

Package ‘pense’

September 13, 2020

Type Package

Title Penalized Elastic Net S/MM-Estimator of Regression

Version 2.0.1

Date 2020-09-13

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Encoding UTF-8

Biarch true

SystemRequirements C++11

URL <https://dakep.github.io/pense-rpkg>,
<https://github.com/dakep/pense-rpkg>

BugReports <https://github.com/dakep/pense-rpkg/issues>

Description Robust penalized (adaptive) elastic net S and M estimators for linear regression. The methods are proposed in Cohen Freue, G. V., Keppinger, D., Salibián-Barrera, M., and Smucler, E. (2019) <<https://projecteuclid.org/euclid.aoas/1574910036>>. The package implements the extensions and algorithms described in Keppinger, D. (2020) <[doi:10.14288/1.0392915](https://doi.org/10.14288/1.0392915)>.

Depends R (>= 3.5.0),
Matrix

Imports Rcpp,
methods,
parallel,
lifecycle (>= 0.2.0),
rlang (>= 0.4.0)

LinkingTo Rcpp,
RcppArmadillo (>= 0.9.600)

Suggests testthat (>= 2.1.0)

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NeedsCompilation yes

RoxygenNote 7.1.1

Roxygen list(markdown = TRUE, load = ``source``)

RdMacros lifecycle

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coef.pense_cvfit	<i>Extract Coefficient Estimates</i>
------------------	--------------------------------------

Description

Extract coefficients from a PENSE (or LS-EN) regularization path with hyper-parameters chosen by cross-validation.

Usage

```
## S3 method for class 'pense_cvfit'
coef(
  object,
  lambda = c("min", "se"),
  se_mult = 1,
  sparse = NULL,
  exact = deprecated(),
  correction = deprecated(),
  ...
)
```

Arguments

object	PENSE with cross-validated hyper-parameters to extract coefficients from.
lambda	either a string specifying which penalty level to use ("min" or "se") or a single numeric value of the penalty parameter. See details.
se_mult	If lambda = "se", the multiple of standard errors to tolerate.
sparse	should coefficients be returned as sparse or dense vectors? Defaults to the sparse argument supplied to pense_cv() . Can also be set to sparse = 'matrix', in which case a sparse matrix is returned instead of a sparse vector.
exact	deprecated. Always gives a warning if lambda is not part of the fitted sequence and coefficients are interpolated.
correction	defunct.
...	currently not used.

Details

If lambda = "se" and object contains fitted estimates for every penalization level in the sequence, extract the coefficients of the most parsimonious model with prediction performance statistically indistinguishable from the best model. This is determined to be the model with prediction performance within $se_mult * cv_se$ from the best model.

Value

either a numeric vector or a sparse vector of type [dsparseVector](#) of size $p + 1$, depending on the sparse argument. Note: prior to version 2.0.0 sparse coefficients were returned as sparse matrix of type *dgCMatrix*. To get a sparse matrix, use sparse = 'matrix'.

See Also

Other functions for extracting components: [coef.pense_fit\(\)](#), [predict.pense_cvfit\(\)](#), [predict.pense_fit\(\)](#), [residuals.pense_cvfit\(\)](#), [residuals.pense_fit\(\)](#)

Examples

```
# Compute the PENSE regularization path for Freeny's revenue data
# (see ?freeny)
data(freeny)
x <- as.matrix(freeny[, 2:5])

regpath <- pense(x, freeny$y, alpha = 0.5)
```

```

plot(regpath)

# Extract the coefficients at a certain penalization level
coef(regpath, lambda = regpath$lambda[40])

# What penalization level leads to good prediction performance?
cv_results <- pense_cv(x, freeny$y, alpha = 0.5, cv_repl = 2,
                      cv_k = 4)
plot(cv_results, se_mult = 1)

# Extract the coefficients at the penalization level with
# smallest prediction error ...
coef(cv_results)
# ... or at the penalization level with prediction error
# statistically indistinguishable from the minimum.
coef(cv_results, lambda = 'se')

```

coef.pense_fit	<i>Extract Coefficient Estimates</i>
----------------	--------------------------------------

Description

Extract coefficients from a PENSE (or LS-EN) regularization path fitted by [pense\(\)](#) or [elnet\(\)](#).

Usage

```

## S3 method for class 'pense_fit'
coef(
  object,
  lambda,
  sparse = NULL,
  exact = deprecated(),
  correction = deprecated(),
  ...
)

```

Arguments

object	PENSE regularization path to extract coefficients from.
lambda	a single value of the penalty parameter.
sparse	should coefficients be returned as sparse or dense vectors? Defaults to the sparse argument supplied to pense() . Can also be set to sparse = 'matrix', in which case a sparse matrix is returned instead of a sparse vector.
exact	defunct Always gives a warning if lambda is not part of the fitted sequence and hence coefficients are interpolated.
correction	defunct.
...	currently not used.

Value

either a numeric vector or a sparse vector of type [dsparseVector](#) of size $p + 1$, depending on the sparse argument. Note: prior to version 2.0.0 sparse coefficients were returned as sparse matrix of type *dgCMatrix*. To get a sparse matrix, use sparse = 'matrix'.

See Also

[coef.pense_cvfit\(\)](#) for extracting coefficients from a PENSE fit with hyper-parameters chosen by cross-validation

Other functions for extracting components: [coef.pense_cvfit\(\)](#), [predict.pense_cvfit\(\)](#), [predict.pense_fit\(\)](#), [residuals.pense_cvfit\(\)](#), [residuals.pense_fit\(\)](#)

Examples

```
# Compute the PENSE regularization path for Freeny's revenue data
# (see ?freeny)
data(freeny)
x <- as.matrix(freeny[, 2:5])

regpath <- pense(x, freeny$y, alpha = 0.5)
plot(regpath)

# Extract the coefficients at a certain penalization level
coef(regpath, lambda = regpath$lambda[40])

# What penalization level leads to good prediction performance?
cv_results <- pense_cv(x, freeny$y, alpha = 0.5, cv_repl = 2,
                      cv_k = 4)
plot(cv_results, se_mult = 1)

# Extract the coefficients at the penalization level with
# smallest prediction error ...
coef(cv_results)
# ... or at the penalization level with prediction error
# statistically indistinguishable from the minimum.
coef(cv_results, lambda = 'se')
```

consistency_const	<i>Get the Constant for Consistency for the M-Scale</i>
-------------------	---------------------------------------------------------

Description

Get the Constant for Consistency for the M-Scale

Usage

```
consistency_const(delta, rho)
```

Arguments

delta	desired breakdown point (between 0 and 0.5)
rho	the name of the chosen ρ function.

Value

consistency constant

See Also

Other miscellaneous functions: [rho_function\(\)](#)

deprecated_en_options *Deprecated Options for the EN Algorithms*

Description

Deprecated

Usage

```
en_options_aug_lars(use_gram = c("auto", "yes", "no"), eps = 1e-12)
```

```
en_options_dal(
  maxit = 100,
  eps = 1e-08,
  eta_mult = 2,
  eta_start_numerator = 0.01,
  eta_start,
  preconditioner = c("approx", "none", "diagonal"),
  verbosity = 0
)
```

Arguments

use_gram	ignored. Should the Gram matrix be pre-computed.
eps	ignored. Numeric tolerance for convergence.
maxit	maximum number of iterations allowed.
eta_mult	multiplier to increase eta at each iteration.
eta_start_numerator	if eta_start is missing, it is defined by $\text{eta_start} = \text{eta_start_numerator} / \text{lambda}$.
eta_start	ignored. The start value for eta.
preconditioner	ignored. Preconditioner for the numerical solver. If none, a standard solver will be used, otherwise the faster preconditioned conjugate gradient is used.
verbosity	ignored.

Functions

- en_options_aug_lars: Superseded by [en_lars_options\(\)](#).
- en_options_dal: Superseded by [en_dal_options\(\)](#)

See Also

Other deprecated functions: [enpy\(\)](#), [initest_options\(\)](#), [mstep_options\(\)](#), [pense_options\(\)](#), [pensem\(\)](#)

elnet

*Compute the Least Squares (Adaptive) Elastic Net Regularization Path***Description**

Compute least squares EN estimates for linear regression with optional observation weights and penalty loadings.

Usage

```
elnet(
  x,
  y,
  alpha,
  nlambdas = 100,
  lambda_min_ratio,
  lambda,
  penalty_loadings,
  weights,
  intercept = TRUE,
  en_algorithm_opts,
  sparse = FALSE,
  eps = 1e-06,
  standardize = TRUE,
  correction = deprecated(),
  xtest = deprecated(),
  options = deprecated()
)
```

Arguments

<code>x</code>	<code>n</code> by <code>p</code> matrix of numeric predictors.
<code>y</code>	vector of response values of length <code>n</code> .
<code>alpha</code>	elastic net penalty mixing parameter with $0 \leq \alpha \leq 1$. <code>alpha = 1</code> is the LASSO penalty, and <code>alpha = 0</code> the Ridge penalty.
<code>nlambdas</code>	number of penalization levels.
<code>lambda_min_ratio</code>	Smallest value of the penalization level as a fraction of the largest level (i.e., the smallest value for which all coefficients are zero). The default depends on the sample size relative to the number of variables and <code>alpha</code> . If more observations than variables are available, the default is $1e-3 * \alpha$, otherwise $1e-2 * \alpha$.
<code>lambda</code>	optional user-supplied sequence of penalization levels. If given and not <code>NULL</code> , <code>nlambdas</code> and <code>lambda_min_ratio</code> are ignored.
<code>penalty_loadings</code>	a vector of positive penalty loadings (a.k.a. weights) for different penalization of each coefficient.
<code>weights</code>	a vector of positive observation weights.
<code>intercept</code>	include an intercept in the model.

<code>en_algorithm_opts</code>	options for the EN algorithm. See en_algorithm_options for details.
<code>sparse</code>	use sparse coefficient vectors.
<code>eps</code>	numerical tolerance.
<code>standardize</code>	standardize variables to have unit variance. Coefficients are always returned in original scale.
<code>correction</code>	defunct. Correction for EN estimates is not supported anymore.
<code>xtest</code>	deprecated. Instead, extract coefficients with coef.pense_fit() and compute predictions manually.
<code>options</code>	deprecated. Use <code>en_algorithm_opts</code> instead.

Details

The elastic net estimator for the linear regression model solves the optimization problem

$$\operatorname{argmin}_{\mu, \beta} (1/2n) \sum_i w_i (y_i - \mu - x'_i \beta)^2 + \lambda \sum_j 0.5(1 - \alpha) \beta_j^2 + \alpha l_j |\beta_j|$$

with observation weights w_i and penalty loadings l_j .

Value

a list-like object with the following items

`lambda` the sequence of penalization parameters.

`estimates` a list of estimates. Each estimate contains the following information:

`intercept` intercept estimate.

`beta` beta (slope) estimate.

`lambda` penalization level at which the estimate is computed.

`alpha` *alpha* hyper-parameter at which the estimate is computed.

`statuscode` if > 0 the algorithm experienced issues when computing the estimate.

`status` optional status message from the algorithm.

`call` the original call.

`predictions` if `xtest` was given, a matrix of predicted values. Each column corresponds to the predictions from the estimate at the `lambda` value at the same index.

See Also

[pense\(\)](#) for an S-estimate of regression with elastic net penalty.

[coef.pense_fit\(\)](#) for extracting coefficient estimates.

[plot.pense_fit\(\)](#) for plotting the regularization path.

Other functions for computing non-robust estimates: [elnet_cv\(\)](#)

Examples

```
# Compute the LS-EN regularization path for Freeny's revenue data
# (see ?freeny)
data(freeny)
x <- as.matrix(freeny[, 2:5])

regpath <- elnet(x, freeny$y, alpha = 0.75)
plot(regpath)

# Extract the coefficients at a certain penalization level
coef(regpath, lambda = regpath$lambda[5])

# What penalization level leads to good prediction performance?
cv_results <- elnet_cv(x, freeny$y, alpha = 0.75, cv_repl = 10,
                      cv_k = 4, cv_measure = 'tau')
plot(cv_results, se_mult = 1)
plot(cv_results, se_mult = 1, what = 'coef.path')

# Extract the coefficients at the penalization level with
# smallest prediction error ...
coef(cv_results)
# ... or at the penalization level with prediction error
# statistically indistinguishable from the minimum.
coef(cv_results, lambda = 'se')
```

elnet_cv

Cross-validation for Least-Squares (Adaptive) Elastic Net Estimates

Description

Perform (repeated) K-fold cross-validation for [elnet\(\)](#).

Usage

```
elnet_cv(
  x,
  y,
  lambda,
  cv_k,
  cv_repl = 1,
  cv_metric = c("rmspe", "tau_size", "mape"),
  fit_all = TRUE,
  cl = NULL,
  ncores = deprecated(),
  ...
)
```

Arguments

x	n by p matrix of numeric predictors.
y	vector of response values of length n.

<code>lambda</code>	optional user-supplied sequence of penalization levels. If given and not NULL, <code>nlambda</code> and <code>lambda_min_ratio</code> are ignored.
<code>cv_k</code>	number of folds per cross-validation.
<code>cv_repl</code>	number of cross-validation replications.
<code>cv_metric</code>	either a string specifying the performance metric to use, or a function to evaluate prediction errors in a single CV replication. If a function, it is called with a single numeric vector of prediction errors and must return a scalar number.
<code>fit_all</code>	If TRUE, fit the model for all penalization levels. Otherwise, only at penalization level with smallest average CV performance.
<code>cl</code>	a parallel cluster. Can only be used if <code>ncores = 1</code> , because multi-threading can not be used in parallel R sessions on the same host.
<code>ncores</code>	deprecated and not used anymore.
<code>...</code>	Arguments passed on to elnet
<code>alpha</code>	elastic net penalty mixing parameter with $0 \leq \alpha \leq 1$. <code>alpha = 1</code> is the LASSO penalty, and <code>alpha = 0</code> the Ridge penalty.
<code>nlambda</code>	number of penalization levels.
<code>lambda_min_ratio</code>	Smallest value of the penalization level as a fraction of the largest level (i.e., the smallest value for which all coefficients are zero). The default depends on the sample size relative to the number of variables and <code>alpha</code> . If more observations than variables are available, the default is $1e-3 * \alpha$, otherwise $1e-2 * \alpha$.
<code>penalty_loadings</code>	a vector of positive penalty loadings (a.k.a. weights) for different penalization of each coefficient.
<code>standardize</code>	standardize variables to have unit variance. Coefficients are always returned in original scale.
<code>weights</code>	a vector of positive observation weights.
<code>intercept</code>	include an intercept in the model.
<code>sparse</code>	use sparse coefficient vectors.
<code>en_algorithm_opts</code>	options for the EN algorithm. See en_algorithm_options for details.
<code>eps</code>	numerical tolerance.
<code>xtest</code>	deprecated. Instead, extract coefficients with coef.pense_fit() and compute predictions manually.
<code>options</code>	deprecated. Use <code>en_algorithm_opts</code> instead.
<code>correction</code>	defunct. Correction for EN estimates is not supported anymore.

Details

The built-in CV metrics are

"rmspe" Root mean squared prediction error (default).

"mape" Median absolute prediction error.

"tau_size" τ -size of the prediction error, computed by [tau_size\(\)](#).

Value

a list with components:

`lambda` the sequence of penalization levels.

`cvres` data frame of average cross-validated performance.

`cv_replications` matrix of cross-validated performance metrics, one column per replication. Rows are in the same order as in `cvres`.

`call` the original call.

`estimates` the estimates fitted on the full data. Same format as returned by `elnet()`.

See Also

`elnet()` for computing the LS-EN regularization path without cross-validation.

`pense_cv()` for cross-validation of S-estimates of regression with elastic net penalty.

`coef.pense_cvfit()` for extracting coefficient estimates.

`plot.pense_cvfit()` for plotting the CV performance or the regularization path.

Other functions for computing non-robust estimates: `elnet()`

Examples

```
# Compute the LS-EN regularization path for Freeny's revenue data
# (see ?freeny)
data(freeny)
x <- as.matrix(freeny[, 2:5])

regpath <- elnet(x, freeny$y, alpha = 0.75)
plot(regpath)

# Extract the coefficients at a certain penalization level
coef(regpath, lambda = regpath$lambda[5])

# What penalization level leads to good prediction performance?
cv_results <- elnet_cv(x, freeny$y, alpha = 0.75, cv_repl = 10,
                      cv_k = 4, cv_measure = 'tau')
plot(cv_results, se_mult = 1)
plot(cv_results, se_mult = 1, what = 'coef.path')

# Extract the coefficients at the penalization level with
# smallest prediction error ...
coef(cv_results)
# ... or at the penalization level with prediction error
# statistically indistinguishable from the minimum.
coef(cv_results, lambda = 'se')
```

enpy

Deprecated: ENPY Initial Estimates for EN S-Estimators

Description

Deprecated Superseded by `enpy_initial_estimates()`.

Usage

```
enpy(x, y, alpha, lambda, delta, cc, options, en_options)
```

Arguments

x	data matrix with predictors.
y	response vector.
alpha, lambda	EN penalty parameters (NOT adjusted for the number of observations in x).
delta	desired breakdown point of the resulting estimator.
cc	tuning constant for the S-estimator. Default is to chosen based on the breakdown point delta. Should never have to be changed.
options	ignored. Additional options for the initial estimator.
en_options	ignored. Additional options for the EN algorithm.

Value

coeff	A numeric matrix with one initial coefficient per column
objF	A vector of values of the objective function for the respective coefficient

See Also

Other deprecated functions: [deprecated_en_options](#), [initest_options\(\)](#), [mstep_options\(\)](#), [pense_options\(\)](#), [pensem\(\)](#)

enpy_initial_estimates

ENPY Initial Estimates for EN S-Estimators

Description

Compute initial estimates for the EN S-estimator using the EN-PY procedure.

Usage

```
enpy_initial_estimates(
  x,
  y,
  alpha,
  lambda,
  bdp = 0.25,
  cc,
  intercept = TRUE,
  penalty_loadings,
  enpy_opts = enpy_options(),
  mscale_opts = mscale_algorithm_options(),
  eps = 1e-06,
  sparse = FALSE,
  ncores = 1L
)
```

Arguments

x	n by p matrix of numeric predictors.
y	vector of response values of length n.
alpha	elastic net penalty mixing parameter with $0 \leq \alpha \leq 1$. alpha = 1 is the LASSO penalty, and alpha = 0 the Ridge penalty.
lambda	a vector of positive values of penalization levels.
bdp	desired breakdown point of the estimator, between 0 and 0.5.
cc	cutoff value for the bisquare rho function. By default, chosen to yield a consistent estimate for the Normal distribution.
intercept	include an intercept in the model.
penalty_loadings	a vector of positive penalty loadings (a.k.a. weights) for different penalization of each coefficient. Only allowed for alpha > 0.
enpy_opts	options for the EN-PY algorithm, created with the enpy_options() function.
mscale_opts	options for the M-scale estimation. See mscale_algorithm_options() for details.
eps	numerical tolerance.
sparse	use sparse coefficient vectors.
ncores	number of CPU cores to use in parallel. By default, only one CPU core is used. May not be supported on your platform, in which case a warning is given.

Details

If these manually computed initial estimates are intended as starting points for [pense\(\)](#), they are by default *shared* for all penalization levels. To restrict the use of the initial estimates to the penalty level they were computed for, use [as_starting_point\(..., specific = TRUE\)](#). See [as_starting_point\(\)](#) for details.

References

Cohen Freue, G.V.; Kepplinger, D.; Salibián-Barrera, M.; Smucler, E. Robust elastic net estimators for variable selection and identification of proteomic biomarkers. *Ann. Appl. Stat.* **13** (2019), no. 4, 2065–2090 doi: [10.1214/19AOAS1269](#)

See Also

Other functions for initial estimates: [prinsens\(\)](#), [starting_point\(\)](#)

enpy_options

Options for the ENPY Algorithm

Description

Additional control options for the elastic net Peña-Yohai procedure.

Usage

```

enpy_options(
  max_it = 10,
  keep_psc_proportion = 0.5,
  en_algorithm_opts,
  keep_residuals_measure = c("threshold", "proportion"),
  keep_residuals_proportion = 0.5,
  keep_residuals_threshold = 2,
  retain_best_factor = 2,
  retain_max = 500
)

```

Arguments

<code>max_it</code>	maximum number of EN-PY iterations.
<code>keep_psc_proportion</code>	how many observations should to keep based on the Principal Sensitivity Components.
<code>en_algorithm_opts</code>	options for the LS-EN algorithm. See en_algorithm_options for details.
<code>keep_residuals_measure</code>	how to determine what observations to keep, based on their residuals. If <code>proportion</code> , a fixed number of observations is kept. If <code>threshold</code> , only observations with residuals below the threshold are kept.
<code>keep_residuals_proportion</code>	proportion of observations to kept based on their residuals.
<code>keep_residuals_threshold</code>	only observations with (standardized) residuals less than this threshold are kept.
<code>retain_best_factor</code>	only keep candidates that are within this factor of the best candidate. If ≤ 1 , only keep candidates from the last iteration.
<code>retain_max</code>	maximum number of candidates, i.e., only the best <code>retain_max</code> candidates are retained.

Details

The EN-PY procedure for computing initial estimates iteratively cleans the data of observations with possibly outlying residual or high leverage. Least-squares elastic net (LS-EN) estimates are computed on the possibly clean subsets. At each iteration, the Principal Sensitivity Components are computed to remove observations with potentially high leverage. Among all the LS-EN estimates, the estimate with smallest M-scale of the residuals is selected. Observations with largest residual for the selected estimate are removed and the next iteration is started.

Value

options for the ENPY algorithm.

en_admm_options	<i>Use the ADMM Elastic Net Algorithm</i>
-----------------	-------------------------------------------

Description

Use the ADMM Elastic Net Algorithm

Usage

```
en_admm_options(max_it = 1000, step_size, acceleration = 1)
```

Arguments

max_it	maximum number of iterations.
step_size	step size for the algorithm.
acceleration	acceleration factor for linearized ADMM.

Value

options for the ADMM EN algorithm.

See Also

Other EN algorithms: [en_dal_options\(\)](#), [en_lars_options\(\)](#)

en_algorithm_options	<i>Control the Algorithm to Compute (Weighted) Least-Squares Elastic Net Estimates</i>
----------------------	----------------------------------------------------------------------------------------

Description

The package supports different algorithms to compute the EN estimate for weighted LS loss functions. Each algorithm has certain characteristics that make it useful for some problems. To select a specific algorithm and adjust the options, use any of the `en_***_options` functions.

Details

- [en_lars_options\(\)](#): Use the tuning-free LARS algorithm. This computes *exact* (up to numerical errors) solutions to the EN-LS problem. It is not iterative and therefore can not benefit from approximate solutions, but in turn guarantees that a solution will be found.
- [en_admm_options\(\)](#): Use an iterative ADMM-type algorithm which needs $O(np)$ operations per iteration and converges sub-linearly.
- [en_dal_options\(\)](#): Use the iterative Dual Augmented Lagrangian (DAL) method. DAL needs $O(n^3p^2)$ operations per iteration, but converges exponentially.

en_dal_options	<i>Use the DAL Elastic Net Algorithm</i>
----------------	------------------------------------------

Description

Use the DAL Elastic Net Algorithm

Usage

```
en_dal_options(
  max_it = 100,
  max_inner_it = 100,
  eta_multiplier = 2,
  eta_start_conservative = 0.01,
  eta_start_aggressive = 1,
  lambda_relchange_aggressive = 0.25
)
```

Arguments

max_it	maximum number of (outer) iterations.
max_inner_it	maximum number of (inner) iterations in each outer iteration.
eta_multiplier	multiplier for the barrier parameter. In each iteration, the barrier must be more restrictive (i.e., the multiplier must be > 1).
eta_start_conservative	conservative initial barrier parameter. This is used if the previous penalty is undefined or too far away.
eta_start_aggressive	aggressive initial barrier parameter. This is used if the previous penalty is close.
lambda_relchange_aggressive	how close must the lambda parameter from the previous penalty term be to use an aggressive initial barrier parameter (i.e., what constitutes "too far").

Value

options for the DAL EN algorithm.

See Also

Other EN algorithms: [en_admm_options\(\)](#), [en_lars_options\(\)](#)

en_lars_options	<i>Use the LARS Elastic Net Algorithm</i>
-----------------	-------------------------------------------

Description

Use the LARS Elastic Net Algorithm

Usage

```
en_lars_options()
```

See Also

Other EN algorithms: [en_admm_options\(\)](#), [en_dal_options\(\)](#)

initest_options	<i>Deprecated Options for Initial Estimates</i>
-----------------	-------------------------------------------------

Description

Deprecated Superseded by [enpy_options\(\)](#).

Usage

```
initest_options(
  keep_solutions = 5,
  psc_method = c("exact", "rr"),
  maxit = 10,
  maxit_pense_refinement = 5,
  eps = 1e-06,
  psc_keep = 0.5,
  resid_keep_method = c("proportion", "threshold"),
  resid_keep_prop = 0.6,
  resid_keep_thresh = 2,
  mscale_eps = 1e-08,
  mscale_maxit = 200
)
```

Arguments

keep_solutions	how many initial estimates should be kept to perform full PENSE iterations?
psc_method	The method to use for computing the principal sensitivity components. See details for the possible choices.
maxit	maximum number of refinement iterations.
maxit_pense_refinement	ignored. Maximum number of PENSE iterations to refine initial estimator.
eps	ignored. Numeric tolerance for convergence.
psc_keep	proportion of observations to keep based on the PSC scores.

resid_keep_method

How to clean the data based on large residuals. If "proportion", observations with the smallest resid_keep_prop residuals will be retained. If "threshold", all observations with scaled residuals smaller than the threshold resid_keep_thresh will be retained.

resid_keep_prop, resid_keep_thresh

proportion or threshold for observations to keep based on their residual.

mscale_eps, mscale_maxit

ignored. Maximum number of iterations and numeric tolerance for the M-scale.

See Also

Other deprecated functions: [deprecated_en_options](#), [enpy\(\)](#), [mstep_options\(\)](#), [pense_options\(\)](#), [pensem\(\)](#)

mloc	<i>Compute the M-estimate of Location</i>
------	-------------------------------------------

Description

Compute the M-estimate of location using an auxiliary estimate of the scale.

Usage

```
mloc(x, scale, rho, cc, opts = mscale_algorithm_options())
```

Arguments

x	numeric values. Missing values are verbosely ignored.
scale	scale of the x values. If omitted, uses the mad() .
rho	the ρ function to use. See rho_function() for available functions.
cc	value of the tuning constant for the chosen ρ function. By default, chosen to achieve 95% efficiency under the Normal distribution.
opts	a list of options for the M-estimating algorithm, see mscale_algorithm_options() for details.

Value

a single numeric value, the M-estimate of location.

See Also

Other functions to compute robust estimates of location and scale: [mlocscale\(\)](#), [mscale\(\)](#), [tau_size\(\)](#)

mlocscale

Compute the M-estimate of Location and Scale

Description

Simultaneous estimation of the location and scale by means of M-estimates.

Usage

```
mlocscale(
  x,
  bdp = 0.25,
  scale_cc = consistency_const(bdp, "bisquare"),
  location_rho,
  location_cc,
  opts = mscale_algorithm_options()
)
```

Arguments

x	numeric values. Missing values are verbosely ignored.
bdp	desired breakdown point (between 0 and 0.5).
scale_cc	cutoff value for the bisquare ρ function for computing the scale estimate. By default, chosen to yield a consistent estimate for normally distributed values.
location_rho, location_cc	ρ function and cutoff value for computing the location estimate. See rho_function() for a list of available ρ functions.
opts	a list of options for the M-estimating equation, see mscale_algorithm_options() for details.

Value

a vector with 2 elements, the M-estimate of location and the M-scale estimate.

See Also

Other functions to compute robust estimates of location and scale: [mloc\(\)](#), [mscale\(\)](#), [tau_size\(\)](#)

mm_algorithm_options

MM-Algorithm to Compute Penalized Elastic Net S- and M-Estimates

Description

Additional options for the MM algorithm to compute EN S- and M-estimates.

Usage

```
mm_algorithm_options(
  max_it = 500,
  tightening = c("adaptive", "exponential", "none"),
  tightening_steps = 10,
  en_algorithm_opts
)
```

Arguments

`max_it` maximum number of iterations.

`tightening` how to make inner iterations more precise as the algorithm approaches a local minimum.

`tightening_steps` for *adaptive* tightening strategy, how often to tighten until the desired tolerance is attained.

`en_algorithm_opts` options for the inner LS-EN algorithm. See [en_algorithm_options](#) for details.

Value

options for the MM algorithm.

mscale	<i>Compute the M-Scale of Centered Values</i>
--------	-----------------------------------------------

Description

Compute the M-scale without centering the values.

Usage

```
mscale(
  x,
  bdp = 0.25,
  cc = consistency_const(bdp, "bisquare"),
  opts = mscale_algorithm_options(),
  delta = deprecated(),
  rho = deprecated(),
  eps = deprecated(),
  maxit = deprecated()
)
```

Arguments

`x` numeric values. Missing values are verbosely ignored.

`bdp` desired breakdown point (between 0 and 0.5).

`cc` cutoff value for the bisquare rho function. By default, chosen to yield a consistent estimate for the Normal distribution.

opts	a list of options for the M-scale estimation algorithm, see mscale_algorithm_options() for details.
delta	deprecated. Use bpd instead.
rho, eps, maxit	deprecated. Instead set control options for the algorithm with the opts arguments.

Value

the M-estimate of scale.

See Also

Other functions to compute robust estimates of location and scale: [mlocscale\(\)](#), [mloc\(\)](#), [tau_size\(\)](#)

mscale_algorithm_options

Options for the M-scale Estimation Algorithm

Description

Options for the M-scale Estimation Algorithm

Usage

```
mscale_algorithm_options(max_it = 200, eps = 1e-08)
```

Arguments

max_it	maximum number of iterations.
eps	numerical tolerance to check for convergence.

Value

options for the M-scale estimation algorithm.

mstep_options

Deprecated Additional Options for the Penalized EN MM-estimator

Description

Deprecated Superseded by [mm_algorithm_options\(\)](#) and options supplied directly to [pense\(\)](#).

Usage

```
mstep_options(
  cc = 3.44,
  maxit = 1000,
  eps = 1e-06,
  adjust_bdp = FALSE,
  verbosity = 0,
  en_correction = TRUE
)
```

Arguments

<code>cc</code>	ignored. Tuning constant for the M-estimator.
<code>maxit</code>	maximum number of iterations allowed.
<code>eps</code>	ignored. Numeric tolerance for convergence.
<code>adjust_bdp</code>	ignored. Should the breakdown point be adjusted based on the effective degrees of freedom?
<code>verbosity</code>	ignored. Verbosity of the algorithm.
<code>en_correction</code>	ignored. Should the corrected EN estimator be used to choose the optimal lambda with CV. If TRUE, as by default, the estimator is "bias corrected".

See Also

Other deprecated functions: [deprecated_en_options](#), [enpy\(\)](#), [initest_options\(\)](#), [pense_options\(\)](#), [pensem\(\)](#)

pense

Compute (Adaptive) Elastic Net S-Estimates of Regression

Description

Compute elastic net S-estimates (PENSE estimates) along a grid of penalization levels with optional penalty loadings for adaptive elastic net.

Usage

```
pense(
  x,
  y,
  alpha,
  nlambda = 50,
  nlambda_enpy = 10,
  lambda,
  lambda_min_ratio,
  enpy_lambda,
  penalty_loadings,
  intercept = TRUE,
  bdp = 0.25,
  add_zero_based = TRUE,
  enpy_specific = FALSE,
  other_starts,
  eps = 1e-06,
  explore_solutions = 10,
  explore_tol = 0.1,
  max_solutions = 10,
  comparison_tol = sqrt(eps),
  sparse = FALSE,
  ncores = 1,
  standardize = TRUE,
  algorithm_opts = mm_algorithm_options(),
```

```

mscale_opts = mscale_algorithm_options(),
enpy_opts = enpy_options(),
cv_k = deprecated(),
cv_objective = deprecated(),
...
)

```

Arguments

<code>x</code>	<code>n</code> by <code>p</code> matrix of numeric predictors.
<code>y</code>	vector of response values of length <code>n</code> .
<code>alpha</code>	elastic net penalty mixing parameter with $0 \leq \alpha \leq 1$. <code>alpha = 1</code> is the LASSO penalty, and <code>alpha = 0</code> the Ridge penalty.
<code>nlambda</code>	number of penalization levels.
<code>nlambda_enpy</code>	number of penalization levels where the EN-PY initial estimate is computed.
<code>lambda</code>	optional user-supplied sequence of penalization levels. If given and not <code>NULL</code> , <code>nlambda</code> and <code>lambda_min_ratio</code> are ignored.
<code>lambda_min_ratio</code>	Smallest value of the penalization level as a fraction of the largest level (i.e., the smallest value for which all coefficients are zero). The default depends on the sample size relative to the number of variables and <code>alpha</code> . If more observations than variables are available, the default is $1e-3 * \alpha$, otherwise $1e-2 * \alpha$.
<code>enpy_lambda</code>	optional user-supplied sequence of penalization levels at which EN-PY initial estimates are computed. If given and not <code>NULL</code> , <code>nlambda_enpy</code> is ignored.
<code>penalty_loadings</code>	a vector of positive penalty loadings (a.k.a. weights) for different penalization of each coefficient. Only allowed for <code>alpha > 0</code> .
<code>intercept</code>	include an intercept in the model.
<code>bdp</code>	desired breakdown point of the estimator, between 0 and 0.5.
<code>add_zero_based</code>	also consider the 0-based regularization path. See details for a description.
<code>enpy_specific</code>	use the EN-PY initial estimates only at the penalization level they are computed for. See details for a description.
<code>other_starts</code>	a list of other starting points, created by starting_point() . If the output of enpy_initial_estimates() is given, the starting points will be <i>shared</i> among all penalization levels. Note that if a the starting point is <i>specific</i> to a penalization level, this penalization level is added to the grid of penalization levels (either the manually specified grid in <code>lambda</code> or the automatically generated grid of size <code>nlambda</code>). If <code>standardize = TRUE</code> , the starting points are also scaled.
<code>eps</code>	numerical tolerance.
<code>explore_solutions</code>	number of solutions to compute up to the desired precision <code>eps</code> .
<code>explore_tol</code>	numerical tolerance for exploring possible solutions. Should be (much) looser than <code>eps</code> to be useful.
<code>max_solutions</code>	only retain up to <code>max_solutions</code> unique solutions per penalization level.

<code>comparison_tol</code>	numeric tolerance to determine if two solutions are equal. The comparison is first done on the absolute difference in the value of the objective function at the solution. If this is less than <code>comparison_tol</code> , two solutions are deemed equal. If the squared difference of the intercepts is less than <code>comparison_tol</code> and the squared L_2 norm of the difference vector is less than <code>comparison_tol</code> .
<code>sparse</code>	use sparse coefficient vectors.
<code>ncores</code>	number of CPU cores to use in parallel. By default, only one CPU core is used. May not be supported on your platform, in which case a warning is given.
<code>standardize</code>	logical flag to standardize the x variables prior to fitting the PENSE estimates. Coefficients are always returned on the original scale. This can fail for variables with a large proportion of a single value (e.g., zero-inflated data). In this case, either compute with <code>standardize = FALSE</code> or standardize the data manually.
<code>algorithm_opts</code>	options for the MM algorithm to compute the estimates. See mm_algorithm_options() for details.
<code>mscale_opts</code>	options for the M-scale estimation. See mscale_algorithm_options() for details.
<code>enpy_opts</code>	options for the ENPY initial estimates, created with the enpy_options() function. See enpy_initial_estimates() for details.
<code>cv_k, cv_objective</code>	deprecated and ignored. See pense_cv() for estimating prediction performance via cross-validation.
<code>...</code>	ignored. See the section on deprecated parameters below.

Value

a list-like object with the following items

`lambda` the sequence of penalization levels.

`estimates` a list of estimates. Each estimate contains the following information:

`intercept` intercept estimate.

`beta` beta (slope) estimate.

`lambda` penalization level at which the estimate is computed.

`alpha` *alpha* hyper-parameter at which the estimate is computed.

`objf_value` value of the objective function at the solution.

`statuscode` if > 0 the algorithm experienced issues when computing the estimate.

`status` optional status message from the algorithm.

`call` the original call.

Strategies for Using Starting Points

The function supports several different strategies to compute, and use the provided starting points for optimizing the PENSE objective function.

Starting points are computed internally but can also be supplied via `other_starts`. By default, starting points are computed internally by the EN-PY procedure for penalization levels supplied in `enpy_lambda` (or the automatically generated grid of length `nlambda_enpy`). By default, starting points computed by the EN-PY procedure are *shared* for all penalization levels in `lambda` (or the automatically generated grid of length `nlambda`). If the starting points should be *specific* to the penalization level the starting points' penalization level, set the `enpy_specific` argument to `TRUE`.

In addition to EN-PY initial estimates, the algorithm can also use the "0-based" strategy if `add_zero_based = TRUE` (by default). Here, the 0-vector is used to start the optimization at the largest penalization level in `lambda`. At subsequent penalization levels, the solution at the previous penalization level is also used as starting point.

At every penalization level, all starting points are explored using the loose numerical tolerance `explore_tol`. Only the best `explore_solutions` are computed to the stringent numerical tolerance `eps`. Finally, only the best `max_solutions` are retained and carried forward as starting points for the subsequent penalization level.

Deprecated Arguments

Starting with version 2.0.0, cross-validation is performed by separate function `pense_cv()`. Arguments related cross-validation cause an error when supplied to `pense()`. Furthermore, the following arguments are deprecated as of version 2.0.0: `initial`, `warm_reset`, `cl`, `options`, `init_options`, `en_options`. If `pense()` is called with any of these arguments, warnings detail how to replace them.

See Also

`pense_cv()` for selecting hyper-parameters via cross-validation.

`coef.pense_fit()` for extracting coefficient estimates.

`plot.pense_fit()` for plotting the regularization path.

Other functions to compute robust estimates: `regmest()`

Examples

```
# Compute the PENSE regularization path for Freeny's revenue data
# (see ?freeny)
data(freeny)
x <- as.matrix(freeny[, 2:5])

regpath <- pense(x, freeny$y, alpha = 0.5)
plot(regpath)

# Extract the coefficients at a certain penalization level
coef(regpath, lambda = regpath$lambda[40])

# What penalization level leads to good prediction performance?
cv_results <- pense_cv(x, freeny$y, alpha = 0.5, cv_repl = 2,
                      cv_k = 4)
plot(cv_results, se_mult = 1)

# Extract the coefficients at the penalization level with
# smallest prediction error ...
coef(cv_results)
# ... or at the penalization level with prediction error
# statistically indistinguishable from the minimum.
coef(cv_results, lambda = 'se')
```

pensem	<i>Deprecated Alias of pensem_cv</i>
--------	--------------------------------------

Description

[pensem\(\)](#) is a deprecated alias for [pensem_cv\(\)](#).

Usage

```
pensem(x, ...)
```

Arguments

x	either a numeric matrix of predictor values, or a cross-validated PENSE fit from pense_cv() .
...	ignored. See the section on deprecated parameters below.

See Also

Other deprecated functions: [deprecated_en_options](#), [enpy\(\)](#), [initest_options\(\)](#), [mstep_options\(\)](#), [pense_options\(\)](#)

pensem_cv	<i>Compute Penalized Elastic Net M-Estimates from PENSE</i>
-----------	-------------------------------------------------------------

Description

This is a convenience wrapper around [pense_cv\(\)](#) and [regmest_cv\(\)](#), for the common use-case of computing a highly-robust S-estimate followed by a more efficient M-estimate using the scale of the residuals from the S-estimate.

Usage

```
pensem_cv(x, ...)

## Default S3 method:
pensem_cv(
  x,
  y,
  alpha = 0.5,
  nlambdas = 50,
  lambda_min_ratio,
  lambda_m,
  lambda_s,
  standardize = TRUE,
  penalty_loadings,
  intercept = TRUE,
  bdp = 0.25,
  ncores = 1,
```

```

    sparse = FALSE,
    eps = 1e-06,
    cc = 4.7,
    cv_k = 5,
    cv_repl = 1,
    cl = NULL,
    cv_metric = c("tau_size", "mape", "rmspe"),
    add_zero_based = TRUE,
    explore_solutions = 10,
    explore_tol = 0.1,
    max_solutions = 10,
    fit_all = TRUE,
    comparison_tol = sqrt(eps),
    algorithm_opts = mm_algorithm_options(),
    mscale_opts = mscale_algorithm_options(),
    nlambdas_early = 10,
    early_opts = early_options(),
    ...
)

## S3 method for class 'pense_cvfit'
pensem_cv(
  x,
  scale,
  alpha,
  nlambdas = 50,
  lambda_min_ratio,
  lambda_m,
  standardize = TRUE,
  penalty_loadings,
  intercept = TRUE,
  bdp = 0.25,
  ncores = 1,
  sparse = FALSE,
  eps = 1e-06,
  cc = 4.7,
  cv_k = 5,
  cv_repl = 1,
  cl = NULL,
  cv_metric = c("tau_size", "mape", "rmspe"),
  add_zero_based = TRUE,
  explore_solutions = 10,
  explore_tol = 0.1,
  max_solutions = 10,
  fit_all = TRUE,
  comparison_tol = sqrt(eps),
  algorithm_opts = mm_algorithm_options(),
  mscale_opts = mscale_algorithm_options(),
  x_train,
  y_train,
  ...
)

```

Arguments

<code>x</code>	either a numeric matrix of predictor values, or a cross-validated PENSE fit from <code>pense_cv()</code> .
<code>...</code>	ignored. See the section on deprecated parameters below.
<code>y</code>	vector of response values of length <code>n</code> .
<code>alpha</code>	elastic net penalty mixing parameter with $0 \leq \alpha \leq 1$. <code>alpha = 1</code> is the LASSO penalty, and <code>alpha = 0</code> the Ridge penalty.
<code>nlambda</code>	number of penalization levels.
<code>lambda_min_ratio</code>	Smallest value of the penalization level as a fraction of the largest level (i.e., the smallest value for which all coefficients are zero). The default depends on the sample size relative to the number of variables and <code>alpha</code> . If more observations than variables are available, the default is $1e-3 * \alpha$, otherwise $1e-2 * \alpha$.
<code>lambda_m, lambda_s</code>	optional user-supplied sequence of penalization levels for the S- and M-estimates. If given and not NULL, <code>nlambda</code> and <code>lambda_min_ratio</code> are ignored for the respective estimate (S and/or M).
<code>standardize</code>	logical flag to standardize the <code>x</code> variables prior to fitting the PENSE estimates. Coefficients are always returned on the original scale. This can fail for variables with a large proportion of a single value (e.g., zero-inflated data). In this case, either compute with <code>standardize = FALSE</code> or standardize the data manually.
<code>penalty_loadings</code>	a vector of positive penalty loadings (a.k.a. weights) for different penalization of each coefficient. Only allowed for <code>alpha > 0</code> .
<code>intercept</code>	include an intercept in the model.
<code>bdp</code>	desired breakdown point of the estimator, between 0 and 0.5.
<code>ncores</code>	number of CPU cores to use in parallel. By default, only one CPU core is used. May not be supported on your platform, in which case a warning is given.
<code>sparse</code>	use sparse coefficient vectors.
<code>eps</code>	numerical tolerance.
<code>cc</code>	cutoff constant for Tukey's bisquare ρ function in the M-estimation objective function.
<code>cv_k</code>	number of folds per cross-validation.
<code>cv_repl</code>	number of cross-validation replications.
<code>cl</code>	a parallel cluster. Can only be used if <code>ncores = 1</code> , because multi-threading can not be used in parallel R sessions on the same host.
<code>cv_metric</code>	either a string specifying the performance metric to use, or a function to evaluate prediction errors in a single CV replication. If a function, it is called with a single numeric vector of prediction errors and must return a scalar number.
<code>add_zero_based</code>	also consider the 0-based regularization path. See details for a description.
<code>explore_solutions</code>	number of solutions to compute up to the desired precision <code>eps</code> .
<code>explore_tol</code>	numerical tolerance for exploring possible solutions. Should be (much) looser than <code>eps</code> to be useful.
<code>max_solutions</code>	only retain up to <code>max_solutions</code> unique solutions per penalization level.

fit_all	If TRUE, fit the model for all penalization levels. Otherwise, only at penalization level with smallest average CV performance.
comparison_tol	numeric tolerance to determine if two solutions are equal. The comparison is first done on the absolute difference in the value of the objective function at the solution. If this is less than comparison_tol, two solutions are deemed equal. If the squared difference of the intercepts is less than comparison_tol and the squared L_2 norm of the difference vector is less than comparison_tol.
algorithm_opts	options for the MM algorithm to compute the estimates. See mm_algorithm_options() for details.
mscale_opts	options for the M-scale estimation. See mscale_algorithm_options() for details.
nlambda_ncpy	number of penalization levels where the EN-PY initial estimate is computed.
ncpy_opts	options for the ENPY initial estimates, created with the ncpy_options() function. See ncpy_initial_estimates() for details.
scale	initial scale estimate to use in the M-estimation. By default the S-scale from the PENSE fit is used.
x_train, y_train	override arguments x and y as provided in the call to pense_cv() . This is useful if the arguments in the pense_cv() call are not available in the current environment.

Value

an object of cross-validated regularized M-estimates as returned from [regmest_cv\(\)](#).

See Also

[pense_cv\(\)](#) to compute the starting S-estimate.

Other functions to compute robust estimates with CV: [pense_cv\(\)](#), [regmest_cv\(\)](#)

pense_cv

Cross-validation for (Adaptive) PENSE Estimates

Description

Perform (repeated) K-fold cross-validation for [pense\(\)](#).

[adapense_cv\(\)](#) is a convenience wrapper to compute adaptive PENSE estimates.

Usage

```
pense_cv(
  x,
  y,
  standardize = TRUE,
  lambda,
  cv_k,
  cv_repl = 1,
  cv_metric = c("tau_size", "mape", "rmspe"),
  fit_all = TRUE,
```

```

    cl = NULL,
    ...
)

adapense_cv(x, y, alpha, alpha_preliminary = 0, exponent = 1, ...)

```

Arguments

<code>x</code>	<code>n</code> by <code>p</code> matrix of numeric predictors.
<code>y</code>	vector of response values of length <code>n</code> .
<code>standardize</code>	whether to standardize the <code>x</code> variables prior to fitting the PENSE estimates. Can also be set to "cv_only", in which case the input data is not standardized, but the training data in the CV folds is scaled to match the scaling of the input data. Coefficients are always returned on the original scale. This can fail for variables with a large proportion of a single value (e.g., zero-inflated data). In this case, either compute with <code>standardize = FALSE</code> or standardize the data manually.
<code>lambda</code>	optional user-supplied sequence of penalization levels. If given and not <code>NULL</code> , <code>nlambda</code> and <code>lambda_min_ratio</code> are ignored.
<code>cv_k</code>	number of folds per cross-validation.
<code>cv_repl</code>	number of cross-validation replications.
<code>cv_metric</code>	either a string specifying the performance metric to use, or a function to evaluate prediction errors in a single CV replication. If a function, it is called with a single numeric vector of prediction errors and must return a scalar number.
<code>fit_all</code>	If <code>TRUE</code> , fit the model for all penalization levels. Otherwise, only at penalization level with smallest average CV performance.
<code>cl</code>	a parallel cluster. Can only be used if <code>ncores = 1</code> , because multi-threading can not be used in parallel R sessions on the same host.
<code>...</code>	Arguments passed on to pense
<code>nlambda</code>	number of penalization levels.
<code>lambda_min_ratio</code>	Smallest value of the penalization level as a fraction of the largest level (i.e., the smallest value for which all coefficients are zero). The default depends on the sample size relative to the number of variables and <code>alpha</code> . If more observations than variables are available, the default is $1e-3 * \alpha$, otherwise $1e-2 * \alpha$.
<code>nlambda_ency</code>	number of penalization levels where the EN-PY initial estimate is computed.
<code>penalty_loadings</code>	a vector of positive penalty loadings (a.k.a. weights) for different penalization of each coefficient. Only allowed for <code>alpha > 0</code> .
<code>ency_lambda</code>	optional user-supplied sequence of penalization levels at which EN-PY initial estimates are computed. If given and not <code>NULL</code> , <code>nlambda_ency</code> is ignored.
<code>other_starts</code>	a list of other staring points, created by starting_point() . If the output of ency_initial_estimates() is given, the starting points will be <i>shared</i> among all penalization levels. Note that if a the starting point is <i>specific</i> to a penalization level, this penalization level is added to the grid of penalization levels (either the manually specified grid in <code>lambda</code> or the automatically generated grid of size <code>nlambda</code>). If <code>standardize = TRUE</code> , the starting points are also scaled.
<code>intercept</code>	include an intercept in the model.

bdp	desired breakdown point of the estimator, between 0 and 0.5.
eps	numerical tolerance.
explore_solutions	number of solutions to compute up to the desired precision eps.
explore_tol	numerical tolerance for exploring possible solutions. Should be (much) looser than eps to be useful.
max_solutions	only retain up to max_solutions unique solutions per penalization level.
comparison_tol	numeric tolerance to determine if two solutions are equal. The comparison is first done on the absolute difference in the value of the objective function at the solution. If this is less than comparison_tol, two solutions are deemed equal if the squared difference of the intercepts is less than comparison_tol and the squared L_2 norm of the difference vector is less than comparison_tol.
add_zero_based	also consider the 0-based regularization path. See details for a description.
enpy_specific	use the EN-PY initial estimates only at the penalization level they are computed for. See details for a description.
sparse	use sparse coefficient vectors.
ncores	number of CPU cores to use in parallel. By default, only one CPU core is used. May not be supported on your platform, in which case a warning is given.
algorithm_opts	options for the MM algorithm to compute the estimates. See mm_algorithm_options() for details.
mscale_opts	options for the M-scale estimation. See mscale_algorithm_options() for details.
enpy_opts	options for the ENPY initial estimates, created with the enpy_options() function. See enpy_initial_estimates() for details.
cv_objective	deprecated and ignored. See pense_cv() for estimating prediction performance via cross-validation.
alpha	elastic net penalty mixing parameter with $0 \leq \alpha \leq 1$. alpha = 1 is the LASSO penalty, and alpha = 0 the Ridge penalty.
alpha_preliminary	alpha parameter for the preliminary estimate.
exponent	the exponent for computing the penalty loadings based on the preliminary estimate.

Details

The built-in CV metrics are

"tau_size" τ -size of the prediction error, computed by [tau_size\(\)](#) (default).

"mape" Median absolute prediction error.

"rmspe" Root mean squared prediction error.

`adapense_cv()` is a convenience wrapper which performs 3 steps:

1. compute preliminary estimates via `pense_cv(..., alpha = alpha_preliminary)`,
2. computes the penalty loadings from the estimate beta with best prediction performance by `adapense_loadings = 1 / abs(beta)^exponent`, and
3. compute the adaptive PENSE estimates via `pense_cv(..., penalty_loadings = adapense_loadings)`.

Value

a list with components:

`lambda` the sequence of penalization levels.

`cvres` data frame of average cross-validated performance.

`cv_replications` matrix of cross-validated performance metrics, one column per replication.
Rows are in the same order as in `cvres`.

`call` the original call.

`estimates` the estimates fitted on the full data. Same format as returned by `pense()`.

the object returned by `adapense_cv()` has additional components

`preliminary` the CV results for the preliminary estimate.

`penalty_loadings` the penalty loadings used for the adaptive PENSE estimate.

See Also

`pense()` for computing regularized S-estimates without cross-validation.

`coef.pense_cvfit()` for extracting coefficient estimates.

`plot.pense_cvfit()` for plotting the CV performance or the regularization path.

Other functions to compute robust estimates with CV: `pensem_cv()`, `regmest_cv()`

Other functions to compute robust estimates with CV: `pensem_cv()`, `regmest_cv()`

Examples

```
# Compute the adaptive PENSE regularization path for Freeny's
# revenue data (see ?freeny)
data(freeny)
x <- as.matrix(freeny[, 2:5])

## Either use the convenience function directly ...
ada_convenience <- adapense_cv(x, freeny$y, alpha = 0.5,
                               cv_repl = 2, cv_k = 4)

## ... or compute the steps manually:
# Step 1: Compute preliminary estimates with CV
preliminary_estimate <- pense_cv(x, freeny$y, alpha = 0,
                                 cv_repl = 2, cv_k = 4)
plot(preliminary_estimate, se_mult = 1)

# Step 2: Use the coefficients with best prediction performance
# to define the penalty loadings:
prelim_coefs <- coef(preliminary_estimate, lambda = 'min')
pen_loadings <- 1 / abs(prelim_coefs[-1])

# Step 3: Compute the adaptive PENSE estimates and estimate
# their prediction performance.
ada_manual <- pense_cv(x, freeny$y, alpha = 0.5, cv_repl = 2,
                      cv_k = 4, penalty_loadings = pen_loadings)

# Visualize the prediction performance and coefficient path of
# the adaptive PENSE estimates (manual vs. automatic)
def.par <- par(no.readonly = TRUE)
```



```

layout(matrix(1:4, ncol = 2, byrow = TRUE))
plot(ada_convenience$preliminary)
plot(preliminary_estimate)
plot(ada_convenience)
plot(ada_manual)
par(def.par)

```

pense_options

Deprecated Additional Options for the Penalized EN S-estimator

Description

Deprecated Superseded by [mm_algorithm_options\(\)](#) and options supplied directly to [pense\(\)](#).

Usage

```

pense_options(
  delta = 0.25,
  maxit = 1000,
  eps = 1e-06,
  mscale_eps = 1e-08,
  mscale_maxit = 200,
  verbosity = 0,
  cc = NULL,
  en_correction = TRUE
)

```

Arguments

delta	desired breakdown point of the resulting estimator.
maxit	maximum number of iterations allowed.
eps	numeric tolerance for convergence.
mscale_eps, mscale_maxit	maximum number of iterations and numeric tolerance for the M-scale.
verbosity	ignored. Verbosity of the algorithm.
cc	ignored. Tuning constant for the S-estimator. Default is to chosen based on the breakdown point delta. Should never have to be changed.
en_correction	ignored. Should the corrected EN estimator be used to choose the optimal lambda with CV. If TRUE, as by default, the estimator is "bias corrected".

See Also

Other deprecated functions: [deprecated_en_options](#), [enpy\(\)](#), [initest_options\(\)](#), [mstep_options\(\)](#), [pensem\(\)](#)

plot.pense_cvfit

Plot Method for Penalized Estimates With Cross-Validation

Description

Plot the cross-validation performance or the coefficient path for fitted penalized elastic net S- or LS-estimates of regression.

Usage

```
## S3 method for class 'pense_cvfit'
plot(x, what = c("cv", "coef.path"), se_mult = 1, ...)
```

Arguments

x	fitted estimates with cross-validation information.
what	plot either the CV performance or the coefficient path.
se_mult	if plotting CV performance, multiplier of the estimated SE.
...	currently ignored.

See Also

Other functions for plotting and printing: [plot.pense_fit\(\)](#), [prediction_performance\(\)](#), [summary.pense_cvfit\(\)](#)

Examples

```
# Compute the PENSE regularization path for Freeny's revenue data
# (see ?freeny)
data(freeny)
x <- as.matrix(freeny[, 2:5])

regpath <- pense(x, freeny$y, alpha = 0.5)
plot(regpath)

# Extract the coefficients at a certain penalization level
coef(regpath, lambda = regpath$lambda[40])

# What penalization level leads to good prediction performance?
cv_results <- pense_cv(x, freeny$y, alpha = 0.5, cv_repl = 2,
                      cv_k = 4)
plot(cv_results, se_mult = 1)

# Extract the coefficients at the penalization level with
# smallest prediction error ...
coef(cv_results)
# ... or at the penalization level with prediction error
# statistically indistinguishable from the minimum.
coef(cv_results, lambda = 'se')
```

plot.pense_fit

Plot Method for Penalized Estimates

Description

Plot the coefficient path for fitted penalized elastic net S- or LS-estimates of regression.

Usage

```
## S3 method for class 'pense_fit'
plot(x, ...)
```

Arguments

x	fitted estimates.
...	currently ignored.

See Also

Other functions for plotting and printing: [plot.pense_cvfit\(\)](#), [prediction_performance\(\)](#), [summary.pense_cvfit\(\)](#)

Examples

```
# Compute the PENSE regularization path for Freeny's revenue data
# (see ?freeny)
data(freeny)
x <- as.matrix(freeny[, 2:5])

regpath <- pense(x, freeny$y, alpha = 0.5)
plot(regpath)

# Extract the coefficients at a certain penalization level
coef(regpath, lambda = regpath$lambda[40])

# What penalization level leads to good prediction performance?
cv_results <- pense_cv(x, freeny$y, alpha = 0.5, cv_repl = 2,
                      cv_k = 4)
plot(cv_results, se_mult = 1)

# Extract the coefficients at the penalization level with
# smallest prediction error ...
coef(cv_results)
# ... or at the penalization level with prediction error
# statistically indistinguishable from the minimum.
coef(cv_results, lambda = 'se')
```

predict.pense_cvfit *Predict Method for PENSE Fits*

Description

Predict response values using a PENSE (or LS-EN) regularization path with hyper-parameters chosen by cross-validation.

Usage

```
## S3 method for class 'pense_cvfit'
predict(
  object,
  newdata,
  lambda = c("min", "se"),
  se_mult = 1,
  exact = deprecated(),
  correction = deprecated(),
  ...
)
```

Arguments

object	PENSE with cross-validated hyper-parameters to extract coefficients from.
newdata	an optional matrix of new predictor values. If missing, the fitted values are computed.
lambda	either a string specifying which penalty level to use or a single numeric value of the penalty parameter. See details.
se_mult	If lambda = "se", the multiple of standard errors to tolerate.
exact	deprecated. Always gives a warning if lambda is not part of the fitted sequence and coefficients are interpolated.
correction	defunct.
...	currently not used.

Details

If lambda = "se" and object contains fitted estimates for every penalization level in the sequence, extract the residuals of the most parsimonious model with prediction performance statistically indistinguishable from the best model. This is determined to be the model with prediction performance within $se_mult * cv_se$ from the best model.

Value

a numeric vector of residuals for the given penalization level.

See Also

Other functions for extracting components: [coef.pense_cvfit\(\)](#), [coef.pense_fit\(\)](#), [predict.pense_fit\(\)](#), [residuals.pense_cvfit\(\)](#), [residuals.pense_fit\(\)](#)

Examples

```
# Compute the LS-EN regularization path for Freeny's revenue data
# (see ?freeny)
data(freeny)
x <- as.matrix(freeny[, 2:5])

regpath <- elnet(x, freeny$y, alpha = 0.75)

# Predict the response using a specific penalization level
predict(regpath, newdata = freeny[1:5, 2:5],
        lambda = regpath$lambda[10])

# Extract the residuals at a certain penalization level
residuals(regpath, lambda = regpath$lambda[5])

# Select penalization level via cross-validation
cv_results <- elnet_cv(x, freeny$y, alpha = 0.5, cv_repl = 10,
                      cv_k = 4)

# Predict the response using the "best" penalization level
predict(cv_results, newdata = freeny[1:5, 2:5])

# Extract the residuals at the "best" penalization level
residuals(cv_results)^2
# Extract the residuals at a more parsimonious penalization level
residuals(cv_results, lambda = 'se')
```

predict.pense_fit	<i>Predict Method for PENSE Fits</i>
-------------------	--------------------------------------

Description

Predict response values using a PENSE (or LS-EN) regularization path fitted by `pense()` or `elnet()`.

Usage

```
## S3 method for class 'pense_fit'
predict(
  object,
  newdata,
  lambda,
  exact = deprecated(),
  correction = deprecated(),
  ...
)
```

Arguments

object	PENSE regularization path to extract residuals from.
newdata	an optional matrix of new predictor values. If missing, the fitted values are computed.
lambda	a single value of the penalty parameter.

exact	defunct Always gives a warning if <code>lambda</code> is not part of the fitted sequence and coefficients need to be interpolated.
correction	defunct.
...	currently not used.

Value

a numeric vector of residuals for the given penalization level.

See Also

Other functions for extracting components: `coef.pense_cvfit()`, `coef.pense_fit()`, `predict.pense_cvfit()`, `residuals.pense_cvfit()`, `residuals.pense_fit()`

Examples

```
# Compute the LS-EN regularization path for Freeny's revenue data
# (see ?freeny)
data(freeny)
x <- as.matrix(freeny[, 2:5])

regpath <- elnet(x, freeny$y, alpha = 0.75)

# Predict the response using a specific penalization level
predict(regpath, newdata = freeny[1:5, 2:5],
        lambda = regpath$lambda[10])

# Extract the residuals at a certain penalization level
residuals(regpath, lambda = regpath$lambda[5])

# Select penalization level via cross-validation
cv_results <- elnet_cv(x, freeny$y, alpha = 0.5, cv_repl = 10,
                      cv_k = 4)

# Predict the response using the "best" penalization level
predict(cv_results, newdata = freeny[1:5, 2:5])

# Extract the residuals at the "best" penalization level
residuals(cv_results)^2
# Extract the residuals at a more parsimonious penalization level
residuals(cv_results, lambda = 'se')
```

prediction_performance

Prediction Performance of Adaptive PENSE Fits

Description

Extract the prediction performance of one or more (adaptive) PENSE fits.

Usage

```
prediction_performance(..., lambda = c("min", "se"), se_mult = 1)

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

Arguments

<code>...</code>	one or more (adaptive) PENSE fits with cross-validation information.
<code>lambda</code>	a string specifying which penalty level to use ("min" or "se") . See details.
<code>se_mult</code>	If <code>lambda = "se"</code> , the multiple of standard errors to tolerate.
<code>x</code>	an object with information on prediction performance created with <code>prediction_performance()</code> .

Details

If `lambda = "se"` and the cross-validation was performed with multiple replications, use the penalty level with prediction performance within `se_mult` of the best prediction performance.

Value

a data frame with details about the prediction performance of the given PENSE fits. The data frame has a custom print method summarizing the prediction performances.

See Also

[summary.pense_cvfit\(\)](#) for a summary of the fitted model.

Other functions for plotting and printing: [plot.pense_cvfit\(\)](#), [plot.pense_fit\(\)](#), [summary.pense_cvfit\(\)](#)

prinsens

Principal Sensitivity Components

Description

Compute Principal Sensitivity Components for Elastic Net Regression

Usage

```
prinsens(
  x,
  y,
  alpha,
  lambda,
  intercept = TRUE,
  penalty_loadings,
  en_algorithm_opts,
  eps = 1e-06,
  sparse = FALSE,
  ncores = 1L,
  method = deprecated()
)
```

Arguments

<code>x</code>	<code>n</code> by <code>p</code> matrix of numeric predictors.
<code>y</code>	vector of response values of length <code>n</code> .
<code>alpha</code>	elastic net penalty mixing parameter with $0 \leq \alpha \leq 1$. <code>alpha = 1</code> is the LASSO penalty, and <code>alpha = 0</code> the Ridge penalty.
<code>lambda</code>	optional user-supplied sequence of penalization levels. If given and not <code>NULL</code> , <code>nlambda</code> and <code>lambda_min_ratio</code> are ignored.
<code>intercept</code>	include an intercept in the model.
<code>penalty_loadings</code>	a vector of positive penalty loadings (a.k.a. weights) for different penalization of each coefficient. Only allowed for <code>alpha > 0</code> .
<code>en_algorithm_opts</code>	options for the LS-EN algorithm. See en_algorithm_options for details.
<code>eps</code>	numerical tolerance.
<code>sparse</code>	use sparse coefficient vectors.
<code>ncores</code>	number of CPU cores to use in parallel. By default, only one CPU core is used. May not be supported on your platform, in which case a warning is given.
<code>method</code>	defunct. PSCs are always computed for EN estimates. For the PY procedure for unpenalized estimation use package pyinit .

Value

a list of principal sensitivity components, one per element in `lambda`. Each PSC is itself a list with items `lambda`, `alpha`, and `pscs`.

References

- Cohen Freue, G.V.; Kepplinger, D.; Salibián-Barrera, M.; Smucler, E. Robust elastic net estimators for variable selection and identification of proteomic biomarkers. *Ann. Appl. Stat.* **13** (2019), no. 4, 2065–2090 doi: [10.1214/19AOAS1269](#)
- Pena, D., and Yohai, V.J. A Fast Procedure for Outlier Diagnostics in Large Regression Problems. *J. Amer. Statist. Assoc.* **94** (1999). no. 446, 434–445. doi: [10.2307/2670164](#)

See Also

Other functions for initial estimates: [enpy_initial_estimates\(\)](#), [starting_point\(\)](#)

regmest

Compute (Adaptive) Elastic Net M-Estimates of Regression

Description

Compute elastic net M-estimates along a grid of penalization levels with optional penalty loadings for adaptive elastic net.

Usage

```

regmest(
  x,
  y,
  alpha,
  nlambdas = 50,
  lambda,
  lambda_min_ratio,
  scale,
  starting_points,
  penalty_loadings,
  intercept = TRUE,
  cc = 4.7,
  eps = 1e-06,
  explore_solutions = 10,
  explore_tol = 0.1,
  max_solutions = 10,
  comparison_tol = sqrt(eps),
  sparse = FALSE,
  ncores = 1,
  standardize = TRUE,
  algorithm_opts = mm_algorithm_options(),
  add_zero_based = TRUE,
  mscale_bdp = 0.25,
  mscale_opts = mscale_algorithm_options()
)

```

Arguments

<code>x</code>	n by p matrix of numeric predictors.
<code>y</code>	vector of response values of length n.
<code>alpha</code>	elastic net penalty mixing parameter with $0 \leq \alpha \leq 1$. <code>alpha = 1</code> is the LASSO penalty, and <code>alpha = 0</code> the Ridge penalty.
<code>nlambdas</code>	number of penalization levels.
<code>lambda</code>	optional user-supplied sequence of penalization levels. If given and not NULL, <code>nlambdas</code> and <code>lambda_min_ratio</code> are ignored.
<code>lambda_min_ratio</code>	Smallest value of the penalization level as a fraction of the largest level (i.e., the smallest value for which all coefficients are zero). The default depends on the sample size relative to the number of variables and <code>alpha</code> . If more observations than variables are available, the default is $1e-3 * \alpha$, otherwise $1e-2 * \alpha$.
<code>scale</code>	fixed scale of the residuals.
<code>starting_points</code>	a list of starting points, created by starting_point() . The starting points are shared among all penalization levels.
<code>penalty_loadings</code>	a vector of positive penalty loadings (a.k.a. weights) for different penalization of each coefficient. Only allowed for <code>alpha > 0</code> .
<code>intercept</code>	include an intercept in the model.

<code>cc</code>	cutoff constant for Tukey's bisquare ρ function.
<code>eps</code>	numerical tolerance.
<code>explore_solutions</code>	number of solutions to compute up to the desired precision <code>eps</code> .
<code>explore_tol</code>	numerical tolerance for exploring possible solutions. Should be (much) looser than <code>eps</code> to be useful.
<code>max_solutions</code>	only retain up to <code>max_solutions</code> unique solutions per penalization level.
<code>comparison_tol</code>	numeric tolerance to determine if two solutions are equal. The comparison is first done on the absolute difference in the value of the objective function at the solution. If this is less than <code>comparison_tol</code> , two solutions are deemed equal if the squared difference of the intercepts is less than <code>comparison_tol</code> and the squared L_2 norm of the difference vector is less than <code>comparison_tol</code> .
<code>sparse</code>	use sparse coefficient vectors.
<code>ncores</code>	number of CPU cores to use in parallel. By default, only one CPU core is used. May not be supported on your platform, in which case a warning is given.
<code>standardize</code>	logical flag to standardize the x variables prior to fitting the M-estimates. Coefficients are always returned on the original scale. This can fail for variables with a large proportion of a single value (e.g., zero-inflated data). In this case, either compute with <code>standardize = FALSE</code> or standardize the data manually.
<code>algorithm_opts</code>	options for the MM algorithm to compute estimates. See mm_algorithm_options() for details.
<code>add_zero_based</code>	also consider the 0-based regularization path in addition to the given starting points.
<code>mscale_bdp, mscale_opts</code>	options for the M-scale estimate used to standardize the predictors (if <code>standardize = TRUE</code>).

Value

a list-like object with the following items

`lambda` the sequence of penalization levels.

`scale` the used scale of the residuals.

`estimates` a list of estimates. Each estimate contains the following information:

`intercept` intercept estimate.

`beta` beta (slope) estimate.

`lambda` penalization level at which the estimate is computed.

`alpha` *alpha* hyper-parameter at which the estimate is computed.

`objf_value` value of the objective function at the solution.

`statuscode` if > 0 the algorithm experienced issues when computing the estimate.

`status` optional status message from the algorithm.

`call` the original call.

See Also

[regmest_cv\(\)](#) for selecting hyper-parameters via cross-validation.

[coef.pense_fit\(\)](#) for extracting coefficient estimates.

[plot.pense_fit\(\)](#) for plotting the regularization path.

Other functions to compute robust estimates: [pense\(\)](#)

regmest_cv

*Cross-validation for (Adaptive) Elastic Net M-Estimates***Description**

Perform (repeated) K-fold cross-validation for `regmest()`.

`adamest_cv()` is a convenience wrapper to compute adaptive elastic-net M-estimates.

Usage

```
regmest_cv(
  x,
  y,
  standardize = TRUE,
  lambda,
  cv_k,
  cv_repl = 1,
  cv_metric = c("tau_size", "mape", "rmspe"),
  fit_all = TRUE,
  cl = NULL,
  ...
)
```

```
adamest_cv(x, y, alpha, alpha_preliminary = 0, exponent = 1, ...)
```

Arguments

<code>x</code>	<code>n</code> by <code>p</code> matrix of numeric predictors.
<code>y</code>	vector of response values of length <code>n</code> .
<code>standardize</code>	whether to standardize the <code>x</code> variables prior to fitting the PENSE estimates. Can also be set to <code>"cv_only"</code> , in which case the input data is not standardized, but the training data in the CV folds is scaled to match the scaling of the input data. Coefficients are always returned on the original scale. This can fail for variables with a large proportion of a single value (e.g., zero-inflated data). In this case, either compute with <code>standardize = FALSE</code> or standardize the data manually.
<code>lambda</code>	optional user-supplied sequence of penalization levels. If given and not <code>NULL</code> , <code>nlambda</code> and <code>lambda_min_ratio</code> are ignored.
<code>cv_k</code>	number of folds per cross-validation.
<code>cv_repl</code>	number of cross-validation replications.
<code>cv_metric</code>	either a string specifying the performance metric to use, or a function to evaluate prediction errors in a single CV replication. If a function, it is called with a single numeric vector of prediction errors and must return a scalar number.
<code>fit_all</code>	If <code>TRUE</code> , fit the model for all penalization levels. Otherwise, only at penalization level with smallest average CV performance.
<code>cl</code>	a parallel cluster. Can only be used if <code>ncores = 1</code> , because multi-threading can not be used in parallel R sessions on the same host.
<code>...</code>	Arguments passed on to regmest
	<code>scale</code> fixed scale of the residuals.

<code>nlambda</code>	number of penalization levels.
<code>lambda_min_ratio</code>	Smallest value of the penalization level as a fraction of the largest level (i.e., the smallest value for which all coefficients are zero). The default depends on the sample size relative to the number of variables and <code>alpha</code> . If more observations than variables are available, the default is $1e-3 * \alpha$, otherwise $1e-2 * \alpha$.
<code>penalty_loadings</code>	a vector of positive penalty loadings (a.k.a. weights) for different penalization of each coefficient. Only allowed for <code>alpha > 0</code> .
<code>starting_points</code>	a list of starting points, created by <code>starting_point()</code> . The starting points are shared among all penalization levels.
<code>intercept</code>	include an intercept in the model.
<code>add_zero_based</code>	also consider the 0-based regularization path in addition to the given starting points.
<code>cc</code>	cutoff constant for Tukey's bisquare ρ function.
<code>eps</code>	numerical tolerance.
<code>explore_solutions</code>	number of solutions to compute up to the desired precision <code>eps</code> .
<code>explore_tol</code>	numerical tolerance for exploring possible solutions. Should be (much) looser than <code>eps</code> to be useful.
<code>max_solutions</code>	only retain up to <code>max_solutions</code> unique solutions per penalization level.
<code>comparison_tol</code>	numeric tolerance to determine if two solutions are equal. The comparison is first done on the absolute difference in the value of the objective function at the solution. If this is less than <code>comparison_tol</code> , two solutions are deemed equal if the squared difference of the intercepts is less than <code>comparison_tol</code> and the squared L_2 norm of the difference vector is less than <code>comparison_tol</code> .
<code>sparse</code>	use sparse coefficient vectors.
<code>ncores</code>	number of CPU cores to use in parallel. By default, only one CPU core is used. May not be supported on your platform, in which case a warning is given.
<code>algorithm_opts</code>	options for the MM algorithm to compute estimates. See <code>mm_algorithm_options()</code> for details.
<code>mscale_bdp</code>	options for the M-scale estimate used to standardize the predictors (if <code>standardize = TRUE</code>).
<code>mscale_opts</code>	options for the M-scale estimate used to standardize the predictors (if <code>standardize = TRUE</code>).
<code>alpha</code>	elastic net penalty mixing parameter with $0 \leq \alpha \leq 1$. <code>alpha = 1</code> is the LASSO penalty, and <code>alpha = 0</code> the Ridge penalty.
<code>alpha_preliminary</code>	alpha parameter for the preliminary estimate.
<code>exponent</code>	the exponent for computing the penalty loadings based on the preliminary estimate.

Details

The built-in CV metrics are

"tau_size" τ -size of the prediction error, computed by `tau_size()` (default).

"mape" Median absolute prediction error.

"rmspe" Root mean squared prediction error.

adamest_cv() is a convenience wrapper which performs 3 steps:

1. compute preliminary estimates via `regmest_cv(..., alpha = alpha_preliminary)`,
2. computes the penalty loadings from the estimate beta with best prediction performance by `adamest_loadings = 1 / abs(beta)^exponent`, and
3. compute the adaptive PENSE estimates via `regmest_cv(..., penalty_loadings = adamest_loadings)`.

Value

a list with components:

`lambda` the sequence of penalization levels.

`scale` the used scale of the residuals.

`cvres` data frame of average cross-validated performance.

`cv_replications` matrix of cross-validated performance metrics, one column per replication.
Rows are in the same order as in `cvres`.

`call` the original call.

`estimates` the estimates fitted on the full data. Same format as returned by `regmest()`.

the object returned by `adamest_cv()` has additional components

`preliminary` the CV results for the preliminary estimate.

`penalty_loadings` the penalty loadings used for the adaptive elastic net M-estimate.

See Also

`regmest()` for computing regularized S-estimates without cross-validation.

`coef.pense_cvfit()` for extracting coefficient estimates.

`plot.pense_cvfit()` for plotting the CV performance or the regularization path.

Other functions to compute robust estimates with CV: `pense_cv()`, `pensem_cv()`

Other functions to compute robust estimates with CV: `pense_cv()`, `pensem_cv()`

Examples

```
# Compute the adaptive PENSE regularization path for Freeny's
# revenue data (see ?freeny)
data(freeny)
x <- as.matrix(freeny[, 2:5])

## Either use the convenience function directly ...
ada_convenience <- adapense_cv(x, freeny$y, alpha = 0.5,
                               cv_repl = 2, cv_k = 4)

## ... or compute the steps manually:
# Step 1: Compute preliminary estimates with CV
preliminary_estimate <- pense_cv(x, freeny$y, alpha = 0,
                                 cv_repl = 2, cv_k = 4)
plot(preliminary_estimate, se_mult = 1)

# Step 2: Use the coefficients with best prediction performance
# to define the penalty loadings:
```

```

prelim_coefs <- coef(preliminary_estimate, lambda = 'min')
pen_loadings <- 1 / abs(prelim_coefs[-1])

# Step 3: Compute the adaptive PENSE estimates and estimate
# their prediction performance.
ada_manual <- pense_cv(x, freeny$y, alpha = 0.5, cv_repl = 2,
                      cv_k = 4, penalty_loadings = pen_loadings)

# Visualize the prediction performance and coefficient path of
# the adaptive PENSE estimates (manual vs. automatic)
def.par <- par(no.readonly = TRUE)
layout(matrix(1:4, ncol = 2, byrow = TRUE))
plot(ada_convenience$preliminary)
plot(preliminary_estimate)
plot(ada_convenience)
plot(ada_manual)
par(def.par)

```

residuals.pense_cvfit *Extract Residuals*

Description

Extract residuals from a PENSE (or LS-EN) regularization path with hyper-parameters chosen by cross-validation.

Usage

```

## S3 method for class 'pense_cvfit'
residuals(
  object,
  lambda = c("min", "se"),
  se_mult = 1,
  exact = deprecated(),
  correction = deprecated(),
  ...
)

```

Arguments

object	PENSE with cross-validated hyper-parameters to extract coefficients from.
lambda	either a string specifying which penalty level to use or a single numeric value of the penalty parameter. See details.
se_mult	If lambda = "se", the multiple of standard errors to tolerate.
exact	deprecated. Always gives a warning if lambda is not part of the fitted sequence and coefficients are interpolated.
correction	defunct.
...	currently not used.

Details

If `lambda = "se"` and object contains fitted estimates for every penalization level in the sequence, extract the residuals of the most parsimonious model with prediction performance statistically indistinguishable from the best model. This is determined to be the model with prediction performance within `se_mult * cv_se` from the best model.

Value

a numeric vector of residuals for the given penalization level.

See Also

Other functions for extracting components: `coef.pense_cvfit()`, `coef.pense_fit()`, `predict.pense_cvfit()`, `predict.pense_fit()`, `residuals.pense_fit()`

Examples

```
# Compute the LS-EN regularization path for Freeny's revenue data
# (see ?freeny)
data(freeny)
x <- as.matrix(freeny[, 2:5])

regpath <- elnet(x, freeny$y, alpha = 0.75)

# Predict the response using a specific penalization level
predict(regpath, newdata = freeny[1:5, 2:5],
        lambda = regpath$lambda[10])

# Extract the residuals at a certain penalization level
residuals(regpath, lambda = regpath$lambda[5])

# Select penalization level via cross-validation
cv_results <- elnet_cv(x, freeny$y, alpha = 0.5, cv_repl = 10,
                      cv_k = 4)

# Predict the response using the "best" penalization level
predict(cv_results, newdata = freeny[1:5, 2:5])

# Extract the residuals at the "best" penalization level
residuals(cv_results)^2
# Extract the residuals at a more parsimonious penalization level
residuals(cv_results, lambda = 'se')
```

`residuals.pense_fit` *Extract Residuals*

Description

Extract residuals from a PENSE (or LS-EN) regularization path fitted by `pense()` or `elnet()`.

Usage

```
## S3 method for class 'pense_fit'
residuals(object, lambda, exact = deprecated(), correction = deprecated(), ...)
```

Arguments

object	PENSE regularization path to extract residuals from.
lambda	a single value of the penalty parameter.
exact	defunct Always gives a warning if lambda is not part of the fitted sequence and coefficients need to be interpolated.
correction	defunct.
...	currently not used.

Value

a numeric vector of residuals for the given penalization level.

See Also

Other functions for extracting components: `coef.pense_cvfit()`, `coef.pense_fit()`, `predict.pense_cvfit()`, `predict.pense_fit()`, `residuals.pense_cvfit()`

Examples

```
# Compute the LS-EN regularization path for Freeny's revenue data
# (see ?freeny)
data(freeny)
x <- as.matrix(freeny[, 2:5])

regpath <- elnet(x, freeny$y, alpha = 0.75)

# Predict the response using a specific penalization level
predict(regpath, newdata = freeny[1:5, 2:5],
        lambda = regpath$lambda[10])

# Extract the residuals at a certain penalization level
residuals(regpath, lambda = regpath$lambda[5])

# Select penalization level via cross-validation
cv_results <- elnet_cv(x, freeny$y, alpha = 0.5, cv_repl = 10,
                      cv_k = 4)

# Predict the response using the "best" penalization level
predict(cv_results, newdata = freeny[1:5, 2:5])

# Extract the residuals at the "best" penalization level
residuals(cv_results)^2
# Extract the residuals at a more parsimonious penalization level
residuals(cv_results, lambda = 'se')
```

rho_function

List Available Rho Functions

Description

List Available Rho Functions

Usage

```
rho_function(rho)
```

Arguments

rho the name of the ρ function to check for existence.

Value

if rho is missing returns a vector of supported ρ function names, otherwise the internal integer representation of the ρ function.

See Also

Other miscellaneous functions: [consistency_const\(\)](#)

starting_point	<i>Create Starting Points for the PENSE Algorithm</i>
----------------	-------------------------------------------------------

Description

Create a starting point for starting the PENSE algorithm in [pense\(\)](#). Multiple starting points can be created by combining starting points via `c(starting_point_1, starting_point_2, ...)`.

Usage

```
starting_point(beta, intercept, lambda)

as_starting_point(object, specific = FALSE, ...)

## S3 method for class 'enpy_starting_points'
as_starting_point(object, specific = FALSE, ...)

## S3 method for class 'pense_fit'
as_starting_point(object, specific = FALSE, lambda, ...)

## S3 method for class 'pense_cvfit'
as_starting_point(
  object,
  specific = FALSE,
  lambda = c("min", "se"),
  se_mult = 1,
  ...
)
```

Arguments

beta beta coefficients at the starting point. Can be a numeric vector, a sparse vector of class [dsparseVector](#), or a sparse matrix of class [dgCMatrix](#) with a single column.

intercept intercept coefficient at the starting point.

lambda	optionally either a string specifying which penalty level to use ("min" or "se") or a numeric vector of the penalty levels to extract from object. Penalization levels not present in object are ignored with a warning. If NULL, all estimates in object are extracted.
object	an object with estimates to use as starting points.
specific	whether the estimates should be used as starting points only at the penalization level they are computed for. Defaults to using the estimates as starting points for all penalization levels.
...	further arguments passed to or from other methods.
se_mult	If lambda = "se", the multiple of standard errors to tolerate.

Details

A starting points can either be *shared*, i.e., used for every penalization level PENSE estimates are computed for, or *specific* to one penalization level. To create a specific starting point, provide the penalization level as lambda. If lambda is missing, a shared starting point is created. Shared and specific starting points can all be combined into a single list of starting points, with [pense\(\)](#) handling them correctly. Note that specific starting points will lead to the lambda value being added to the grid of penalization levels. See [pense\(\)](#) for details.

Starting points computed via [enpy_initial_estimates\(\)](#) are by default *shared* starting points but can be transformed to *specific* starting points via `enpy_starting_point(..., specific = TRUE)`.

When creating starting points from cross-validated fits, it is possible to extract only the estimate with best CV performance (lambda = "min"), or the estimate with CV performance statistically indistinguishable from the best performance (lambda = "se"). This is determined to be the estimate with prediction performance within `se_mult * cv_se` from the best model.

Value

an object of type `starting_points` to be used as starting point for [pense\(\)](#).

See Also

Other functions for initial estimates: [enpy_initial_estimates\(\)](#), [prinsens\(\)](#)

summary.pense_cvfit	<i>Summarize Cross-Validated PENSE Fit</i>
---------------------	--------------------------------------------

Description

If lambda = "se" and object contains fitted estimates for every penalization level in the sequence, extract the coefficients of the most parsimonious model with prediction performance statistically indistinguishable from the best model. This is determined to be the model with prediction performance within `se_mult * cv_se` from the best model.

Usage

```
## S3 method for class 'pense_cvfit'
summary(object, lambda = c("min", "se"), se_mult = 1, ...)

## S3 method for class 'pense_cvfit'
print(x, lambda = c("min", "se"), se_mult = 1, ...)
```

Arguments

object, x	an (adaptive) PENSE fit with cross-validation information.
lambda	either a string specifying which penalty level to use ("min" or "se") or a single numeric value of the penalty parameter. See details.
se_mult	If lambda = "se", the multiple of standard errors to tolerate.
...	ignored.

See Also

[prediction_performance\(\)](#) for information about the estimated prediction performance.

[coef.pense_cvfit\(\)](#) for extracting only the estimated coefficients.

Other functions for plotting and printing: [plot.pense_cvfit\(\)](#), [plot.pense_fit\(\)](#), [prediction_performance\(\)](#)

tau_size

*Compute the Tau-Scale of Centered Values***Description**

Compute the τ -scale without centering the values.

Usage

```
tau_size(x)
```

Arguments

x numeric values. Missing values are verbosely ignored.

Value

the τ estimate of scale of centered values.

See Also

Other functions to compute robust estimates of location and scale: [mlocscale\(\)](#), [mloc\(\)](#), [mscale\(\)](#)

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