)$. Significance level. Default: 0.05.
<code>c</code>	Numeric > 0 . Sub-exponential scale. Default: 2.

<code>v_opt</code>	Numeric > 0. Default: 10.
<code>null</code>	Character. Null hypothesis type: • "pw" — period-wise weak null (average combination). • "w" — standard weak null (minimum combination).
<code>calibrate</code>	Logical. Apply p-to-e calibration. Default: TRUE.
<code>cal_strategy</code>	Character. "mixture" (default) or "simple".

Details

For $h = 1$: calls `eprocess()` directly and returns its output unchanged.

For $h \geq 2$: 1. Split x_s into h streams 2. Compute e-process on each stream independently 3. Combine using the appropriate null rule 4. Convert to p-process, combine, calibrate back to e-process 5. Unroll to original time scale

The period-wise ("pw") null is less conservative than the standard ("w") null but tests a different (weaker) hypothesis. See CR24 Section 4.4.

Value

data.frame with columns `t`, `e_pq`, `e_qp`, `log_e_pq`, `log_e_qp`.

Examples

```
scores1 <- c(-0.04, -0.09, -0.01, -0.16, -0.04, -0.09)
scores2 <- c(-0.09, -0.16, -0.04, -0.25, -0.09, -0.16)
ep <- eprocess_lag(scores1, scores2, h = 2, alpha = 0.05)
head(ep)
```

`eprocess_predictable` *Fixed-lambda e-process with predictable bounds (Proposition EC.7)*

Description

Constructs a valid e-process when score difference bounds vary over time but are predictable (known at time $i-1$ before observing $\hat{\Delta}_i$).

Usage

```
eprocess_predictable(
  scores1,
  scores2,
  c_seq,
  lambda = NULL,
  alpha = 0.05,
  gammas = NULL,
  clip_max = 1e+07,
  strict = FALSE
)
```

Arguments

scores1	Numeric vector. Scores for forecaster 1.
scores2	Numeric vector. Scores for forecaster 2.
c_seq	Numeric vector. Predictable bound sequence (c_i), same length as scores1. Must satisfy $ \text{scores1}[i] - \text{scores2}[i] \leq c_i/2$ and $c_i > 0$ for all i .
lambda	Numeric in $[0, 1/c_0]$. Betting parameter. Must be strictly less than $1/c_0$ where $c_0 = \max(c_seq)$. If NULL, uses the recommended default $\lambda = 0.5/c_0$.
alpha	Numeric in $(0,1)$. Significance level for rejection rule. Default: 0.05. Not used in computation, only for API consistency. Pass the same value to predictable_rejections() when evaluating rejection.
gammas	Numeric vector or NULL. Predictable centering sequence. If NULL, constructed as lagged running mean.
clip_max	Numeric. Maximum e-process value. Default: $1e7$.
strict	Logical. If TRUE, any violation of the bound condition at any time point will stop execution with an error. If FALSE (default), a warning is issued but the e-process is still computed. Note that violations invalidate the e-process guarantee, so <code>strict = TRUE</code> is recommended for formal inference.

Details

The e-process is computed as:

$$\log E_t(\lambda) = \sum_{i=1}^t \left[\lambda \hat{\delta}_i - \psi_{E, c_i}(\lambda) (\hat{\delta}_i - \gamma_i)^2 \right]$$

where

$$\psi_{E, c}(\lambda) = \frac{-\log(1 - c\lambda) - c\lambda}{c^2}$$

is evaluated at each step with the current c_i .

LAMBDA CHOICE: $\lambda = 0.5/c_0$ is a conservative default that stays well within the valid domain $[0, 1/c_0]$. For better power, λ can be tuned to the expected signal size, but must never reach $1/c_0$.

VALIDITY CHECK: The function verifies $|\hat{\delta}_i - \gamma_i| \leq c_i/2$ at each step and warns if violated. Violations invalidate the e-process guarantee.

Value

data.frame with columns: t, e_pq, e_qp, log_e_pq, log_e_qp, c_seq, lambda_used

Predictability

The bound sequence c_seq (and the centering sequence $gammas$) must be predictable: c_i is fixed and known at time $i - 1$, before $\text{scores1}[i]/\text{scores2}[i]$ (and hence $\hat{\delta}_i$) are observed — formally, c_i is $\mathcal{F}_{i-1}$ -measurable. A bound chosen after seeing $\hat{\delta}_i$ (e.g. derived from the realised data range) invalidates the e-process guarantee, even if it numerically satisfies $|\hat{\delta}_i - \gamma_i| \leq c_i/2$.

Examples

```
scores1 <- c(0.10, 0.20, 0.15, 0.25)
scores2 <- c(0.05, 0.10, 0.10, 0.20)
c_seq <- rep(1, length(scores1))
ep <- eprocess_predictable(scores1, scores2, c_seq = c_seq)
head(ep)
```

eprocess_rejections *Determine rejection times for an e-process output*

Description

Determine rejection times for an e-process output

Usage

```
eprocess_rejections(ep, alpha = 0.05)
```

Arguments

ep	data.frame. Output of eprocess().
alpha	Numeric. Significance level. Threshold is $2/\alpha$.

Value

Named list with elements:

- threshold — rejection threshold ($2 / \alpha$).
- tau_pq — first t where $e_{pq} \geq \text{threshold}$ (NA if never crossed).
- tau_qp — first t where $e_{qp} \geq \text{threshold}$ (NA if never crossed).
- reject_pq — logical: was $H_0^w(p, q)$ ever rejected?
- reject_qp — logical: was $H_0^w(q, p)$ ever rejected?

Examples

```
scores1 <- c(-0.04, -0.09, -0.01, -0.16)
scores2 <- c(-0.09, -0.16, -0.04, -0.25)
ep <- eprocess(scores1, scores2, alpha = 0.05)
eprocess_rejections(ep, alpha = 0.05)
```

ge_boundary	<i>Gamma-exponential mixture boundary</i>
-------------	---

Description

Gamma-exponential mixture boundary

Usage

```
ge_boundary(v, alpha, rho, c, s_lo = -10, s_hi = 500)
```

Arguments

v	Numeric vector ≥ 0 . Intrinsic time values.
alpha	Numeric in (0,1). ONE-SIDED significance level.
rho	Numeric > 0 . From rho_from_vopt().
c	Numeric > 0 . Sub-exponential scale.
s_lo	Numeric. Lower search bound for uniroot. Default: -10.
s_hi	Numeric. Upper search bound for uniroot. Default: 500.

Details

Computes $u_{GE}(v; \alpha, \rho, c) = \sup\{s : m(s, v) < 1/\alpha\}$ by solving $m(s, v) = 1/\alpha$ numerically for s via [uniroot\(\)](#), separately for each v_i .

Root-finding fallback: the search starts in $[s_lo, s_hi]$; if $m(s_hi, v_i)$ has not yet crossed the target, s_hi is doubled once and retried. If it still fails, a warning is issued and s_hi is returned as a conservative fallback value. Increase s_hi directly if this warning appears often (e.g. at large v or small α); increase $\text{abs}(s_lo)$ if no root is found at small v .

Computed elementwise; can be slow for long vectors — consider caching boundary values when the same (α , ρ , c) are reused.

Value

Numeric vector of boundary values (cumulative-sum scale).

Examples

```
rho <- rho_from_vopt(v_opt = 10, alpha = 0.025)
ge_boundary(v = 1:3, alpha = 0.025, rho = rho, c = 2)
```

lambda_betting_agrapa *Approximate GRAPA (aGRAPA) betting fractions for a constant bound*

Description

Constructs a predictable betting-fraction sequence for the product-form strong-null e-process (`eprocess_betting()`), by mapping the constant-bound score-difference stream onto $[0, 1]$ and applying the aGRAPA plug-in estimator of Waudby-Smith & Ramdas (2024), Online Supplementary Material, Section B.3, evaluated at the fixed null $m = 1/2$.

Usage

```
lambda_betting_agrapa(
  xs,
  c,
  kappa = 0.5,
  prior_mean = 0.5,
  prior_variance = 0.25,
  fake_obs = 1
)
```

Arguments

xs	Numeric vector. Score-difference stream $\hat{\delta}_t = S(p_t, y_t) - S(q_t, y_t)$.
c	Numeric > 0 . Either a scalar (constant bound) or a length-length(xs) predictable vector c_t (must satisfy $ xs_t \leq c_t/2$ pointwise, with c_t known before xs_t is observed). Scalars are recycled to a vector of the same length as xs.
kappa	Numeric in $(0, 1]$. WSR's internal truncation safeguard on the Y-scale $[-2*kappa, 2*kappa]$. Default 0.5. Note on saturation: Because seqcomp applies a strict $[0, 1]$ explicit projection <i>after</i> WSR's native truncation to ensure one-sided null validity, any $kappa \geq 0.5$ is mathematically equivalent. The lower bound $-2*kappa$ is always superseded by the 0 floor, and for $kappa \geq 0.5$, the upper bound $2*kappa \geq 1$ is superseded by the 1 cap. kappa only actively restricts the bet size when $kappa < 0.5$ (e.g., $kappa = 0.1$).
prior_mean	Numeric. Regularization prior for the running mean estimator on the Y-scale. Default 0.5 (WSR's own default; also happens to equal the fixed null $m = 1/2$ here, which is why the very first returned value is exactly 0).
prior_variance	Numeric in $(0, 0.25]$. Regularization prior for the running variance estimator. Default 0.25 (WSR's own default).
fake_obs	Numeric > 0 . Number of "fake observations" for regularization. Default 1 (WSR's own default).

Details

Derivation (constant c only): map $Y_t = \hat{\delta}_t/c + 1/2 \in [0, 1]$, so the null $\mu_t \leq 0$ becomes $\mathbb{E}[Y_t \mid \mathcal{F}_{t-1}] \leq 1/2$. WSR's aGRAPA plug-in is run exactly as published, fixed at $m = 1/2$, on the Y -scale running mean/variance. Two clips are then applied in sequence: WSR's own native truncation $[-2\kappa, 2\kappa]$, followed by an explicit floor/cap onto $[0, 1]$. The floor is important for the one-sided composite null $\mu \leq 0$ to remain valid, the played λ_t must be nonnegative on *every* round. WSR's raw aGRAPA value routinely goes negative whenever the running mean currently sits below the null, which happens routinely under the null itself. The cap enforces Arnold's own Prop 3.2 bound $\lambda_t \leq 1/c$, which is strictly tighter than WSR's native bankruptcy-avoidance bound.

The floor/cap projection and the $m = 1/2$ null hold identically for a predictable, time-varying c_t . Because c_t is $\mathcal{F}_{t-1}$ -measurable, $Y_t = x_t/c_t + 1/2$ still satisfies $\mathbb{E}[Y_t \mid \mathcal{F}_{t-1}] \leq 1/2$ under the strong null, and the same aGRAPA plug-in can be run unmodified at the fixed null $m = 1/2$; only the final $Y \mapsto d$ conversion ($\lambda_t Y / c$) becomes pointwise. This time-varying extension is original to this package.

Value

Numeric vector of length `length(xs)`: predictable λ_t values in $[0, 1/c]$, for use as `eprocess_betting()`'s `lambda_t` argument together with `c_t = c`.

References

Waudby-Smith, I. and Ramdas, A. (2024). Estimating means of bounded random variables by betting. *Journal of the Royal Statistical Society Series B: Statistical Methodology*, 86(1), 1–27.

Examples

```
xs <- c(0.6, -0.2, 0.4)
lambda_betting_agrapa(xs, c = 2)
```

lambda_betting_ons	<i>Online Newton Step (ONS-m) betting fractions for a constant bound</i>
--------------------	--

Description

Constructs a predictable betting-fraction sequence for the product-form strong-null e-process (`eprocess_betting()`), by mapping the constant-bound score-difference stream onto $[0, 1]$ and running WSR's ONS-m algorithm (Waudby-Smith & Ramdas 2024, Online Supplementary Material, Algorithm 1, Section B.5), projected each round onto $[0, 1]$ instead of their native symmetric box, and evaluated at the fixed null $m = 1/2$.

Usage

```
lambda_betting_ons(xs, c, eta = 2/(2 - log(3)))
```

Arguments

xs	Numeric vector. Score-difference stream $\hat{\delta}_t = S(p_t, y_t) - S(q_t, y_t)$.
c	Numeric > 0 . Either a scalar (constant bound) or a length-length(xs) predictable vector c_t (must satisfy $ xs_t \leq c_t/2$ pointwise, with c_t known before xs_t is observed). Scalars are recycled to a vector of the same length as xs.
eta	Numeric > 0 . WSR's ONS step-size constant. Default $2 / (2 - \log(3))$, WSR's own stated value (natural log).

Details

Same $Y_t = xs_t/c + 1/2$ mapping as [lambda_betting_agrapa\(\)](#). WSR's Algorithm 1 is run exactly as stated, but using the standard OCO gradient for minimizing the negative log-wealth $-\log(1 + \lambda y_t)$, and replacing WSR's own projection target $[-c/(1-m), c/m]$ with $[0, 1]$. Following generic online-convex-optimization theory, λ_t^O is projected onto $[0, 1]$ before being used to compute the next round's gradient. Validity of the resulting e-process only requires the played $\lambda_{d,t}$ to lie in $[0, 1/c]$ regardless of derivation.

This modification is original to this package.

Value

Numeric vector of length length(xs): predictable λ_t values in $[0, 1/c]$, for use as eprocess_betting()'s lambda_t argument together with c_t = c.

References

Waudby-Smith, I. and Ramdas, A. (2024). Estimating means of bounded random variables by betting. *Journal of the Royal Statistical Society Series B: Statistical Methodology*, 86(1), 1–27.

Examples

```
xs <- c(0.6, -0.2, 0.4)
lambda_betting_ons(xs, c = 2)
```

lambda_betting_quantile

Adaptive betting fraction for quantile-forecast strong-null tests

Description

Implements the quantile-specific adaptive betting scheme of Arnold et al. (2026), translating their log-scale loss convention to seqcomp's positively-oriented scores.

Usage

```
lambda_betting_quantile(p_t, q_t, tau, delta_hat_lag1 = NULL, eps = 1e-08)
```

Arguments

p_t	Numeric vector. Quantile forecasts of the first forecaster.
q_t	Numeric vector. Quantile forecasts of the second forecaster.
tau	Numeric scalar in (0, 1). The quantile level.
delta_hat_lag1	Numeric vector. The <i>previous</i> step's score difference. Must have <code>delta_hat_lag1[1] = 0</code> .
eps	Positive numeric scalar. Used as a denominator safeguard against zero. Default: <code>1e-8</code> .

Details

If two forecasts are identical at a time point, their tick-loss difference and its analytic bound are both zero. To accommodate the strictly positive bound required by `eprocess_betting()` and the adaptive betting rules, this function replaces such bounds (and any smaller bounds) by `eps`. This is a conservative predictable enlargement of the bound; the corresponding Arnold betting fraction is capped at `1 / c_safe`.

Value

A list with `c_t` and `lambda_t` vectors for `eprocess_betting()`.

Note

Scale Translation: This function assumes `p_t`, `q_t`, and `delta_hat_lag1` are calculated on the raw, linear scale. To replicate the exact log-scale bounds used in the Arnold et al. (2026) Covid-19 case study, the forecast vectors passed to this function must be log-transformed prior to evaluation.

Examples

```
set.seed(456)
T_sim <- 100
y <- rnorm(T_sim)
tau <- 0.90

# Forecaster 1 correctly predicts the true 90th percentile (~1.28)
p_t <- rep(qnorm(tau), T_sim)
# Forecaster 2 is biased and incorrectly predicts the median (0.0)
q_t <- rep(0.0, T_sim)

# 1. Compute pointwise tick loss (positively oriented)
s_p <- tick_loss(p_t, y, tau)
s_q <- tick_loss(q_t, y, tau)

# 2. Compute the lagged score difference required by Arnold's heuristic
xs <- s_p - s_q
delta_lag1 <- c(0, head(xs, -1))

# 3. Generate the predictable bounds (c_t) and betting fractions (lambda_t)
bnds <- lambda_betting_quantile(p_t, q_t, tau, delta_hat_lag1 = delta_lag1)
```

```
# 4. Plug these directly into the betting e-process
res <- eprocess_betting(
  scores1 = s_p,
  scores2 = s_q,
  c_t = bnds$c_t,
  lambda_t = bnds$lambda_t
)

# View the final evidence accumulation
tail(res[, c("t", "e_pq", "e_qp")], 3)
```

log_score

*Logarithmic score for binary and categorical forecasts***Description**

Computes the positively oriented logarithmic score. Vector probability input is treated as binary; matrix probability input is treated as categorical.

Usage

```
log_score(p, y, eps = 1e-15)
```

Arguments

p	Numeric vector in $[0, 1]$ for binary forecasts, or a numeric matrix whose rows are probability vectors for categorical forecasts.
y	For binary vector input, numeric vector in $\{0, 1\}$. For categorical matrix input, integer vector in $\{1, \dots, K\}$, where $K = \text{ncol}(p)$.
eps	Numeric. Probability floor used before taking logarithms. Default is $1e-15$. Set to 0 to disable clipping.

Details

For binary forecasts, this computes

$$S(p, y) = y \log(p) + (1 - y) \log(1 - p).$$

For categorical forecasts, this computes

$$S(\mathbf{p}, y) = \log(p_y),$$

where p_y is the forecast probability assigned to the realised category.

Value

Numeric vector of scores in $(-\text{Inf}, 0]$. Higher is better.

Use with seqcomp

The logarithmic score is unbounded below. It should not be used directly with the finite-sample bounded-difference confidence sequences or e-processes. For binary outcomes, use `winkler_score()` and `winkler_cs()` when the Winkler construction is appropriate. For unbounded score differences, use `cs_asymptotic()` or supply genuine predictable bounds to `eprocess_predictable()`.

Examples

```
p <- c(0.2, 0.7, 0.9)
y <- c(0, 1, 1)
log_score(p, y)
```

predictable_rejections

Summarise predictable bounds e-process

Description

Summarise predictable bounds e-process

Usage

```
predictable_rejections(ep, alpha = 0.05)
```

Arguments

<code>ep</code>	data.frame. Output of <code>eprocess_predictable()</code> .
<code>alpha</code>	Numeric. Significance level.

Value

Named list matching the `eprocess_rejections()` format (`threshold`, `tau_pq`, `tau_qp`, `reject_pq`, `reject_qp`), plus:

- `c_range` — range of `c_seq` used.
- `lambda` — lambda value used.

Examples

```
scores1 <- c(0.10, 0.20, 0.15, 0.25)
scores2 <- c(0.05, 0.10, 0.10, 0.20)
c_seq <- rep(1, length(scores1))
ep <- eprocess_predictable(scores1, scores2, c_seq = c_seq)
predictable_rejections(ep, alpha = 0.05)
```

ps_boundary	<i>Polynomial stitched (PS) boundary</i>
-------------	--

Description

Alternative boundary for both Theorem 1 and Theorem 2 constructions, included for completeness and for cross-checking CR24 Table results.

Usage

```
ps_boundary(v, alpha, v_opt = 10, c = 1, s = 1.4, eta = 2)
```

Arguments

v	Numeric vector ≥ 0 . Intrinsic time values.
alpha	Numeric in (0,1). ONE-SIDED significance level.
v_opt	Numeric > 0 . Optimal intrinsic time (= m in H21). Default: 10.
c	Numeric > 0 . Sub-exponential scale.
s	Numeric > 1 . Stitching parameter. Default: 1.4.
eta	Numeric > 1 . Geometric spacing. Default: 2.

Details

This is **not** the recommended primary boundary: the CM/GE mixture boundaries (`cm_boundary()`, `ge_boundary()`) are tighter in CR24 and are used by default throughout seqcomp. Use `ps_boundary()` only when you specifically need the polynomial-stitched construction.

Value

Numeric vector of boundary values (cumulative-sum scale).

Examples

```
ps_boundary(v = 1:5, alpha = 0.025, v_opt = 10, c = 1)
```

qlike_score	<i>Negated QLIKE score for variance forecasts</i>
-------------	---

Description

Computes the positively oriented (negated) QLIKE quasi-likelihood loss for variance forecasts.

Usage

```
qlike_score(sigma2_hat, sigma2)
```

Arguments

sigma2_hat	Numeric vector. Forecast variance (strictly positive).
sigma2	Numeric vector. Realised variance (strictly positive).

Details

Standard QLIKE loss is

$$L_{QL}(\hat{\sigma}^2, \sigma^2) = \frac{\sigma^2}{\hat{\sigma}^2} - \log \frac{\sigma^2}{\hat{\sigma}^2} - 1.$$

This is loss-oriented (lower = better, minimum 0 at a perfect forecast), so the function negates it: $S_{QL} = -L_{QL}$.

Literature note: some sources define QLIKE as $\log(\text{sigma2_hat}) + \text{sigma2} / \text{sigma2_hat}$, which differs by constants from the form above. Here the loss is normalised to have minimum 0 and is then negated for positive orientation.

Value

Numeric vector of negated QLIKE scores. Higher is better. Maximum value is 0, achieved at a perfect forecast $\text{sigma2_hat} = \text{sigma2}$. Unbounded below.

Unbounded below

QLIKE is unbounded below. It should not be used directly with the finite-sample bounded-difference confidence sequences or e-processes. Use `cs_asymptotic()` for QLIKE-based confidence sequences, or use `eprocess_predictable()` only when genuine ex ante predictable bounds are available. QLIKE is not compatible with the Winkler construction because Winkler scores are restricted to binary outcomes and probability forecasts.

Examples

```
sigma2_hat <- c(1.0, 1.5, 2.0)
sigma2 <- c(1.1, 1.4, 2.2)
qlike_score(sigma2_hat, sigma2)
```

rho_from_vopt	<i>Convert optimal intrinsic time to rho</i>
---------------	--

Description

Maps the user-specified intrinsic time v_{opt} (the point at which a boundary is tightest) to the ρ tuning parameter, via the Lambert W formula of Howard et al. (2021), Proposition 3 (paper-exact).

Usage

```
rho_from_vopt(v_opt = 10, alpha = 0.025)
```

Arguments

v_{opt}	Numeric > 0. Intrinsic time at which the boundary is tightest. Recommended default from CR24: 10.
alpha	Numeric in (0,1). Significance level (one-sided). For a two-sided boundary at level alpha, pass alpha/2 here.

Details

$$\rho = \frac{v_{\text{opt}}}{-W_{-1}(-\alpha^2/e) - 1}$$

The lower branch W_{-1} is defined for x in $[-1/e, 0)$ and returns values ≤ -1 . For alpha in $(0, 1)$, $-\alpha^2/e$ is always in $(-1/e, 0)$, so the branch is well-defined.

Value

Numeric > 0. The ρ tuning parameter.

Examples

```
rho_from_vopt(v_opt = 10, alpha = 0.025)
```

score_bounds	<i>Score difference bounds for a named scoring rule</i>
--------------	---

Description

Returns lo, hi and the derived scale parameters c_{thm1} , c_{thm23} for the score difference process $\hat{\Delta}_t = S(p, y) - S(q, y)$, in those cases where a genuine, theorem-valid bound is available.

Usage

```
score_bounds(scoring_rule)
```

Arguments

`scoring_rule` Character. One of:

- "brier", "spherical" — bounded, exact finite-sample c .
- "winkler" — descriptive helper for the one-sided CS on the log score.
- "tick" — unbounded; returns NULL with guidance.
- "crps", "crps_normal", "crps_empirical", "crps_std" — unbounded; returns NULL with guidance.
- "log", "qlike" — unbounded; returns NULL with guidance.

Details

Convention (`utils.R::score_diff_scales`): $c_{\text{thm1}} = (hi - lo) / 2$ # Theorem 1: $|\delta_{il}| \leq c_{\text{thm23}}$
 $= hi - lo$ # Theorems 2 & 3: $|\delta_{il}| \leq c/2$

Value

Named list with elements `lo`, `hi`, `c_thm1`, `c_thm23` for bounded rules, or NULL for unbounded rules (with an informative message).

Per-rule notes

- **Brier / Spherical** — individual scores lie in $[-1, 0]$ (Brier) or $[0, 1]$ (Spherical), so score differences lie in $[-1, 1]$ either way. This bound is exact and yields finite-sample anytime-valid CS via Hoeffding/Bernstein.
- **Winkler** — bounded above by 1; the lower bound is problem-dependent, so `lo` = `-Inf` and only `hi` = 1 is used, as a descriptive helper for the one-sided CS wrapper `winkler_cs()`. Not intended for generic Hoeffding/Bernstein use (Theorem 1 requires a finite symmetric interval).
- **Tick loss** — unbounded on general financial returns. Any bound derived from an empirical data range is ex-post and not filtration-respecting, so it cannot justify finite-sample anytime validity. Use `cs_asymptotic()` for tick comparisons.
- **CRPS** (normal, t, empirical) — unbounded, since both the predictive distributions and the realised outcomes are unbounded. A historical data range is again an ex-post surrogate and does not provide a theorem-valid c for Hoeffding/Bernstein. Use `cs_asymptotic()`, or supply genuine ex ante bounds in problem-specific code if available.
- **Log / QLIKE** — both unbounded. For binary log-score comparisons, use `winkler_score()` + `winkler_cs()` when the Winkler construction is appropriate. For categorical log-score, QLIKE, and other unbounded score differences, use `cs_asymptotic()`, or `eprocess_predictable()` only with genuine ex ante predictable bounds.

Examples

```
score_bounds("brier")
score_bounds("winkler")
```

score_diff_scales	<i>Score difference bounds -> sub-Gaussian / sub-exponential scale</i>
-------------------	---

Description

Given score-difference bounds $[lo, hi]$, computes the two scale constants used elsewhere in seqcomp:

- $c_thm1 = (hi - lo) / 2$ — sub-Gaussian scale for Theorem 1, where $|\delta_{t,i}| \leq c_thm1$.
- $c_thm23 = hi - lo$ — sub-exponential scale for Theorems 2 & 3, where $|\delta_{t,i}| \leq c_thm23 / 2$.

Usage

```
score_diff_scales(lo, hi)
```

Arguments

lo	Numeric. Lower bound of score difference (usually $a - b$).
hi	Numeric. Upper bound of score difference (usually $b - a$).

Details

Both conventions bound the same quantity: after centering, $\delta_{t,i}$ lies in $[-(hi-lo)/2, (hi-lo)/2]$, so $\max |\delta_{t,i}| = (hi-lo)/2$, which equals both c_thm1 and $c_thm23 / 2$.

Value

Named list with elements `c_thm1` and `c_thm23`.

Examples

```
score_diff_scales(lo = -1, hi = 1)
```

smcs_compare	<i>Compare Multiple Sequential Forecasters (SMCS)</i>
--------------	---

Description

Evaluates an arbitrary number of candidate forecasters simultaneously, constructing Sequential Model Confidence Sets (SMCS) that maintain family-wise error rate control over time.

Usage

```
smcs_compare(
  forecasts,
  outcomes,
  scoring_rule = c("brier", "spherical", "tick"),
  cs_method = c("bernstein", "hoeffding"),
  tau = NULL,
  alpha = 0.05,
  v_opt = 10,
  clip_max = 1e+07,
  betting_rule = c("default", "naive", "agrapa", "ons", "arnold"),
  period = 1,
  ...
)
```

Arguments

forecasts	A $T \times m$ matrix of forecasts. For binary/categorical probability forecasts, these should be matrices of probabilities. For quantile forecasts, these should be raw predicted quantiles. <i>Note:</i> to replicate the log-scale bounds of the Arnold et al. (2026) Covid-19 study, pass log-transformed forecasts and outcomes.
outcomes	A numeric vector of T realised outcomes.
scoring_rule	Character. Scoring rule used to compare forecasts. Currently supports "brier", "spherical", and "tick".
cs_method	Character. Confidence sequence method for the weak null: "bernstein" or "hoeffding". Default is "bernstein".
tau	Numeric in $(0, 1)$. The quantile level. Required only if <code>scoring_rule = "tick"</code> .
alpha	Numeric in $(0, 1)$. Family-wise significance level. Default is 0.05.
v_opt	Numeric > 0 . Intrinsic time at which the weak-null confidence sequence is tuned to be tightest. Default is 10.
clip_max	Numeric. Maximum e-process value before clipping in the strong-null test. Default is 1e7.
betting_rule	Character. Strong-null betting-fraction rule. "default" preserves the existing behavior: "naive" for "brier" and "spherical", and "arnold" for "tick". "naive", "agrapa", and "ons" are available for all scoring rules; "arnold" is available only for "tick".
period	Positive integer. Number of interleaved periodic sub-streams used by "agrapa" or "ons". Must be 1 for "naive" and "arnold".
...	Additional arguments passed to build_agrapa_betting_array() or build_ons_betting_array() for the selected adaptive rule, such as <code>kappa</code> , <code>prior_mean</code> , <code>prior_variance</code> , <code>fake_obs</code> , or <code>eta</code> .

Details

This is a high-level wrapper that automates pointwise score calculation, boundary generation, and multiplicity corrections via `smcs_strong()` and `smcs_weak()`. For uniformly bounded scoring rules, it returns SMCSs under the strong, uniformly weak, and time-varying weak null hypotheses. For conditionally bounded rules like "tick" loss, it automatically builds the dynamic 3D arrays required for strong-null adaptive betting.

Value

A list containing:

`scores` A $T \times m$ matrix of evaluated pointwise scores.

`smcs_strong` A $T \times m$ logical matrix tracking inclusion in the strong-null SMCS over time (permanent exclusions).

`smcs_uniform_weak` A $T \times m$ logical matrix tracking inclusion in the uniformly weak SMCS over time (permanent exclusions). Currently NULL for "tick" loss.

`smcs_weak` A $T \times m$ logical matrix tracking inclusion in the time-varying weak-null SMCS over time (models can exit and re-enter). Currently NULL for "tick" loss.

`betting_rule` The resolved strong-null betting rule actually used.

`period` The period supplied for an adaptive betting rule.

`betting_args` The additional adaptive-rule settings supplied through ...

Examples

```
set.seed(42)
T_sim <- 100
y <- rbinom(T_sim, 1, 0.5)

# Create 3 forecasters:
# M1 is a perfect oracle (always predicts the true y)
# M2 is slightly noisy (adds small uniform noise to y)
# M3 is an anti-oracle (predicts the exact opposite of y)
fcsts <- matrix(NA, nrow = T_sim, ncol = 3)
fcsts[, 1] <- y
fcsts[, 2] <- abs(y - runif(T_sim, 0, 0.1))
fcsts[, 3] <- 1 - y
colnames(fcsts) <- c("M1", "M2", "M3")

out <- smcs_compare(fcsts, y, scoring_rule = "brier")

# Print the object to see exclusions (M3 will be dropped rapidly)
out

# View how the set sizes shrink over time
summary(out)

# Additionally, use an adaptive strong-null betting rule
out_agrapa <- smcs_compare(
  fcsts, y, scoring_rule = "brier", betting_rule = "agrapa"
```

```
)
out_agrapa
```

smcs_strong

Sequential Model Confidence Set (Strong & Uniformly Weak Null)

Description

Constructs a Sequential Model Confidence Set (SMCS) evaluating a family-wise intersection null hypothesis. By maintaining a running intersection over time, any model excluded from the set is permanently eliminated.

Usage

```
smcs_strong(
  scores,
  alpha = 0.05,
  method = c("betting", "mixture"),
  c_param = NULL,
  lambda_param = NULL,
  ...
)
```

Arguments

scores	A $T \times m$ matrix of positively-oriented scores.
alpha	Numeric in $(0, 1)$. Family-wise significance level. Default is 0.05.
method	Character. "betting" (uses eprocess_betting() , true strong null) or "mixture" (uses eprocess() , tests uniformly weak null). Default is "betting".
c_param	Numeric scalar, $m \times m$ matrix, or $T \times m \times m$ array. The predictable bound parameter. Required. Time-varying arrays are only allowed if method = "betting".
lambda_param	Optional parameter for betting fractions, matching the shape allowed for c_param. Only used if method = "betting".
...	Additional arguments passed to the underlying pairwise e-process function (e.g., v_opt and clip_max for "mixture"; clip_max for "betting").

Details

Depending on the method chosen, this function tests different hypotheses:

- method = "betting": Tests the **Strong Null** hypothesis (conditional step-by-step superiority). Uses a product-form betting martingale.

- `method = "mixture"`: Tests the **Uniformly Weak Null** hypothesis (average superiority over time). Uses an exponential-mixture martingale. Because strong superiority implies uniform weak superiority, feeding "mixture" into this closed-testing machinery yields a valid (though strictly testing the uniformly weak null) SMCS.

The function computes pairwise e-processes between all models, constructs an intersection e-process for each model, applies a closed-testing multiplicity adjustment via `vovk_wang_merge()`, and permanently excludes models when their adjusted e-value exceeds $1/\alpha$.

Value

A list containing:

`E_i_dot` A $T \times m$ matrix of unadjusted intersection e-processes.

`E_star` A $T \times m$ matrix of closed-testing adjusted e-processes.

`smcs` A $T \times m$ logical matrix. TRUE indicates the model remains in the SMCS at time t .

Examples

```
set.seed(1)
# 3 models, 100 time steps. Model 3 is artificially much worse.
scores <- matrix(runif(300, -0.5, 0), nrow = 100, ncol = 3)
scores[, 3] <- scores[, 3] - 0.5
colnames(scores) <- c("M1", "M2", "M3")

# Using the betting method with a global bound c = 2
res <- smcs_strong(scores, alpha = 0.05, method = "betting", c_param = 2)
tail(res$smcs)
```

smcs_weak

Sequential Model Confidence Set (Weak Null)

Description

Constructs a Sequential Model Confidence Set (SMCS) evaluating the weak null hypothesis that a model outperforms all other candidate models on average over time.

Usage

```
smcs_weak(
  scores,
  alpha = 0.05,
  cs_method = c("bernstein", "hoeffding"),
  c_param = NULL,
  ...
)
```

Arguments

scores	A $T \times m$ matrix of positively-oriented scores.
alpha	Numeric in $(0, 1)$. Family-wise significance level. Default is 0.05.
cs_method	Character. "bernstein" (uses <code>cs_bernstein()</code>) or "hoeffding" (uses <code>cs_hoeffding()</code>).
c_param	Numeric scalar or $m \times m$ matrix. The uniform bound parameter. Must be constant over time.
...	Additional arguments passed to the chosen CS function (e.g., <code>v_opt</code>).

Details

Uses a joint confidence sequence decoupling result: a model i remains in the SMCS at time t if and only if, for every competitor j , the pairwise $(1 - \alpha/(m(m - 1)))$ -confidence sequence for the average score difference does not strictly rule out that i is better than j .

Unlike the strong null, this SMCS does not maintain a strict running intersection; a model's average score can recover over time, allowing it to dynamically exit and re-enter the confidence set.

Value

A list containing:

`smcs` A $T \times m$ logical matrix. TRUE indicates the model is in the weakly superior set at time t .
`alpha_adjusted` The Bonferroni-adjusted significance level applied to each pairwise sequence.

Examples

```
set.seed(2)
scores <- matrix(runif(300, -0.5, 0), nrow = 100, ncol = 3)
scores[, 3] <- scores[, 3] - 0.5
colnames(scores) <- c("M1", "M2", "M3")

res <- smcs_weak(scores, alpha = 0.05, cs_method = "bernstein", c_param = 2)
tail(res$smcs)
```

spherical_score	<i>Spherical score for binary and categorical forecasts</i>
-----------------	---

Description

Computes the positively oriented spherical score. Vector probability input is treated as binary; matrix probability input is treated as categorical.

Usage

```
spherical_score(p, y)
```

Arguments

p	Numeric vector in $[0, 1]$ for binary forecasts, or a numeric matrix whose rows are probability vectors for categorical forecasts.
y	For binary vector input, numeric vector in $\{0, 1\}$. For categorical matrix input, integer vector in $\{1, \dots, K\}$, where $K = \text{ncol}(p)$.

Details

For binary forecasts, this computes

$$S(p, y) = \frac{py + (1 - p)(1 - y)}{\sqrt{p^2 + (1 - p)^2}}.$$

For categorical forecasts, this computes

$$S(\mathbf{p}, y) = \frac{p_y}{\|\mathbf{p}\|_2},$$

where p_y is the forecast probability assigned to the realised category.

Score differences lie in $[-1, 1]$, so use $c = 1$ for Theorem 1 and $c = 2$ for Theorems 2 and 3.

Value

Numeric vector of scores in $[0, 1]$. Higher is better.

Examples

```
p <- c(0.2, 0.7, 0.9)
y <- c(0, 1, 1)
spherical_score(p, y)
```

split_streams

Split a sequence into h interleaved lag streams

Description

For lag h , the k -th stream ($k = 1, \dots, h$) contains indices $\{k, k + h, k + 2h, \dots\}$, following the CR24 convention.

Usage

```
split_streams(xs, h)
```

Arguments

xs	Numeric vector. Score differences $\hat{\Delta}_t$.
h	Integer ≥ 1 . Lag (number of steps ahead).

Value

List of length h . Each element is a numeric vector containing the score differences for that stream.

Examples

```
split_streams(1:10, h = 3)
# stream 1: indices 1, 4, 7, 10
# stream 2: indices 2, 5, 8
# stream 3: indices 3, 6, 9
```

tick_loss	<i>Negated tick loss for quantile forecasts</i>
-----------	---

Description

Computes the positively oriented (negated) tick/quantile loss (Koenker & Bassett, 1978).

Usage

```
tick_loss(q, y, tau)
```

Arguments

q	Numeric vector. Quantile forecasts at level tau.
y	Numeric vector. Realised outcomes.
tau	Numeric in (0,1). Quantile level.

Details

The standard tick loss is

$$\rho_{\tau}(u) = u (\tau - \mathbb{I}(u < 0)) ,$$

where $u = y - q_{\tau}$ is the forecast error. This is loss-oriented (lower = better), so the function negates it:

$$S_T(q, y; \tau) = -(y - q) (\tau - \mathbb{I}(y < q)) .$$

Tick loss is unbounded on general real-valued outcomes. Bounds derived from an empirical data range are ex-post and do not provide theorem-valid constants for finite-sample Hoeffding/Bernstein confidence sequences or e-processes.

Sign convention: the negation means $\text{hat_delta_t} > 0$ when forecaster p has smaller tick loss, hence a better quantile forecast, than forecaster q .

Value

Numeric vector of negated tick loss scores. Higher = better.

Examples

```
q <- c(1.0, 1.5, 2.0)
y <- c(1.2, 1.4, 2.3)
tick_loss(q, y, tau = 0.5)
```

unroll_stream

Unroll a stream-wise quantity back to the original time scale

Description

After computing a per-stream cumulative quantity (e.g. e-process values), restores them to the original length T by zero-padding the first $k - 1$ positions, repeating each stream value h times, then truncating to length T .

Usage

```
unroll_stream(stream_vals, k, h, T_)
```

Arguments

stream_vals	Numeric vector. Values for stream k (length $\sim T/h$).
k	Integer. Stream index (1-based).
h	Integer. Lag.
T_	Integer. Total original sequence length.

Value

Numeric vector of length $T_$.

Alignment only, not theoretical updating

For lagged forecasts ($h \geq 2$), the returned series is aligned to the evaluated score-difference index after stream splitting. It should **not** be interpreted as a process that updates at the original forecast-issuance time. The unrolled process is for visualization and alignment only; the theoretical validity argument relies strictly on the streamwise sub-filtrations, not on this unrolled representation.

Examples

```
unroll_stream(c(1, 2, 3), k = 2, h = 2, T_ = 6)
```

vovk_wang_merge

Arithmetic-mean closed-testing e-value merge

Description

Given a vector of e-values $e_1, \dots, e_m$ (where each e_i represents the evidence against the intersection null hypothesis for model i), computes the closed-testing adjusted e-values using the arithmetic mean as the e-merging function.

Usage

```
vovk_wang_merge(e_values)
```

Arguments

`e_values` Numeric vector of non-negative e-values, one per model.

Details

This implements the Accelerated E-Value Calibration algorithm (Tim Stephan, ETH Zurich; Section H of the supplementary material to Arnold et al., 2026), which solves the Vovk & Wang (2021) closed-testing minimization in $O(m \log m)$ time rather than the naive $O(m^2)$.

Value

A numeric vector of the same length, containing the closed-testing adjusted e-values e_i^* , in the original (unsorted) order.

Examples

```
raw_values <- c(10, 2, 1)
# Model 1 has strong evidence against it, Model 3 has none.
vovk_wang_merge(raw_values)
```

winkler_compare

Full Winkler comparison pipeline (Proposition EC.4)

Description

Convenience wrapper that computes Winkler scores, one-sided CS, and e-process in a single call.

Usage

```
winkler_compare(
  p,
  q,
  y,
  alpha = 0.05,
  base_score = log_score,
  v_opt = 10,
  lower_bound = NULL
)
```

Arguments

p	Numeric vector in (0,1).
q	Numeric vector in (0,1).
y	Numeric vector containing only 0 and 1. Binary outcomes.
alpha	Numeric in (0,1). Default: 0.05.
base_score	Function. Default: log_score.
v_opt	Numeric > 0. Default: 10.
lower_bound	Numeric or NULL. See winkler_cs().

Value

Named list with elements:

- `winkler_scores` — raw Winkler score vector.
- `cs` — `data.frame` from `winkler_cs()`.
- `etest_p_worse` — one-sided e-process testing whether p is worse than q.
- `etest_q_worse` — one-sided e-process testing whether q is worse than p.
- `rejections` — list of one-sided rejection summaries.

Examples

```
p <- c(0.7, 0.6, 0.8, 0.65)
q <- c(0.5, 0.7, 0.6, 0.55)
y <- c(1, 1, 0, 1)
winkler_compare(p, q, y, alpha = 0.05)
```

winkler_cs	<i>One-sided empirical Bernstein CS for Winkler scores (Proposition EC.4)</i>
------------	---

Description

Applies the Winkler normalisation and constructs a one-sided upper confidence sequence for the mean Winkler score $W_t = (1/t) \sum w_i$. The CS takes the form $(-\text{Inf}, U_t]$, valid uniformly over all $t \geq 1$.

Usage

```
winkler_cs(
  p,
  q,
  y,
  alpha = 0.05,
  base_score = log_score,
  v_opt = 10,
  lower_bound = NULL
)
```

Arguments

p	Numeric vector in (0,1). Forecasts from model 1.
q	Numeric vector in (0,1). Forecasts from model 2.
y	Numeric vector containing only 0 and 1. Binary outcomes.
alpha	Numeric in (0,1). Significance level. Default: 0.05.
base_score	Function. Underlying scoring rule. Default: log_score.
v_opt	Numeric > 0. Optimal intrinsic time. Default: 10.
lower_bound	Numeric or NULL. Analytical lower bound on w_i for two-sided CS via Corollary 2. If NULL (default), returns one-sided CS only. If supplied, must satisfy $w_i \geq \text{lower_bound}$ for all i almost surely.

Details

Scale convention: Winkler score bounded above by 1, so $c/2 = 1$, $c = 2$. This is hardcoded — do not change c without re-deriving the bound.

Value

data.frame with columns t, estimate, lower, upper. lower = -Inf always (one-sided) unless lower_bound is supplied.

Interpretation

If $U_t < 0$ for some t , this is time-uniform evidence that forecaster 1 (p) is worse than forecaster 2 (q) on average — i.e. a rejection is evidence against p , not for it. More generally, $W_t > 0$ suggests p outperforms q ; $W_t < 0$ suggests q outperforms p .

Examples

```
p <- c(0.7, 0.6, 0.8, 0.65)
q <- c(0.5, 0.7, 0.6, 0.55)
y <- c(1, 1, 0, 1)
winkler_cs(p, q, y, alpha = 0.05)
```

winkler_etest

*E-process for Winkler scores (Proposition EC.4 + Theorem 3)***Description**

Tests whether the mean Winkler score $W_t \geq 0$ for all t . A rejection provides time-uniform evidence that forecaster 1 (p) is worse than forecaster 2 (q) under the base scoring rule.

Usage

```
winkler_etest(
  p,
  q,
  y,
  alpha = 0.05,
  base_score = log_score,
  v_opt = 10,
  clip_max = 1e+07
)
```

Arguments

<code>p</code>	Numeric vector in (0,1). Forecasts from model 1.
<code>q</code>	Numeric vector in (0,1). Forecasts from model 2.
<code>y</code>	Numeric vector containing only 0 and 1. Binary outcomes.
<code>alpha</code>	Numeric in (0,1). Significance level. Default: 0.05.
<code>base_score</code>	Function. Underlying scoring rule. Default: <code>log_score</code> .
<code>v_opt</code>	Numeric > 0. Default: 10.
<code>clip_max</code>	Numeric. Maximum e-process value before clipping. Default: 1e7.

Value

data.frame with columns `t`, `e`, `log_e`.

Rejection rule

Reject at level α when $e \geq 1 / \alpha$; this provides time-uniform evidence that p is worse than q .

Examples

```
p <- c(0.7, 0.6, 0.8, 0.65)
q <- c(0.5, 0.7, 0.6, 0.55)
y <- c(1, 1, 0, 1)
winkler_etest(p, q, y, alpha = 0.05)
```

winkler_score	<i>Winkler-normalized binary score</i>
---------------	--

Description

Normalises the score difference $S(p,y) - S(q,y)$ by the maximum possible score difference given the forecaster ordering, mapping the result to $(-\text{Inf}, 1]$ (Proposition EC.4, Choe & Ramdas 2024). Used to apply Theorems 2 & 3 to unbounded scoring rules on binary outcomes.

Usage

```
winkler_score(p, q, y, base_score = log_score, eps = 1e-08)
```

Arguments

<code>p</code>	Numeric vector in $(0,1)$. Forecasts from model 1.
<code>q</code>	Numeric vector in $(0,1)$. Forecasts from model 2.
<code>y</code>	Numeric vector containing only 0 and 1. Binary outcomes.
<code>base_score</code>	Function. The underlying scoring rule $S(p, y)$. Must accept two arguments: forecast probability and outcome. Default: <code>log_score</code> (with <code>eps</code> clipping).
<code>eps</code>	Numeric. Zero-protection for the normaliser denominator. Default: <code>1e-8</code> (matches Python comparecast convention).

Details

$$w(p, q, y) = \frac{S(p, y) - S(q, y)}{S(p, \mathbb{I}(p > q)) - S(q, \mathbb{I}(p > q))}$$

with the convention $0/0 := 0$.

The lower bound is problem-dependent (depends on how extreme p and q can be). For a two-sided CS via Corollary 2, the user must establish a finite lower bound analytically. If no finite lower bound can be guaranteed, use the one-sided (upper) CS only, as in the CR24 MLB experiments.

Value

Numeric vector. Winkler scores in $(-\text{Inf}, 1]$. Upper bound of 1 is tight: $w = 1$ when $y = 1(p > q)$.

When to use

Strictly limited to binary outcomes y in $\{0, 1\}$ and probability forecasts p, q in $(0, 1)$. Not applicable to QLIKE or other continuous-outcome scoring rules. See CR24 EC.7 for discussion.

For use in Theorems 2 & 3: upper bound = 1 implies $c/2 = 1$, so use $c = 2$ in all GE boundary and e-process calls.

Examples

```
p <- c(0.7, 0.6, 0.8, 0.65)
q <- c(0.5, 0.7, 0.6, 0.55)
y <- c(1, 1, 0, 1)
winkler_score(p, q, y)
```

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