

Package ‘L1pack’

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Title Routines for L1 Estimation

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Description L1 estimation for linear regression using Barrodale and Roberts' method
<[doi:10.1145/355616.361024](https://doi.org/10.1145/355616.361024)> and the EM algorithm <[doi:10.1023/A:1020759012226](https://doi.org/10.1023/A:1020759012226)>,
density, distribution function, quantile function and random number generation
for univariate and multivariate Laplace distribution.

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LinkingTo fastmatrix

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confint.lad	<i>Confidence intervals from lad models</i>
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Description

Computes confidence intervals for one or more parameters in a fitted model corresponding to a lad object.

Usage

```
## S3 method for class 'lad'
confint(object, parm, level = 0.95, ...)
```

Arguments

object	a fitted model object.
parm	a specification of which parameters are to be given confidence intervals, either a vector of numbers or a vector of names. If missing, all parameters are considered.
level	the confidence level required.
...	additional argument(s) for methods.

Details

confint is a generic function. Confidence intervals associated to lad objects are asymptotic, and needs suitable [coef](#) and [vcov](#) methods to be available.

Value

A matrix (or vector) with columns giving lower and upper confidence limits for each parameter. These will be labelled as $(1-\text{level})/2$ and $1 - (1-\text{level})/2$ in % (by default 2.5% and 97.5%).

See Also

[confint.glm](#) and [confint.nls](#) in package **MASS**.

Examples

```
fm <- lad(stack.loss ~ ., data = stackloss, method = "BR")
confint(fm) # based on asymptotic normality
```

l1fit

Minimum absolute residual (L1) regression

Description

Performs an L1 regression on a matrix of explanatory variables and a vector of responses.

Usage

```
l1fit(x, y, intercept = TRUE, tolerance = 1e-07, print.it = TRUE)
```

Arguments

x	vector or matrix of explanatory variables. Each row corresponds to an observation and each column to a variable. The number of rows of x should equal the number of data values in y, and there should be fewer columns than rows. Missing values are not allowed.
y	numeric vector containing the response. Missing values are not allowed.
intercept	logical flag. If TRUE, an intercept term is included in the regression model.
tolerance	numerical value used to test for singularity in the regression.
print.it	logical flag. If TRUE, then warnings about non-unique solutions and rank deficiency are given.

Details

The Barrodale-Roberts algorithm, which is a specialized linear programming algorithm, is used.

Value

	list defining the regression (compare with function lsfit).
coefficients	vector of coefficients.
residuals	residuals from the fit.
message	vector of one or two character strings stating whether a non-unique solution is possible, or if the x matrix was found to be rank deficient.

References

Barrodale, I., and Roberts, F.D.K. (1973). An improved algorithm for discrete L1 linear approximations. *SIAM Journal of Numerical Analysis* **10**, 839-848.

Barrodale, I., and Roberts, F.D.K. (1974). Solution of an overdetermined system of equations in the L1 norm. *Communications of the ACM* **17**, 319-320.

Bloomfield, P., and Steiger, W.L. (1983). *Least Absolute Deviations: Theory, Applications, and Algorithms*. Birkhauser, Boston, Mass.

Examples

```
l1fit(stack.x, stack.loss)
```

 lad

Least absolute deviations regression

Description

This function is used to fit linear models considering Laplace errors.

Usage

```
lad(formula, data, subset, na.action, method = "BR", tol = 1e-7, maxiter = 500,
    model = TRUE, x = FALSE, y = FALSE, contrasts = NULL)
```

Arguments

formula	an object of class "formula": a symbolic description of the model to be fitted.
data	an optional data frame containing the variables in the model. If not found in data, the variables are taken from <code>environment(formula)</code> , typically the environment from which <code>lad</code> is called.
subset	an optional expression indicating the subset of the rows of data that should be used in the fit.
na.action	a function that indicates what should happen when the data contain NAs.
method	character string specifying the fitting method to be used; the options are "BR" Barrodale and Roberts' method (the default) and "EM" for an EM algorithm using IRLS.
tol	the relative tolerance for the iterative algorithm. Default is <code>tol = 1e-7</code> .
maxiter	The maximum number of iterations for the EM method. Default to 500.
model, x, y	logicals. If TRUE the corresponding components of the fit (the model frame, the model matrix, the response) are returned.
contrasts	an optional list. See the <code>contrasts.arg</code> of <code>model.matrix.default</code> .

Value

an object of class `lad` representing the linear model fit. Generic function `print`, show the results of the fit.

The functions `print` and `summary` are used to obtain and print a summary of the results. The generic accessor functions `coefficients`, `fitted.values` and `residuals` extract various useful features of the value returned by `lad`.

Author(s)

The design was inspired by the R function [lm](#).

References

- Barrodale, I., and Roberts, F.D.K. (1974). Solution of an overdetermined system of equations in the L1 norm. *Communications of the ACM* **17**, 319-320.
- Phillips, R.F. (2002). Least absolute deviations estimation via the EM algorithm. *Statistics and Computing* **12**, 281-285.

Examples

```
fm <- lad(stack.loss ~ ., data = stackloss, method = "BR")
summary(fm)
```

lad.fit	<i>Fitter functions for least absolute deviation (LAD) regression</i>
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Description

This function is a *switcher* among various numerical fitting functions ([lad.fit.BR](#), and [lad.fit.EM](#)). The argument method does the switching: "BR" for [lad.fit.BR](#), etc. This should usually *not* be used directly unless by experienced users.

Usage

```
lad.fit(x, y, method = "BR", tol = 1e-7, maxiter = 500)
```

Arguments

x	design matrix of dimension $n \times p$.
y	vector of observations of length n .
method	currently, methods "BR" (default), and "EM" are supported.
tol	the relative tolerance for the iterative algorithm. Default is $\text{tol} = 1e-7$.
maxiter	The maximum number of iterations for the EM method. Default to 500.

Value

a [list](#) with components:

coefficients	a named vector of coefficients.
scale	final scale estimate of the random error.
residuals	the residuals, that is response minus fitted values.
fitted.values	the fitted values.
SAD	the sum of absolute deviations.
weights	estimated EM weights.
basic	basic observations, that is observations with zero residuals.
logLik	the log-likelihood at convergence.

See Also

[lad.fit.BR](#), [lad.fit.EM](#).

Examples

```
x <- cbind(1, stack.x)
fm <- lad.fit(x, stack.loss, method = "BR")
fm
```

lad.fit-methods

Fit a least absolute deviation (LAD) regression model

Description

Fits a linear model using LAD methods, returning the bare minimum computations.

Usage

```
lad.fit.BR(x, y, tol = 1e-7)
lad.fit.EM(x, y, tol = 1e-7, maxiter = 500)
```

Arguments

<code>x, y</code>	numeric vectors or matrices for the predictors and the response in a linear model. Typically, but not necessarily, <code>x</code> will be constructed by one of the fitting functions.
<code>tol</code>	the relative tolerance for the iterative algorithm. Default is <code>tol = 1e-7</code> .
<code>maxiter</code>	The maximum number of iterations for the EM method. Default to 500.

Value

The bare bones of a `lad` object: the coefficients, residuals, fitted values, and some information used by `summary.lad`.

See Also

[lad](#), [lad.fit](#), [lm](#)

Examples

```
x <- cbind(1, stack.x)
z <- lad.fit.BR(x, stack.loss)
z
```

Laplace

*The Laplace distribution***Description**

Density, distribution function, quantile function and random generation for the Laplace distribution with location parameter *location* and scale parameter *scale*.

Usage

```
dlaplace(x, location = 0, scale = 1, log = FALSE)
plaplace(q, location = 0, scale = 1, lower.tail = TRUE, log.p = FALSE)
qlaplace(p, location = 0, scale = 1, lower.tail = TRUE, log.p = FALSE)
rlaplace(n, location = 0, scale = 1)
```

Arguments

<i>x</i> , <i>q</i>	vector of quantiles.
<i>location</i> , <i>scale</i>	location and scale parameters. Scale must be positive.
<i>log</i> , <i>log.p</i>	logical; if TRUE, probabilities <i>p</i> are given as $\log(p)$.
<i>lower.tail</i>	logical; if TRUE (default), probabilities are $P[X \leq x]$, otherwise, $P[X > x]$.
<i>p</i>	vector of probabilities.
<i>n</i>	number of observations. If $\text{length}(n) > 1$, the length is taken to be the number required.

Details

If *location* or *scale* are not specified, they assume the default values of 0 and 1 respectively.

The Laplace distribution with location μ and scale ϕ has density

$$f(x) = \frac{1}{\sqrt{2}\phi} \exp(-\sqrt{2}|x - \mu|/\phi)$$

Value

dlaplace, *plaplace*, and *qlaplace* are respectively the density, distribution function and quantile function of the Laplace distribution. *rlaplace* generates random deviates from the Laplace.

The length of the result is determined by *n* for *rlaplace*, and is the maximum of the lengths of the numerical parameters for the other functions.

Author(s)

Felipe Osorio and Tymoteusz Wolodzko

References

Kotz, S., Kozubowski, T.J., and Podgorski, K. (2001). *The Laplace Distributions and Generalizations*. Birkhauser, Boston.

Phillips, R.F. (2002). Least absolute deviations estimation via the EM algorithm. *Statistics and Computing* **12**, 281-285.

See Also

[Distributions](#) for other standard distributions and [rmLaplace](#) for the random generation from the multivariate Laplace distribution.

Examples

```
x <- rlaplace(1000)
## Q-Q plot for Laplace data against true theoretical distribution:
qqplot(qlaplace(ppoints(1000)), x, main = "Laplace Q-Q plot",
       xlab = "Theoretical quantiles", ylab = "Sample quantiles")
abline(c(0,1), col = "red", lwd = 2)
```

rmLaplace

Multivariate Laplace random deviates

Description

Random number generation from the multivariate Laplace distribution.

Usage

```
rmLaplace(n = 1, center = rep(0, nrow(Scatter)), Scatter = diag(length(center)))
```

Arguments

n	the number of samples requested
center	a vector giving the locations of each variable
Scatter	a positive-definite dispersion matrix

Details

The function rmLaplace is an interface to C routines, which make calls to subroutines from LAPACK. The matrix decomposition is internally done using the Cholesky decomposition. If Scatter is not non-negative definite then there will be a warning message.

Value

If n = 1 a vector of the same length as center, otherwise a matrix of n rows of random vectors.

References

Gomez, E., Gomez-Villegas, M.A., and Marin, J.M. (1998). A multivariate generalization of the power exponential family of distributions. *Communications in Statistics - Theory and Methods* **27**, 589-600.

Kotz, S., Kozubowski, T.J., and Podgorski, K. (2001). *The Laplace Distributions and Generalizations*. Birkhauser, Boston.

Examples

```
# dispersion parameters
Scatter <- matrix(c(1,.5,.5,1), ncol = 2)
Scatter

# generate the sample
y <- rmLaplace(n = 2000, Scatter = Scatter)

# scatterplot of a random bivariate Laplace sample with center
# vector zero and scale matrix 'Scatter'
par(pty = "s")
plot(y, xlab = "", ylab = "")
title("bivariate Laplace sample", font.main = 1)
```

simulate.lad

Simulate responses from lad models

Description

Simulate one or more responses from the distribution corresponding to a fitted lad object.

Usage

```
## S3 method for class 'lad'
simulate(object, nsim = 1, seed = NULL, ...)
```

Arguments

object	an object representing a fitted model.
nsim	number of response vectors to simulate. Defaults to 1.
seed	an object specifying if and how the random number generator should be initialized ('seeded'). For the "lad" method, either NULL or an integer that will be used in a call to <code>set.seed</code> before simulating the response vectors. If set, the value is saved as the "seed" attribute of the returned value. The default, NULL will not change the random generator state, and return <code>.Random.seed</code> as the "seed" attribute, see 'Value'.
...	additional optional arguments.

Value

For the "lad" method, the result is a data frame with an attribute "seed". If argument seed is NULL, the attribute is the value of `.Random.seed` before the simulation was started.

Author(s)

Tymoteusz Wolodzko and Felipe Osorio

Examples

```
fm <- lad(stack.loss ~ ., data = stackloss)
sm <- simulate(fm, nsim = 4)
```

vcov.lad

Calculate variance-covariance matrix from lad models

Description

Returns the variance-covariance matrix of the main parameters of a fitted model for lad objects. The "main" parameters of model correspond to those returned by `coef`, and typically do not contain a nuisance scale parameter.

Usage

```
## S3 method for class 'lad'
vcov(object, ...)
```

Arguments

object	an object representing a fitted model.
...	additional arguments for method functions.

Value

A matrix of the estimated covariances between the parameter estimates in the linear regression model. This should have row and column names corresponding to the parameter names given by the `coef` method.

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