

# Package ‘wle’

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**Title** Weighted Likelihood Estimation

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**LazyData** yes

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**Description**

Compute a robust analysis of deviance table for one or more generalized linear model fits.

**Usage**

```
## S3 method for class 'wle.glm.root'
anova(object, ..., dispersion = NULL, test = NULL)
```

**Arguments**

object, ...	objects of class <code>wle.glm.root</code> , typically the result of a call to <code>extractRoot.wle.glm</code> , or a list of objects each of which a result of a call to <code>"extractRoot.wle.glm"</code> method.
dispersion	the dispersion parameter for the fitting family. By default it is obtained from the object(s).
test	a character string, (partially) matching one of <code>"Chisq"</code> , <code>"F"</code> or <code>"Cp"</code> . See <a href="#">stat.anova</a> .

**Details**

Specifying a single object gives a sequential analysis of deviance table for that fit. That is, the reductions in the residual deviance as each term of the formula is added in turn are given in as the rows of a table, plus the residual deviances themselves.

If more than one object is specified, the table has a row for the residual degrees of freedom and deviance for each model. For all but the first model, the change in degrees of freedom and deviance is also given. (This only makes statistical sense if the models are nested.) It is conventional to list the models from smallest to largest, but this is up to the user.

The table will optionally contain test statistics (and P values) comparing the reduction in deviance for the row to the residuals. For models with known dispersion (e.g., binomial and Poisson fits) the robust chi-squared test is most appropriate, and for those with dispersion estimated by moments (e.g., gaussian, quasibinomial and quasipoisson fits) the Robust F test is most appropriate. Robust Mallows'  $C_p$  statistic is the residual weighted deviance plus twice the estimate of  $\sigma^2$  times the residual (weighted) degrees of freedom, which is closely related to Robust AIC (and a multiple of it if the dispersion is known).

The dispersion estimate will be taken from the largest model, using the value returned by [summary.wle.glm](#). As this will in most cases use a Chisquared-based estimate, the F tests are not based on the residual deviance in the analysis of deviance table shown.

**Value**

An object of class "anova" inheriting from class "data.frame".

**Warning**

The comparison between two or more models by `anova.wle.glm.root` or `anova.wleglm1ist` will only be valid if they are fitted to the same dataset. This may be a problem if there are missing values and R's default of `na.action = na.omit` is used, and `anova.wleglm1ist` will detect this with an error.

Since in a model selection procedure and/or on an ANOVA table the weights of the WLE procedure must be that of the FULL model (and not that of the actual model) statistics on degrees of freedom, deviance and AIC are valid only if object is the FULL model.

**References**

Agostinelli, C. and Markatou, M. (2001) Test of hypotheses based on the Weighted Likelihood Methodology, *Statistica Sinica*, vol. 11, n. 2, 499-514.

Agostinelli, C. (2002) Robust model selection in regression via weighted likelihood methodology *Statistics and Probability Letters*, 56, 289-300.

Agostinelli, C. and Al-quallaf, F. (2009) Robust inference in Generalized Linear Models. Manuscript in preparation.

Hastie, T. J. and Pregibon, D. (1992) *Generalized linear models*. Chapter 6 of *Statistical Models in S* eds J. M. Chambers and T. J. Hastie, Wadsworth & Brooks/Cole.

**See Also**

[extractRoot.wle.glm](#), [wle.glm](#), [anova](#).

**Examples**

```
## --- Continuing the Example from '?wle.glm':  
  
anova(extractRoot(wle.glm.D93))  
anova(extractRoot(wle.glm.D93), test = "Cp")  
anova(extractRoot(wle.glm.D93), test = "Chisq")
```

---

artificial

*Hawkins, Bradu, Kass's Artificial Data*

---

**Description**

This data set was generated by Hawkins, Bradu and Kass in the 1984 for illustrating some of the merits of a robust technique. The data set consists of 75 observations in four dimensions (one response and three explanatory variables). The first 10 observations are bad leverage points, and the next four points are good leverage points (i.e., their  $x$  are outlying, but the corresponding  $y$  fit the model quite well).

**Usage**

```
data(artificial)
```

**Format**

`artificial` is a data frame with 75 cases (rows) and 4 variables (columns) where the last column is the dependent variable, `y.artificial` and `x.artificial` (as a matrix) are also available.

**Source**

Hawkins, D.M., Bradu, D., and Kass, G.V. (1984) Location of several outliers in multiple regression data using elemental sets. *Technometrics*, **26**, 197–208.

**See Also**

Rousseeuw, P.J., and Leroy, A.M. (1987) *Robust regression and outliers detection*, Wiley.

---

binary	<i>Convert decimal base number to binary base</i>
--------	---

---

**Description**

Convert decimal base number to binary base.

**Usage**

```
binary(x, dim)
```

**Arguments**

x	a number in decimal base.
dim	the number of digits, if missing the right number of digits is evaluated.

**Value**

binary	a vector representing the 'x' number in binary base.
dicotomy	the same as 'binary' but 'TRUE' and 'FALSE' instead of 1 and 0.

**Note**

the elements of 'binary' and 'dicotomy' are in reverse order.

**Author(s)**

Claudio Agostinelli

**Examples**

```
binary(2)
binary(10,dim=5)
```

---

cavendish

*Cavendish's determinations of the mean density of the earth Data*


---

**Description**

The Cavendish's determinations of the mean density of the earth data (relative to that of water) are 29 measures performed in the 1798 using a torsion balance devised earlier by Michell. After the sixth of these determinations, Cavendish changed his experimental apparatus by replacing a suspension wire by one that was stiffer. To further complicate matters, Cavendish erred in taking the mean of all 29 determinations by treating the value 4.88 as if it were in fact 5.88.

**Usage**

```
data(cavendish)
```

**Format**

cavendish is a vector of 29 observations, the first six are made before the apparatus replacement.

**Source**

Cavendish, H. (1900) Experiments to determine the density of the earth, *Philosophical Transactions of the Royal Society of London for the year 1798 (Part II)* **88**, 469–526, Reprinted in *The law of gravitation* (A.S. Mackenzie, ed.) American, New York.

**See Also**

Stigler, S.M. (1977) Do robust estimators work with *real* data? *Annals of Statistics*, **5**, 1055–1098, (with discussion).

---

extractRoot

*Extract a Root from a result of a wle function*


---

**Description**

This function extract the information regarding one solution of the Weighted Likelihood Estimating Equation.

**Usage**

```
## S3 method for class 'wle.glm'
extractRoot(object, root=1, ...)
```

**Arguments**

object	an object of class "wle.glm", usually, a result of a call to <code>wle.glm</code> .
root	an integer number to specify which root should be extract.
...	further arguments passed to or from other methods.

**Value**

`extract.wle.glm` returns an object of class "extract.wle.glm.root", a (variable length) list containing at least the following components:

coefficients	a named vector of coefficients
residuals	the <i>working</i> residuals, that is the residuals in the final iteration of the IWLS fit. Since cases with zero weights are omitted, their working residuals are NA.
fitted.values	the fitted mean values, obtained by transforming the linear predictors by the inverse of the link function.
rank	the numeric rank of the fitted linear model.
family	the <code>family</code> object used.
linear.predictors	the linear fit on link scale.
deviance	up to a constant, minus twice the maximized log-likelihood. Where sensible, the constant is chosen so that a saturated model has deviance zero.
aic	<i>Akaike's An Information Criterion</i> , minus twice the maximized log-likelihood plus twice the number of coefficients (so assuming that the dispersion is known).
null.deviance	The deviance for the null model, comparable with deviance. The null model will include the offset, and an intercept if there is one in the model. Note that this will be incorrect if the link function depends on the data other than through the fitted mean: specify a zero offset to force a correct calculation.
iter	the number of iterations of IWLS used.
weights	the <i>working</i> weights, that is the weights in the final iteration of the IWLS fit.
prior.weights	the weights initially supplied, a vector of 1s if none were.
df.residual	the residual degrees of freedom.
df.null	the residual degrees of freedom for the null model.
y	if requested (the default) the y vector used. (It is a vector even for a binomial model.)
x	if requested, the model matrix.
model	if requested (the default), the model frame.
converged	logical. Was the IWLS algorithm judged to have converged?
boundary	logical. Is the fitted value on the boundary of the attainable values?
wle.weights	final (robust) weights based on the WLE approach.
wle.asymptotic	logicals. If TRUE asymptotic weight based on Anscombe residual is used for the corresponding observation.

In addition, non-empty fits will have components `qr`, `R`, `qraux`, `pivot` and `effects` relating to the final weighted linear fit.

<code>family</code>	the <a href="#">family</a> object used.
<code>call</code>	the matched call.
<code>formula</code>	the formula supplied.
<code>terms</code>	the <a href="#">terms</a> object used.
<code>data</code>	the <code>data</code> argument.
<code>offset</code>	the offset vector used.
<code>control</code>	the value of the <code>control</code> argument used.
<code>method</code>	the name of the fitter function used, currently always <code>"wle.glm.fit"</code> .
<code>contrasts</code>	(where relevant) the contrasts used.
<code>xlevels</code>	(where relevant) a record of the levels of the factors used in fitting.
<code>tot.sol</code>	the number of solutions found.
<code>not.conv</code>	the number of starting points that does not converge after the <code>max.iter</code> (defined using <code>wle.glm.control</code> ) iterations are reached.
<code>na.action</code>	(where relevant) information returned by <a href="#">model.frame</a> on the special handling of NAs.

If a [binomial](#) `wle.glm` model was specified by giving a two-column response, the weights returned by `prior.weights` are the total numbers of cases (factored by the supplied case weights) and the component `y` of the result is the proportion of successes.

### Author(s)

Claudio Agostinelli and Fatemah Al-quallaf

### See Also

[anova.wle.glm.root](#)

### Examples

```
## --- Continuing the Example from '?wle.glm':
anova(extractRoot(wle.glm.D93))
```



---

hald

*Hald Data*


---

### Description

Montgomery and Peck (1982) illustrated variable selection techniques on the Hald cement data and gave several references to other analysis. The response variable  $y$  is the *heat evolved* in a cement mix. The four explanatory variables are ingredients of the mix, i.e.,  $x_1$ : *tricalcium aluminate*,  $x_2$ : *tricalcium silicate*,  $x_3$ : *tetracalcium alumino ferrite*,  $x_4$ : *dicalcium silicate*. An important feature of these data is that the variables  $x_1$  and  $x_3$  are highly correlated ( $\text{corr}(x_1, x_3) = -0.824$ ), as well as the variables  $x_2$  and  $x_4$  (with  $\text{corr}(x_2, x_4) = -0.975$ ). Thus we should expect any subset of  $(x_1, x_2, x_3, x_4)$  that includes one variable from highly correlated pair to do as any subset that also includes the other member.

### Usage

```
data(hald)
```

### Format

`hald` is a matrix with 13 observations (rows) and 5 variables (columns), the first column is the dependent variable. `y.hald` and `x.hald` are also availables.

### Source

Montgomery, D.C., Peck, E.A. (1982) *Introduction to linear regression analysis*, John Wiley, New York.

---

mde.vonmises

*von Mises Minimum Distance Estimates*


---

### Description

Computes the minimum distance estimates for the parameters of a von Mises distribution: the mean direction and the concentration parameter.

### Usage

```
mde.vonmises(x, bw, mu = NULL, kappa = NULL, n = 512,
  from = circular(0), to = circular(2 * pi), lower = NULL,
  upper = NULL, method = "L-BFGS-B", lower.kappa = .Machine$double.eps,
  upper.kappa = Inf, alpha = NULL, p = 2, control.circular = list(), ...)
## S3 method for class 'mde.vonmises'
print(x, digits = max(3, getOption("digits") - 3), ...)
```

**Arguments**

<code>x</code>	a vector. The object is coerced to class <code>circular</code> .
<code>bw</code>	the value of the smoothing parameter.
<code>mu</code>	initial value for the mean direction. Default: maximum likelihood estimate.
<code>kappa</code>	initial value for the concentration parameter. Default: maximum likelihood estimate.
<code>n</code>	number of points used to approximate the density.
<code>from</code>	from which point in the circle the density is approximate.
<code>to</code>	to which point in the circle the density is approximate.
<code>lower</code>	a 2 elements vector passed to <code>optim</code> used to constrained optimization. First element for the mean direction, second element for the concentration.
<code>upper</code>	a 2 elements vector passed to <code>optim</code> used to constrained optimization. First element for the mean direction, second element for the concentration.
<code>method</code>	passed to <code>optim</code> .
<code>lower.kappa</code>	if <code>lower</code> is NULL this parameter is used to constrained optimization for the concentration parameter.
<code>upper.kappa</code>	if <code>upper</code> is NULL this parameter is used to constrained optimization for the concentration parameter.
<code>alpha</code>	if not NULL overrides the value of <code>p</code> . See the next argument <code>p</code> . This is a different parameterization, <code>alpha=-1/2</code> provides Hellinger distance, <code>alpha=-1</code> provides Kullback-Leibler distance and <code>alpha=-2</code> provides Neyman's Chi-Square distance.
<code>p</code>	<code>p=2</code> provides Hellinger distance, <code>p=-1</code> provides Kullback-Leibler distance and <code>p=Inf</code> provides Neyman's Chi-Square distance. It is ignored if <code>alpha</code> is not NULL.
<code>control.circular</code>	the attribute of the resulting object ( <code>mu</code> )
<code>digits</code>	integer indicating the precision to be used.
<code>...</code>	further parameters in <code>print.mde.vonmises</code> .

**Details**

The distance from an estimated density (by the non parametric kernel density estimator) and the model is evaluated by simple rectangular approximation. `optim` is used to performs minimization.

**Value**

Returns a list with the following components:

<code>call</code>	the <code>match.call()</code> .
<code>mu</code>	the estimate of the mean direction.
<code>kappa</code>	the estimate of the concentration parameter.
<code>dist</code>	the distance between the estimated density and the model.

data	the original supplied data converted in radians, clockwise and zero at 0.
x	the 'n' coordinates of the points where the density is estimated.
y	the estimated density values.
k	the density at the model.

**Author(s)**

Claudio Agostinelli

**References**

C. Agostinelli. Robust estimation for circular data. *Computational Statistics & Data Analysis*, 51(12):5867-5875, 2007.

**See Also**

[circular](#), [mle.vonmises](#) and [wle.vonmises](#).

**Examples**

```
set.seed(1234)
x <- c(rvonmises(n=200, mu=circular(0), kappa=10), rvonmises(n=20, mu=circular(pi/2), kappa=20))
res <- mde.vonmises(x, bw=500, mu=circular(0), kappa=10)
res
plot(circular(0), type='n', xlim=c(-1, 1.75), shrink=1.2)
lines(circular(res$x), res$y)
lines(circular(res$x), res$k, col=2)
legend(1,1.5, legend=c('estimated density', 'MDE'), lty=c(1, 1), col=c(1, 2))
```

---

mde.wrappednormal

*Wrapped Normal Minimum Distance Estimates*


---

**Description**

Computes the minimum distance estimates for the parameters of a Wrapped Normal distribution: the mean direction and the concentration parameter (and the scale parameter).

**Usage**

```
mde.wrappednormal(x, bw, mu = NULL, rho = NULL, sd = NULL,
  alpha = NULL, p = 2, tol = 1e-05, n = 512, from = circular(0),
  to = circular(2 * pi), lower = NULL, upper = NULL,
  method = "L-BFGS-B", lower.rho = 1e-06, upper.rho = 1 - 1e-06,
  min.sd = 0.001, K = NULL, min.k = 10, control.circular = list(), ...)
## S3 method for class 'mde.wrappednormal'
print(x, digits = max(3, getOption("digits") - 3), ...)
```

**Arguments**

x	a vector. The object is coerced to class <a href="#">circular</a> .
bw	the value of the smoothing parameter.
mu	initial value for the mean direction. Default: maximum likelihood estimate.
rho	initial value for the concentration parameter. Default: maximum likelihood estimate.
sd	initial value for the standard deviation parameter. This value is used only if rho is NULL. Default: maximum likelihood estimate.
alpha	if not NULL overrides the value of p. See the next argument p. This is a different parameterization, $\alpha=-1/2$ provides Hellinger distance, $\alpha=-1$ provides Kullback-Leibler distance and $\alpha=-2$ provides Neyman's Chi-Square distance.
p	$p=2$ provides Hellinger distance, $p=-1$ provides Kullback-Leibler distance and $p=Inf$ provides Neyman's Chi-Square distance. It is ignored if alpha is not NULL.
tol	the absolute accuracy to be used to achieve convergence of the algorithm. This argument is passed to the function which determined the Maximum Likelihood estimates of the parameters. See <a href="#">mle.wrappednormal</a> .
n	number of points used to approximate the density.
from	from which point in the circle the density is approximate.
to	to which point in the circle the density is approximate.
lower	a 2 elements vector passed to <code>optim</code> used to constrained optimization. First element for the mean direction, second element for the concentration.
upper	a 2 elements vector passed to <code>optim</code> used to constrained optimization. First element for the mean direction, second element for the concentration.
method	passed to <code>optim</code> .
lower.rho	if lower is NULL this parameter is used to constrained optimization for the concentration parameter.
upper.rho	if upper is NULL this parameter is used to constrained optimization for the concentration parameter.
min.sd	minimum value for the sd parameter. This argument is passed to the function which determined the Maximum Likelihood estimates of the parameters. See <a href="#">mle.wrappednormal</a> .
K	number of elements used to approximate the density of the wrapped normal.
min.k	minimum number of elements used to approximate the density of the wrapped normal.
control.circular	the attribute of the resulting object (mu)
digits	integer indicating the precision to be used.
...	further parameters in <code>print.mde.wrappednormal</code> .

**Details**

The distance from an estimated density (by the non parametric kernel density estimator) and the model is evaluated by simple rectangular approximation. `optim` is used to performs minimization.

**Value**

Returns a list with the following components:

<code>call</code>	the <code>match.call()</code> .
<code>mu</code>	the estimate of the mean direction.
<code>rho</code>	the estimate of the concentration parameter.
<code>sd</code>	the estimate of the standard deviation parameter.
<code>dist</code>	the distance between the estimated density and the model.
<code>data</code>	the original supplied data converted in radians, clockwise and zero at 0.
<code>x</code>	the 'n' coordinates of the points where the density is estimated.
<code>y</code>	the estimated density values.
<code>k</code>	the density at the model.

**Author(s)**

Claudio Agostinelli

**References**

C. Agostinelli. Robust estimation for circular data. *Computational Statistics & Data Analysis*, 51(12):5867-5875, 2007.

**See Also**

[circular](#), [mle.wrappednormal](#) and [wle.wrappednormal](#).

**Examples**

```
set.seed(1234)
x <- c(rwrappednormal(n=200, mu=circular(0), sd=0.6),
      rwrappednormal(n=20, mu=circular(pi/2), sd=0.1))
res <- mde.wrappednormal(x, bw=0.08, mu=circular(0), sd=0.6)
res
plot(circular(0), type='n', xlim=c(-1, 1.75), shrink=1.2)
lines(circular(res$x), res$y)
lines(circular(res$x), res$k, col=2)
legend(1,1.5, legend=c('estimated density', 'MDE'), lty=c(1, 1), col=c(1, 2))
```

---

mle.aic *Akaike Information Criterion*

---

**Description**

The Akaike Information Criterion is evaluated for each submodel.

**Usage**

```
mle.aic(formula, data=list(), model=TRUE, x=FALSE,
        y=FALSE, var.full=0, alpha=2, contrasts = NULL,
        se=FALSE, verbose=FALSE)
```

**Arguments**

formula	a symbolic description of the model to be fit. The details of model specification are given below.
data	an optional data frame containing the variables in the model. By default the variables are taken from the environment which <code>mle.aic</code> is called from.
model, x, y	logicals. If TRUE the corresponding components of the fit (the model frame, the model matrix, the response.)
var.full	the value of variance to be used, if 0 the variance estimated from the full model is used.
alpha	the penalized constant.
contrasts	an optional list. See the <code>contrasts.arg</code> of <code>model.matrix.default</code> .
se	logical. if TRUE the returning object contains standard errors for the parameters of every model.
verbose	if TRUE warnings are printed.

**Details**

Models for `mle.aic` are specified symbolically. A typical model has the form `response ~ terms` where `response` is the (numeric) response vector and `terms` is a series of terms which specifies a linear predictor for response. A terms specification of the form `first+second` indicates all the terms in `first` together with all the terms in `second` with duplicates removed. A specification of the form `first:second` indicates the the set of terms obtained by taking the interactions of all terms in `first` with all terms in `second`. The specification `first*second` indicates the *cross* of `first` and `second`. This is the same as `first+second+first:second`.

**Value**

`mle.aic` returns an object of class "mle.aic".

The function `summary` is used to obtain and print a summary of the results. The generic accessor functions `coefficients` and `residuals` extract coefficients and residuals returned by `mle.aic`. The object returned by `mle.aic` are:

aic	the AIC for each submodels
coefficients	the parameters estimator, one row vector for each submodel.
scale	an estimation of the error scale, one value for each submodel.
residuals	the residuals from the estimated model, one column vector for each submodel.
call	the match.call().
contrasts	
xlevels	
terms	the model frame.
model	if model=TRUE a matrix with first column the dependent variable and the remain column the explanatory variables for the full model.
x	if x=TRUE a matrix with the explanatory variables for the full model.
y	if y=TRUE a vector with the dependent variable.
info	not well working yet, if 0 no error occurred.
se	standard errors of the parameters, one row vector for each submodel. Available only if se is TRUE.

### Author(s)

Claudio Agostinelli

### References

Akaike, H., (1973) Information theory and an extension of the maximum likelihood principle, in: B.N. Petrov and F. Csaki, eds., *Proc. 2nd International Symposium of Information Theory*, Akad'emiai Kiad'o, Budapest, 267-281.

### Examples

```
library(wle)

data(hald)

cor(hald)

result <- mle.aic(y.hald~x.hald)

summary(result,num.max=10)
```

---

mle.aic.summaries      *Summaries and methods for mle.aic*

---

## Description

All these functions are [methods](#) for class `mle.aic` or `summary.mle.aic`.

## Usage

```
## S3 method for class 'mle.aic'
summary(object, num.max=20, verbose=FALSE, ...)

## S3 method for class 'mle.aic'
print(x, digits = max(3, getOption("digits") - 3),
      num.max=max(1, nrow(x$aic)), ...)

## S3 method for class 'summary.mle.aic'
print(x, digits = max(3, getOption("digits") - 3), ...)
```

## Arguments

<code>object</code>	an object of class <code>mle.aic</code> .
<code>x</code>	an object of class <code>mle.aic</code> or <code>summary.mle.aic</code> .
<code>num.max</code>	the max number of models should be reported.
<code>digits</code>	number of digits to be used for most numbers.
<code>verbose</code>	if TRUE warnings are printed.
<code>...</code>	additional arguments affecting the summary produced (in <code>summary.mle.aic</code> ) or further arguments passed to or from other methods (in <code>print.mle.aic</code> and <code>print.summary.mle.aic</code> ).

## Value

`summary.mle.aic` returns a list:

<code>aic</code>	the first <code>num.max</code> best models with their AIC.
<code>num.max</code>	the number of models reported.
<code>call</code>	

## Author(s)

Claudio Agostinelli

## See Also

[mle.aic](#) a function for evaluate the Akaike Information Criterion.



mle.cp

*Mallows Cp***Description**

The Mallows Cp is evaluated for each submodel.

**Usage**

```
mle.cp(formula, data=list(), model=TRUE, x=FALSE,
        y=FALSE, var.full=0, contrasts=NULL, verbose=FALSE)
```

**Arguments**

formula	a symbolic description of the model to be fit. The details of model specification are given below.
data	an optional data frame containing the variables in the model. By default the variables are taken from the environment which <code>mle.cp</code> is called from.
model, x, y	logicals. If TRUE the corresponding components of the fit (the model frame, the model matrix, the response).
var.full	the value of variance to be used in the denominator of the Mallows Cp, if 0 the variance estimated from the full model is used.
contrasts	an optional list. See the <code>contrasts.arg</code> of <code>model.matrix.default</code> .
verbose	if TRUE warnings are printed.

**Details**

Models for `mle.cp` are specified symbolically. A typical model has the form `response ~ terms` where `response` is the (numeric) response vector and `terms` is a series of terms which specifies a linear predictor for response. A terms specification of the form `first+second` indicates all the terms in `first` together with all the terms in `second` with duplicates removed. A specification of the form `first:second` indicates the set of terms obtained by taking the interactions of all terms in `first` with all terms in `second`. The specification `first*second` indicates the *cross* of `first` and `second`. This is the same as `first+second+first:second`.

**Value**

`mle.cp` returns an object of class "mle.cp".

The function `summary` is used to obtain and print a summary of the results, only models below the bisector are reported. The generic accessor functions `coefficients` and `residuals` extract coefficients and residuals returned by `mle.cp`. The object returned by `mle.cp` are:

cp	Mallows Cp for each submodels
coefficients	the parameters estimator, one row vector for eac submodel.
scale	an estimation of the error scale, one value for each submodel.

residuals	the residuals from the estimated model, one column vector for each submodel.
call	the match.call().
contrasts	
xlevels	
terms	the model frame.
model	if model=TRUE a matrix with first column the dependent variable and the remain column the explanatory variables for the full model.
x	if x=TRUE a matrix with the explanatory variables for the full model.
y	if y=TRUE a vector with the dependent variable.
info	not well working yet, if 0 no error occurred.

**Author(s)**

Claudio Agostinelli

**References**

Mallows, C.L., (1973) Some comments on Cp, *Technometrics*, 15, 661-675.

**Examples**

```
library(wle)
data(hald)
cor(hald)
result <- mle.cp(y.hald~x.hald)
summary(result)
plot(result)
```

---

mle.cp.summaries

*Summaries and methods for mle.cp*

---

**Description**

All these functions are [methods](#) for class `mle.cp` or `summary.mle.cp`.

**Usage**

```
## S3 method for class 'mle.cp'  
summary(object, num.max=20, verbose=FALSE, ...)  
  
## S3 method for class 'mle.cp'  
print(x, digits = max(3, getOption("digits") - 3),  
      num.max=max(1, nrow(x$cp)), ...)  
  
## S3 method for class 'summary.mle.cp'  
print(x, digits = max(3, getOption("digits") - 3), ...)
```

**Arguments**

object	an object of class <code>mle.cp</code> .
x	an object of class <code>mle.cp</code> or <code>summary.mle.cp</code> .
digits	number of digits to be used for most numbers.
num.max	the max number of models should be reported.
verbose	if TRUE warnings are printed.
...	additional arguments affecting the summary produced (in <code>summary.mle.cp</code> ) or further arguments passed to or from other methods (in <code>print.mle.cp</code> and <code>print.summary.mle.cp</code> ).

**Value**

`summary.mle.cp` returns a list:

cp	the first <code>num.max</code> best models with their Mallows Cp.
num.max	the number of models reported.
call	

**Author(s)**

Claudio Agostinelli

**See Also**

[mle.cp](#) a function for evaluate the Mallows Cp.

---

mle.cv

*Cross Validation Selection Method*


---

### Description

The Cross Validation selection method is evaluated for each submodel.

### Usage

```
mle.cv(formula, data=list(), model=TRUE, x=FALSE,
        y=FALSE, monte.carlo=500, split,
        contrasts=NULL, verbose=FALSE)
```

### Arguments

formula	a symbolic description of the model to be fit. The details of model specification are given below.
data	an optional data frame containing the variables in the model. By default the variables are taken from the environment which <code>mle.cv</code> is called from.
model, x, y	logicals. If TRUE the corresponding components of the fit (the model frame, the model matrix, the response.)
monte.carlo	the number of Monte Carlo replication we use to estimate the average prediction error.
split	the size of the construction sample. When the suggested value is outside the possible range, the split size is let equal to $\max(\text{round}(\text{size}^{(3/4)}), \text{nvar} + 2)$ .
contrasts	an optional list. See the <code>contrasts.arg</code> of <code>model.matrix.default</code> .
verbose	if TRUE warnings are printed.

### Details

Models for `mle.cv` are specified symbolically. A typical model has the form `response ~ terms` where `response` is the (numeric) response vector and `terms` is a series of terms which specifies a linear predictor for response. A terms specification of the form `first+second` indicates all the terms in `first` together with all the terms in `second` with duplicates removed. A specification of the form `first:second` indicates the the set of terms obtained by taking the interactions of all terms in `first` with all terms in `second`. The specification `first*second` indicates the *cross* of `first` and `second`. This is the same as `first+second+first:second`.

### Value

`mle.cv` returns an object of class "mle.cv".

The function `summary` is used to obtain and print a summary of the results.

The object returned by `mle.cv` are:

cv	the estimated prediction error for each submodels
----	---

call	the match.call().
contrasts	
xlevels	
terms	the model frame.
model	if model=TRUE a matrix with first column the dependent variable and the remain column the explanatory variables for the full model.
x	if x=TRUE a matrix with the explanatory variables for the full model.
y	if y=TRUE a vector with the dependent variable.
info	not well working yet, if 0 no error occurred.

**Author(s)**

Claudio Agostinelli

**References**

Shao, J., (1993) Linear model selection by Cross-Validation. *Journal American Statistical Association*, 88, 486-494.

**Examples**

```
library(wle)

data(hald)

cor(hald)

result <- mle.cv(y.hald~x.hald)

summary(result)
```

---

mle.cv.summaries      *Summaries and methods for mle.cv*

---

**Description**

All these functions are [methods](#) for class `mle.cv` or `summary.mle.cv`.

**Usage**

```
## S3 method for class 'mle.cv'
summary(object, num.max=20, verbose=FALSE, ...)

## S3 method for class 'mle.cv'
print(x, digits = max(3, getOption("digits") - 3), num.max=max(1, nrow(x$cv)), ...)

## S3 method for class 'summary.mle.cv'
print(x, digits = max(3, getOption("digits") - 3), ...)
```

**Arguments**

object	an object of class <code>mle.cv</code> .
x	an object of class <code>mle.cv</code> or <code>summary.mle.cv</code> .
digits	number of digits to be used for most numbers.
num.max	the max number of models should be reported.
verbose	if TRUE warnings are printed.
...	additional arguments affecting the summary produced (in <code>summary.mle.cv</code> ) or further arguments passed to or from other methods (in <code>print.mle.cv</code> and <code>print.summary.mle.cv</code> ).

**Value**

`summary.mle.cv` returns a list:

cv	the first <code>num.max</code> best models with their estimated prediction error using CV.
num.max	the number of models reported.
call	

**Author(s)**

Claudio Agostinelli

**See Also**

[mle.cv](#) a function for evaluate the Cross-Validation selection criterion for linear models.

---

mle.stepwise

*Stepwise, Backward and Forward selection methods*

---

**Description**

This function performs Stepwise, Forward and Backward model selection.

**Usage**

```
mle.stepwise(formula, data=list(), model=TRUE, x=FALSE,
              y=FALSE, type="Forward", f.in=4.0, f.out=4.0,
              contrans=NULL, verbose=FALSE)
```

**Arguments**

formula	a symbolic description of the model to be fit. The details of model specification are given below.
data	an optional data frame containing the variables in the model. By default the variables are taken from the environment which <code>mle.stepwise</code> is called from.
model, x, y	logicals. If TRUE the corresponding components of the fit (the model frame, the model matrix, the response.)
type	type="Stepwise": the stepwise methods is used, type="Forward": the forward methods is used, type="Backward": the backward method is used.
f.in	the in value
f.out	the out value
contranst	an optional list. See the <code>contrasts.arg</code> of <code>model.matrix.default</code> .
verbose	if TRUE warnings are printed.

**Details**

Models for `mle.stepwise` are specified symbolically. A typical model has the form `response ~ terms` where `response` is the (numeric) response vector and `terms` is a series of terms which specifies a linear predictor for response. A terms specification of the form `first+second` indicates all the terms in `first` together with all the terms in `second` with duplicates removed. A specification of the form `first:second` indicates the the set of terms obtained by taking the interactions of all terms in `first` with all terms in `second`. The specification `first*second` indicates the *cross* of `first` and `second`. This is the same as `first+second+first:second`.

**Value**

`mle.stepwise` returns an object of class "mle.stepwise".

The function `summary` is used to obtain and print a summary of the results.

The object returned by `mle.stepwise` are:

step	the selected models
type	the type o model selection procedure was used.
f.in	the value of f.in used.
f.out	the value of f.out used.
call	the <code>match.call()</code> .
contrasts	
xlevels	
terms	the model frame.
model	if <code>model=TRUE</code> a matrix with first column the dependent variable and the remain column the explanatory variables for the full model.
x	if <code>x=TRUE</code> a matrix with the explanatory variables for the full model.
y	if <code>y=TRUE</code> a vector with the dependent variable.
info	not well working yet, if 0 no error occurred.

**Author(s)**

Claudio Agostinelli

**References**

Beale, E.M.L., Kendall, M.G., Mann, D.W., (1967) The discarding of variables in multivariate analysis, *Biometrika*, 54, 357-366.

Efroymsen, (1960) Multiple regression analysis, in *Mathematical Methods for Digital Computers*, eds. A. Ralston and H.S. Wilf, 191-203, Wiley, New York.

Garside, M.J., (1965) The best sub-set in multiple regression analysis, *Applied Statistics*, 14, 196-200.

Goldberger, A.S, and Jochems, D.B., (1961) Note on stepwise least squares, *Journal of the American Statistical Association*, 56, 105-110.

Goldberger, A.S., (1961) Stepwise least squares: Residual analysis and specification error, *Journal of the American Statistical Association*, 56, 998-1000.

**Examples**

```
library(wle)

data(hald)

cor(hald)

result <- mle.stepwise(y.hald~x.hald)

summary(result)
```

---

```
mle.stepwise.summaries
```

*Accessing summaries for mle.stepwise*

---

**Description**

All these functions are [methods](#) for class `mle.stepwise` or `summary.mle.stepwise`.

**Usage**

```
## S3 method for class 'mle.stepwise'
summary(object, num.max=20, verbose=FALSE, ...)

## S3 method for class 'mle.stepwise'
print(x, digits = max(3, getOption("digits") - 3), num.max=max(1,nrow(x$step)), ...)

## S3 method for class 'summary.mle.stepwise'
print(x, digits = max(3, getOption("digits") - 3), ...)
```



**Arguments**

object	an object of class <code>mle.stepwise</code> .
x	an object of class <code>mle.stepwise</code> or <code>summary.mle.stepwise</code> .
digits	number of digits to be used for most numbers.
num.max	the number of the last iterations reported.
verbose	if TRUE warnings are printed.
...	additional arguments affecting the summary produced (in <code>summary.mle.stepwise</code> ) or further arguments passed to or from other methods (in <code>print.mle.stepwise</code> and <code>print.summary.mle.stepwise</code> ).

**Value**

The function `summary.mle.stepwise` returns the last `num.max` iterations, call plus:

step	the model for each iteration reported.
num.max	the number of iterations reported.
type	the type of selection procedure used.
f.in	the in value
f.out	the out value

**Author(s)**

Claudio Agostinelli

---

plot.mle.cp

*Plot the Mallows Cp*

---

**Description**

Plot the Mallows Cp.

**Usage**

```
## S3 method for class 'mle.cp'
plot(x, base.line=0, num.max=20,
      plot.it=TRUE, log.scale=FALSE,
      xlab="Number of Predictors", ylab=NULL,
      verbose=FALSE, ...)
```

**Arguments**

<code>x</code>	an object of class <code>mle.cp</code> .
<code>base.line</code>	the intercept of the line to split the submodels in acceptable (good) and not-acceptable (bad), (the slope is always one).
<code>num.max</code>	maximum number of submodels plotted.
<code>plot.it</code>	if TRUE the graph is plotted.
<code>log.scale</code>	if TRUE the y-axis as log10 scale.
<code>xlab</code>	a title for the x axis.
<code>ylab</code>	a title for the y axis.
<code>verbose</code>	if TRUE warnings are printed.
<code>...</code>	graphical parameters can be given as arguments.

**Value**

<code>num.good</code>	number of submodels below the <code>base.line</code>
<code>num.bad</code>	number of submodels above the <code>base.line</code>
<code>cp.good</code>	list of the submodels below the <code>base.line</code> with their <code>Cp</code> .
<code>cp.bad</code>	list of the submodels above the <code>base.line</code> with their <code>Cp</code> .

**Author(s)**

Claudio Agostinelli

**See Also**

[mle.cp](#) a function to calculate the Mallows `Cp`.

**Examples**

```
library(wle)

data(hald)

result <- mle.cp(y.hald~x.hald)

plot(result,num.max=7)
```

---

plot.wle.cp

*Plot the Weighted Mallows Cp*


---

**Description**

Plot the weighted Mallows Cp based on weighted likelihood.

**Usage**

```
## S3 method for class 'wle.cp'
plot(x, base.line=0, num.max=20,
      plot.it=TRUE, log.scale=FALSE,
      xlab="Number of Predictors", ylab=NULL,
      verbose=FALSE, ...)
```

**Arguments**

x	an object of class wle.cp.
base.line	the intercept of the line to split the submodels in acceptable (good) and not-acceptable (bad), (the slope is always one).
num.max	maximum number of submodels plotted.
plot.it	if TRUE the graph is plotted.
log.scale	if TRUE the y-axis as log10 scale.
xlab	a title for the x axis.
ylab	a title for the y axis.
verbose	if TRUE warnings are printed.
...	graphical parameters can be given as arguments.

**Value**

num.good	number of submodels below the base.line
num.bad	number of submodels above the base.line
wcp.good	list of the submodels below the base.line with their WCp.
wcp.bad	list of the submodels above the base.line with their WCp.

**Author(s)**

Claudio Agostinelli

## References

Agostinelli, C., (1999) Robust model selection in regression via weighted likelihood methodology, *Working Paper n. 1999.4*, Department of Statistics, University of Padova.

Agostinelli, C., (1999) Robust model selection in regression via weighted likelihood methodology, submitted to *Statistics & Probability Letters*, revised december 1999.

Agostinelli, C., (1998) Inferenza statistica robusta basata sulla funzione di verosimiglianza pesata: alcuni sviluppi, *Ph.D Thesis*, Department of Statistics, University of Padova.

Agostinelli, C., (1998) Verosimiglianza pesata nel modello di regressione lineare, *XXXIX Riunione scientifica della Societ a Italiana di Statistica*, Sorrento 1998.

## See Also

[wle.cp](#) a function to calculate the Weighted Mallows Cp, [wle.lm](#) a function for estimating linear models with normal distribution error and normal kernel.

## Examples

```
library(wle)
x.data <- c(runif(60,20,80),runif(5,73,78))
e.data <- rnorm(65,0,0.6)
y.data <- 8*log(x.data+1)+e.data
y.data[61:65] <- y.data[61:65]-4
z.data <- c(rep(0,60),rep(1,5))
plot(x.data, y.data, xlab="X", ylab="Y")
xx.data <- cbind(x.data, x.data^2, x.data^3, log(x.data+1))
result <- wle.cp(y.data~xx.data)
plot(result,num.max=15)
```

---

plot.wle.lm

*Plots for the Linear Model*

---

## Description

The `plot.wle.lm` function plots a separate graph windows for each root. In each windows four plots are printed: residuals vs fitted, normal qq plot of the residuals, weighted residuals vs weighted fitted, normal qq plot of the weighted residuals. A summary plot is also printed: in the diagonal, the value of the weights vs position of the observations for each root; in the upper diagonal residuals vs residuals of two different roots; in the lower diagonal weights vs weights of two different roots. The roots and the graphs can be chosen by the arguments `roots`, `which.main` and `which`.

## Usage

```
## S3 method for class 'wle.lm'
plot(x, roots, which=1:4, which.main, level.weight=0.5,
     ask = dev.interactive(), col=c(2, 1, 3), id.n=3, labels.id,
     cex.id = 0.75, verbose=FALSE, ...)
```

**Arguments**

x	an object of class <code>wle.lm</code> .
roots	a vector specify for which roots the plots are required.
which	if a subset of the plots for each root is required, specify a subset of the numbers $0:4$ , 0 means no plots.
which.main	if a subset of the plots for the main graphic is required, specify a subset of the numbers $0:\text{roots}^2$ , 0 means no plots. The plots are specified by columns.
level.weight	value of the weight under which an observations is marked with different color.
ask	logical; if TRUE, the user is <i>asked</i> before each plot, see <code>par(ask=.)</code> .
col	a vector of 3 elements, to specify colors for the plots.
id.n	number of points to be labelled in some plots, starting with the ones with less weight.
labels.id	vector of labels, from which the labels for less weighted points will be chosen. If missing uses observation numbers.
cex.id	magnification of point labels.
verbose	if TRUE warnings are printed.
...	graphical parameters can be given as arguments.

**Author(s)**

Claudio Agostinelli

**See Also**

[wle.lm](#) a function for estimating linear models with normal distribution error and normal kernel.

**Examples**

```
library(wle)

data(artificial)

result <- wle.lm(y~x1+x2+x3, data=artificial, boot=40, group=6, num.sol=2)

result

plot(result) # all plots, default behavior

plot(result, roots=1) # only first root, one plot for window

par(mfcol=c(2,2))
plot(result, roots=1) # only first root, as usual

plot(result, roots=2, which=1, which.main=0)
# only second root, only residual vs fitted values plot

plot(result, which=1)
```

```
# main plot + residual vs fitted values plot for each root

par(mfcol=c(3,2))
plot(result, which=1)
# main plot + residual vs fitted values plot for each root all in the same window
```

---

residualsAnscombe      *Anscombe residuals*

---

### Description

Evaluate the Anscombe residuals for a given type of `family` in GLM.

### Usage

```
residualsAnscombe(y, mu, family, ...)
```

### Arguments

`y`                      vector of the response variable  
`mu`                      vector of the same length as `y` with the corresponding fitted values.  
`family`                an object of class `family`.  
`...`                    not used yet.

### Details

The function performs the Anscombe transformation to obtain residuals that are asymptotically normal distributed. For the Binomial family (see Con and Snell 1968) the transformation is

$$\text{beta}(2/3, 2/3) * (\text{pbeta}(y/m, 2/3, 2/3) - \text{pbeta}(\mu - (1 - 2 * \mu) / (6 * m), 2/3, 2/3)) / ((\mu^{1/6}) * (1 - \mu)^{1/6}) / \text{sqrt}(m)$$

where `m` is the number of trial and `y` the number of successes. For the Poisson family (see Con and Snell 1968) the transformation is

$$(3/2 * (y^{2/3}) - (\mu - 1/6)^{2/3}) / (\mu^{1/6})$$

while for the Gamma family (see McCullagh and Nelder 1989) the transformation is

$$3 * (y^{1/3}) - \mu^{1/3} / (\mu^{1/3})$$

and for the Inverse Gaussian family (see McCullagh and Nelder 1989) the transformation is

$$(\ln(y) - \ln(\mu)) / \sqrt{\mu}$$

### Value

It return a vector with the Anscombe residuals.

**Author(s)**

Claudio Agostinelli and Fatemah Al-quallaf

**References**

Agostinelli, C. and Al-quallaf, F. (2009) Robust inference in Generalized Linear Models. Manuscript in preparation.

D. R. Cox and E. J. Snell. A general definition of residuals. *Journal of the Royal Statistical Society. Series B (Methodological)*, 30(2):248-275, 1968.

R. M. Loynes. On cox and snell's general definition of residuals. *Journal of the Royal Statistical Society. Series B (Methodological)*, 31(1):103-106, 1969.

D. A. Pierce and D. W. Schafer. Residuals in generalized linear models. *Journal of the American Statistical Association*, 81(396):977-986, 1986.

Rollin Brant. Residual components in generalized linear models. *The Canadian Journal of Statistics*, 15(2):115-126, 1987.

**See Also**

[wle.glm](#)

**Examples**

```
## Dobson (1990) Page 93: Randomized Controlled Trial :
counts <- c(18,17,15,20,10,20,25,13,12)
outcome <- gl(3,1,9)
treatment <- gl(3,3)
print(d.AD <- data.frame(treatment, outcome, counts))
wle.glm.D93 <- wle.glm(counts ~ outcome + treatment, family=poisson())
res <- residualsAnscombe(counts, mu=wle.glm.D93$root1$fitted.values, family=poisson())
qqnorm(res)
qqline(res)
```

---

rocky

*Rockwell hardness, 100 coils produced in sequence at a Chicago Steel Mill Data*

---

**Description**

This data set is the Rockwell Hardness (measured on Rockwell "B" scale) of a sample of 100 steel coins produced in sequence in a Chicago steel mill.

**Usage**

`data(rocky)`

**Format**

rocky is a time serie with 100 elements.

**Source**

K.W. Hipel and A.I. McLeod (1994), Time series modelling of water resources and environmental systems, Elsevier, Amsterdam.

---

selection

*Selection's Data*

---

**Description**

This data set consists of 60 observations for two variables (ydata and xdata). The appropriate regression model for the first fiftyfourth observations should be like  $y = x^2 + e$  the last sixth comes from a different model.

**Usage**

```
data(selection)
```

**Format**

xdata is a vector which contains  $x, x^2$ , while ydata contains the dependent variable's observations.

**Source**

Agostinelli, C. (1999) Robust model selection in regression via weighted likelihood methodology, submitted to *Statistics & Probability Letters*, revised december 1999.

---

summary.wle.glm

*Summarizing Generalized Linear Model Robust Fits*

---

**Description**

These functions are all [methods](#) for class wle.glm or summary.wle.glm objects.

**Usage**

```
## S3 method for class 'wle.glm'
summary(object, root = 1, dispersion = NULL,
         correlation = FALSE, symbolic.cor = FALSE, ...)

## S3 method for class 'summary.wle.glm'
print(x, digits = max(3, getOption("digits") - 3),
      symbolic.cor = x$symbolic.cor,
      signif.stars = getOption("show.signif.stars"), ...)
```



**Arguments**

object	an object of class "wle.glm", usually, a result of a call to <a href="#">wle.glm</a> .
root	an integer number to specify for which root the summary should be reported.
x	an object of class "summary.wle.glm", usually, a result of a call to <a href="#">summary.glm</a> .
dispersion	the dispersion parameter for the family used. Either a single numerical value or NULL (the default), when it is inferred from object (see 'Details').
correlation	logical; if TRUE, the correlation matrix of the estimated parameters is returned and printed.
digits	the number of significant digits to use when printing.
symbolic.cor	logical. If TRUE, print the correlations in a symbolic form (see <a href="#">symnum</a> ) rather than as numbers.
signif.stars	logical. If TRUE, 'significance stars' are printed for each coefficient.
...	further arguments passed to or from other methods.

**Details**

`print.summary.wle.glm` tries to be smart about formatting the coefficients, standard errors, etc. and additionally gives 'significance stars' if `signif.stars` is TRUE. The `coefficients` component of the result gives the estimated coefficients and their estimated standard errors, together with their ratio. This third column is labelled `t_ratio` if the dispersion is estimated, and `z_ratio` if the dispersion is known (or fixed by the family). A fourth column gives the two-tailed p-value corresponding to the t or z ratio based on a Student t or Normal reference distribution. (It is possible that the dispersion is not known and there are no residual degrees of freedom from which to estimate it. In that case the estimate is NaN.)

Aliased coefficients are omitted in the returned object but restored by the `print` method.

Correlations are printed to two decimal places (or symbolically): to see the actual correlations print `summary(object)$correlation` directly.

The dispersion of a GLM is not used in the fitting process, but it is needed to find standard errors. If dispersion is not supplied or NULL, the dispersion is taken as 1 for the binomial and Poisson families, and otherwise estimated by the residual Chisquared statistic (calculated from cases with non-zero weights) divided by the residual degrees of freedom.

`summary` can be used with Gaussian `wle.glm` fits to handle the case of a linear regression with known error variance, something not handled by [summary.wle.lm](#).

**Value**

`summary.wle.glm` returns an object of class "summary.wle.glm", a list with components

call	the component from object.
family	the component from object.
deviance	the component from object.
contrasts	the component from object.
df.residual	the component from object.

null.deviance	the component from object.
df.null	the component from object.
deviance.resid	the deviance residuals: see <a href="#">residuals.glm</a> .
coefficients	the matrix of coefficients, standard errors, z-values and p-values. Aliased coefficients are omitted.
aliased	named logical vector showing if the original coefficients are aliased.
dispersion	either the supplied argument or the inferred/estimated dispersion if the latter is NULL.
df	a 3-vector of the rank of the model and the number of residual degrees of freedom, plus number of non-aliased coefficients.
cov.unscaled	the unscaled (dispersion = 1) estimated covariance matrix of the estimated coefficients.
cov.scaled	ditto, scaled by dispersion.
correlation	(only if correlation is true.) The estimated correlations of the estimated coefficients.
symbolic.cor	(only if correlation is true.) The value of the argument symbolic.cor.

### Warnings

Since in a model selection procedure and/or on an ANOVA table the weights of the WLE procedure must be that of the FULL model (and not that of the actual model) statistics on degrees of freedom, deviance and AIC are valid only if this is the FULL model.

### See Also

[wle.glm](#), [summary](#).

### Examples

```
## --- Continuing the Example from '?wle.glm':
summary(wle.glm.D93)
```

---

wle.aic

*Weighted Akaike Information Criterion*

---

### Description

The Weighted Akaike Information Criterion.

### Usage

```
wle.aic(formula, data=list(), model=TRUE, x=FALSE,
         y=FALSE, boot=30, group, var.full=0, num.sol=1,
         raf="HD", smooth=0.031, tol=10^(-6),
         equal=10^(-3), max.iter=500, min.weight=0.5,
         method="full", alpha=2, contrasts=NULL, verbose=FALSE)
```

**Arguments**

formula	a symbolic description of the model to be fit. The details of model specification are given below.
data	an optional data frame containing the variables in the model. By default the variables are taken from the environment which <code>wle.aic</code> is called from.
model, x, y	logicals. If TRUE the corresponding components of the fit (the model frame, the model matrix, the response.)
boot	the number of starting points based on bootstrap subsamples to use in the search of the roots.
group	the dimension of the bootstrap subsamples. The default value is $\max(\text{round}(\text{size}/4), \text{var})$ where <i>size</i> is the number of observations and <i>var</i> is the number of variables.
var.full	the value of variance to be used in the denominator of the WAIC, if 0 the variance estimated from the full model is used.
num.sol	maximum number of roots to be searched.
raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
smooth	the value of the smoothing parameter.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.
equal	the absolute value for which two roots are considered the same. (This parameter must be greater than <code>tol</code> ).
max.iter	maximum number of iterations.
min.weight	see details.
method	see details.
alpha	penalty value.
contrasts	an optional list. See the <code>contrasts.arg</code> of <code>model.matrix.default</code> .
verbose	if TRUE warnings are printed.

**Details**

Models for `wle.aic` are specified symbolically. A typical model has the form `response ~ terms` where `response` is the (numeric) response vector and `terms` is a series of terms which specifies a linear predictor for response. A terms specification of the form `first+second` indicates all the terms in `first` together with all the terms in `second` with duplicates removed. A specification of the form `first:second` indicates the the set of terms obtained by taking the interactions of all terms in `first` with all terms in `second`. The specification `first*second` indicates the *cross* of `first` and `second`. This is the same as `first+second+first:second`.

`min.weight`: the weighted likelihood equation could have more than one solution. These roots appear for particular situation depending on contamination level and type. The presence of multiple roots in the full model can create some problem in the set of weights we should use. Actually, the selection of the root is done by the minimum scale error provided. Since this choice is not always

the one would choose, we introduce the `min.weight` parameter in order to choose only between roots that do not down weight everything. This is not still the optimal solution, and perhaps, in the new release, this part will be change.

`method`: this parameter, when set to "reduced", allows to use weights based on the reduced model. This is strongly discourage since the robust and asymptotic property of this kind of weighted AIC are not as good as the one based on `method="full"`.

## Value

`wle.aic` returns an object of `class` "wle.aic".

The function `summary` is used to obtain and print a summary of the results. The generic accessor functions `coefficients` and `residuals` extract coefficients and residuals returned by `wle.aic`. The object returned by `wle.aic` are:

<code>waic</code>	Weighted Akaike Information Criterion for each submodels
<code>coefficients</code>	the parameters estimator, one row vector for each root found and each submodel.
<code>scale</code>	an estimation of the error scale, one value for each root found and each submodel.
<code>residuals</code>	the unweighted residuals from the estimated model, one column vector for each root found and each submodel.
<code>tot.weights</code>	the sum of the weights divide by the number of observations, one value for each root found and each submodel.
<code>weights</code>	the weights associated to each observation, one column vector for each root found and each submodel.
<code>freq</code>	the number of starting points converging to the roots.
<code>call</code>	the <code>match.call()</code> .
<code>contrasts</code>	
<code>xlevels</code>	
<code>terms</code>	the model frame.
<code>model</code>	if <code>model=TRUE</code> a matrix with first column the dependent variable and the remain column the explanatory variables for the full model.
<code>x</code>	if <code>x=TRUE</code> a matrix with the explanatory variables for the full model.
<code>y</code>	if <code>y=TRUE</code> a vector with the dependent variable.
<code>info</code>	not well working yet, if 0 no error occurred.

## Author(s)

Claudio Agostinelli

## References

Agostinelli, C., (1999) Robust model selection in regression via weighted likelihood methodology, *Working Paper n. 1999.4*, Department of Statistics, University of Padova.

Agostinelli, C., (2002) Robust model selection in regression via weighted likelihood methodology, *Statistics & Probability Letters*, 56, 289-300.

Agostinelli, C., (1998) Inferenza statistica robusta basata sulla funzione di verosimiglianza pesata: alcuni sviluppi, *Ph.D Thesis*, Department of Statistics, University of Padova.

Agostinelli, C., Markatou, M., (1998) A one-step robust estimator for regression based on the weighted likelihood reweighting scheme, *Statistics & Probability Letters*, Vol. 37, n. 4, 341-350.

Agostinelli, C., (1998) Verosimiglianza pesata nel modello di regressione lineare, *XXXIX Riunione scientifica della Societa' Italiana di Statistica*, Sorrento 1998.

## Examples

```
library(wle)

x.data <- c(runif(60,20,80),runif(5,73,78))
e.data <- rnorm(65,0,0.6)
y.data <- 8*log(x.data+1)+e.data
y.data[61:65] <- y.data[61:65]-4
z.data <- c(rep(0,60),rep(1,5))

plot(x.data,y.data,xlab="X",ylab="Y")

xx.data <- cbind(x.data,x.data^2,x.data^3,log(x.data+1))
colnames(xx.data) <- c("X","X^2","X^3","log(X+1)")

result <- wle.aic(y.data~xx.data,boot=10,group=10,num.sol=2)

summary(result)

result <- wle.aic(y.data~xx.data+z.data,boot=10,group=10,num.sol=2)

summary(result)
```

## Description

The function evaluate the Weighted Akaike Information Criterion for AutoRegressive Models. This is a robust model selection method to choose the order of an AutoRegressive model.

**Usage**

```
wle.aic.ar(x, order = c(1, 0), seasonal = list(order = c(0, 0),
  period = NA), group, group.start, group.step = group.start,
  xreg = NULL, include.mean = TRUE, na.action = na.fail,
  tol = 10-6, tol.step = tol, equal = 10-3, equal.step = equal,
  raf = "HD", var.full = 0, smooth = 0.0031, smooth.a0 = smooth,
  boot = 10, boot.start = 10, boot.step = boot.start, num.sol = 1,
  x.init = 0, x.seasonal.init = 0, max.iter.out = 20, max.iter.in = 50,
  max.iter.start = 200, max.iter.step = 500, verbose = FALSE,
  w.level = 0.4, min.weights = 0.5, population.size = 10,
  population.choose = 5, elements.random = 2, wle.start = FALSE,
  init.values = NULL, num.max = NULL, num.sol.step = 2,
  min.weights.aic = 0.5, approx.w = TRUE, ask = TRUE, alpha = 2,
  method = "WLS")
```

**Arguments**

x	a univariate time series.
order	maximum order to investigate. A specification of the non-seasonal part of the ARI model: the two components (p,d) are the AR order and the degree of differencing.
seasonal	a specification of the seasonal part of the ARI model, plus the period (which defaults to frequency(x)).
group	the dimension of the bootstrap subsamples.
group.start	the dimension of the bootstrap subsamples used in the starting process if wle.init=TRUE.
group.step	the dimension of the bootstrap subsamples used in a step, it must be less than group.
xreg	optionally, a vector or matrix of external regressors, which must have the same number of rows as x.
include.mean	Should the ARI model include a mean term? The default is TRUE for undifferenced series, FALSE for differenced ones (where a mean would not affect the fit nor predictions).
na.action	function to be applied to remove missing values.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.
tol.step	the absolute accuracy to be used to achieve convergence in a step.
equal	the absolute value for which two roots are considered the same. (This parameter must be greater than tol).
equal.step	the absolute value for which two roots are considered the same in a step. (This parameter must be greater than tol.step).
raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
var.full	An estimate of the residual variance for the full model.
smooth	the value of the smoothing parameter.

smooth.ao	the value of the smoothing parameter used in the outliers classification, default equal to smooth.
boot	the number of starting points based on bootstrap subsamples to use in the search of the roots.
boot.start	the number of starting points based on bootstrap subsamples to use in the search of the roots in the starting process.
boot.step	the number of starting points based on bootstrap subsamples to use in the search of the roots in a step.
num.sol	maximum number of roots to be searched.
x.init	initial values, a vector with the same length of the AR order, or a number, default is 0.
x.seasonal.init	initial values, a vector with the same length of the SAR order, or a number, default is 0.
max.iter.out	maximum number of iterations in the outer loop.
max.iter.in	maximum number of iterations in the inner loop.
max.iter.start	maximum number of iterations in the starting process.
max.iter.step	maximum number of iterations in a step.
verbose	if TRUE warnings are printed.
w.level	the threshold used to decide if an observation could be an additive outlier.
min.weights	see details.
population.size	see details.
population.choose	see details.
elements.random	see details.
wle.start	if TRUE a weighted likelihood estimation is used to have a starting value.
init.values	a vector with initial values for the AR and seasonal AR coefficients and the innovations variance.
num.max	maximum number of observations can be considered as possible additive outliers.
num.sol.step	maximum number of roots to be searched in a step.
min.weights.aic	see details.
approx.w	logical: if TRUE an approximation is used to evaluate the weights in the outlier identification procedure.
ask	logical. If TRUE, in the case of multiple roots in the full model, the users is asked for selecting the root.
alpha	penalty value.
method	if "WLE" the parameters are estimated using weighted likelihood estimating equations in the reduced models, otherwise if "WLS" a weighted least squares approach is used with weights based on the full model.

## Details

`min.weights`: the weighted likelihood equation could have more than one solution. These roots appear for particular situation depending on contamination level and type. We introduce the `min.weight` parameter in order to choose only between roots that do not down weight everything. This is not still the optimal solution, and perhaps, in the new release, this part will be change.

`min.weights.aic` is used as `min.weights` but in the full model. The algorithm used to classify the observations as additive outliers is made by a genetic algorithm. The `population.size`, `population.choose` and `elements.random` are parameters related to this algorithm.

The function `wle.ar.wls` is used to estimate the parameter of an autoregressive model by weighted least squares where the weights are those from the weighted likelihood estimating equation of the full model (the model with the highest order).

## Value

A list of class `wle.aic.ar` with the following components:

<code>full.model</code>	the results for the full model, that is an object of class <code>wle.arima</code> see <a href="#">wle.ar</a> help for further details.
<code>waic</code>	Weighted Akaike Information Criterion for each submodels.
<code>call</code>	<code>match.call</code> result.

## Author(s)

Claudio Agostinelli

## References

Agostinelli C, (2004) Robust Akaike Information Criterion for ARMA models, *Rendiconti per gli Studi Economici Quantitativi*, 1-14, isbn: 88-88037-10-1.

Agostinelli C., (2003) Robust time series estimation via weighted likelihood, in: *Development in Robust Statistics. International Conference on Robust Statistics 2001*, Eds. Dutter, R. and Filzmoser, P. and Rousseeuw, P. and Gather, U., Physica Verlag.

## See Also

[wle.ar](#)

## Examples

```
data(rocky)

res <- wle.aic.ar(x=rocky, order=c(6,0), group=50, group.start=30, method="WLS")
res
plot(res$full.model$weights)
```



---

wle.aic.ar.summaries *Summaries and methods for wle.aic.ar*


---

## Description

All these functions are [methods](#) for class `wle.aic.ar` or `summary.wle.aic.ar`.

## Usage

```
## S3 method for class 'wle.aic.ar'
summary(object, num.max=20, verbose=FALSE, ...)

## S3 method for class 'wle.aic.ar'
print(x, digits = max(3, getOption("digits") - 3),
      num.max=max(1, nrow(x$waic)), ...)

## S3 method for class 'summary.wle.aic.ar'
print(x, digits = max(3, getOption("digits") - 3), ...)
```

## Arguments

<code>object</code>	an object of class <code>wle.aic.ar</code> .
<code>x</code>	an object of class <code>wle.aic</code> or <code>summary.wle.aic.ar</code> .
<code>digits</code>	number of digits to be used for most numbers.
<code>num.max</code>	the max number of models should be reported.
<code>verbose</code>	if TRUE warnings are printed.
<code>...</code>	additional arguments affecting the summary produced (in <code>summary.wle.aic.ar</code> ) or further arguments passed to or from other methods (in <code>print.wle.aic.ar</code> and <code>print.summary.wle.aic.ar</code> ).

## Value

`summary.wle.aic.ar` returns a list:

<code>waic</code>	the first <code>num.max</code> best models with their Weighted Akaike Information Criterion.
<code>num.max</code>	the number of models reported.
<code>call</code>	

## Author(s)

Claudio Agostinelli

## See Also

[wle.aic.ar](#) a function for evaluate the Weighted Akaike Information Criterion for autoregressive models.

---

wle.aic.summaries      *Summaries and methods for wle.aic*

---

### Description

All these functions are [methods](#) for class `wle.aic` or `summary.wle.aic`.

### Usage

```
## S3 method for class 'wle.aic'
summary(object, num.max=20, verbose=FALSE, ...)

## S3 method for class 'wle.aic'
print(x, digits = max(3, getOption("digits") - 3),
      num.max=max(1, nrow(x$waic)), ...)

## S3 method for class 'summary.wle.aic'
print(x, digits = max(3, getOption("digits") - 3), ...)
```

### Arguments

<code>object</code>	an object of class <code>wle.aic</code> .
<code>x</code>	an object of class <code>wle.aic</code> or <code>summary.wle.aic</code> .
<code>digits</code>	number of digits to be used for most numbers.
<code>num.max</code>	the max number of models should be reported.
<code>verbose</code>	if TRUE warnings are printed.
<code>...</code>	additional arguments affecting the summary produced (in <code>summary.wle.aic</code> ) or further arguments passed to or from other methods (in <code>print.wle.aic</code> and <code>print.summary.wle.aic</code> ).

### Value

`summary.wle.aic` returns a list:

<code>waic</code>	the first <code>num.max</code> best models with their Weighted Akaike Information Criterion.
<code>num.max</code>	the number of models reported.
<code>call</code>	

### Author(s)

Claudio Agostinelli

### See Also

[wle.aic](#) a function for evaluate the Weighted Akaike Information Criterion in the linear models.

**Description**

This is a preliminary version of functions for the estimation of the autoregressive parameters via Weighted Likelihood Estimating Equations and a classification algorithm. The main function is `wle.ar`, the remain functions are for internal use and they should not call by the users. They are not documented here.

**Usage**

```
wle.ar(x, order=c(1, 0), seasonal=list(order = c(0, 0),
  period = NA), group, group.start, group.step=group.start,
  xreg=NULL, include.mean=TRUE, na.action=na.fail,
  tol=10-6, tol.step=tol, equal=10-3, equal.step=equal,
  raf="HD", smooth=0.0031, smooth.ao=smooth, boot=10,
  boot.start=10, boot.step=boot.start, num.sol=1, x.init=0,
  x.seasonal.init=0, max.iter.out=20, max.iter.in=50,
  max.iter.start=200, max.iter.step=500, verbose=FALSE,
  w.level=0.4, min.weights=0.5, population.size=10,
  population.choose=5, elements.random=2, wle.start=FALSE,
  init.values=NULL, num.max=NULL, num.sol.step=2, approx.w=TRUE)
```

**Arguments**

<code>x</code>	a univariate time series.
<code>order</code>	a specification of the non-seasonal part of the ARI model: the two components ( $p, d$ ) are the AR order and the degree of differencing.
<code>seasonal</code>	a specification of the seasonal part of the ARI model, plus the period (which defaults to frequency( $x$ )).
<code>group</code>	the dimension of the bootstrap subsamples.
<code>group.start</code>	the dimension of the bootstrap subsamples used in the starting process if <code>wle.init=TRUE</code> .
<code>group.step</code>	the dimension of the bootstrap subsamples used in a step, it must be less than <code>group</code> .
<code>xreg</code>	optionally, a vector or matrix of external regressors, which must have the same number of rows as <code>x</code> .
<code>include.mean</code>	Should the ARI model include a mean term? The default is <code>TRUE</code> for undifferenced series, <code>FALSE</code> for differenced ones (where a mean would not affect the fit nor predictions).
<code>na.action</code>	function to be applied to remove missing values.
<code>tol</code>	the absolute accuracy to be used to achieve convergence of the algorithm.
<code>tol.step</code>	the absolute accuracy to be used to achieve convergence in a step.

equal	the absolute value for which two roots are considered the same. (This parameter must be greater than tol).
equal.step	the absolute value for which two roots are considered the same in a step. (This parameter must be greater than tol.step).
raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
smooth	the value of the smoothing parameter.
smooth.ao	the value of the smoothing parameter used in the outliers classificaton, default equal to smooth.
boot	the number of starting points based on bootstrap subsamples to use in the search of the roots.
boot.start	the number of starting points based on bootstrap subsamples to use in the search of the roots in the starting process.
boot.step	the number of starting points based on bootstrap subsamples to use in the search of the roots in a step.
num.sol	maximum number of roots to be searched.
num.sol.step	maximum number of roots to be searched in a step.
x.init	initial values, a vector with the same length of the AR order, or a number, default is 0.
x.seasonal.init	initial values, a vector with the same length of the SAR order, or a number, default is 0.
max.iter.out	maximum number of iterations in the outer loop.
max.iter.in	maximum number of iterations in the inner loop.
max.iter.start	maximum number of iterations in the starting process.
max.iter.step	maximum number of iterations in a step.
w.level	the threshold used to decide if an observation could be an additive outlier.
population.size	see details.
population.choose	see details.
elements.random	see details.
num.max	maximum number of observations can be considered as possible additive outliers.
wle.start	if TRUE a weighted likelihood estimation is used to have a starting value.
init.values	a vector with initial values for the AR and seasonal AR coefficients and the innovations variance.
verbose	if TRUE warnings are printed.
min.weights	see details.
approx.w	logical: if TRUE an approximation is used to evaluate the weights in the outlier identification procedure.

## Details

`min.weight`: the weighted likelihood equation could have more than one solution. These roots appear for particular situation depending on contamination level and type. We introduce the `min.weight` parameter in order to choose only between roots that do not down weight everything. This is not still the optimal solution, and perhaps, in the new release, this part will be change.

The algorithm used to classify the observations as additive outliers is made by a genetic algorithm. The `population.size`, `population.choose` and `elements.random` are parameters related to this algorithm.

## Value

<code>coef</code>	a vector of AR and regression coefficients.
<code>sigma2.coef</code>	the estimated variance matrix of the coefficients <code>coef</code> .
<code>sigma2</code>	the WLE of the innovations variance.
<code>arma</code>	a compact form of the specification, as a vector giving the number of AR, MA=0, seasonal AR and seasonal MA=0 coefficients, plus the period and the number of non-seasonal and seasonal differences.
<code>resid</code>	the residuals.
<code>resid.with.ao</code>	the residuals with the additive outliers effects.
<code>resid.without.ao</code>	the residuals without the additive outliers effects.
<code>x.ao</code>	the time series without the additive outliers effects.
<code>call</code>	the matched call.
<code>series</code>	the name of the series <code>x</code> .
<code>weights</code>	the weights.
<code>weights.with.ao</code>	the weights with the additive outliers effects.
<code>weights.without.ao</code>	the weights without the additive outliers effects
<code>tot.sol</code>	the number of solutions found.
<code>not.conv</code>	the number of starting points that does not converge after the <code>max.iter.out</code> iteration are reached.
<code>ao.position</code>	the position of the additive outliers.

## Author(s)

Claudio Agostinelli

## References

- Agostinelli C., (2001) Robust time series estimation via weighted likelihood: some preliminary results, *Working Paper n. 2001.3* Department of Statistics, University of Padova.
- Agostinelli C., (2003) Robust time series estimation via weighted likelihood, in: *Development in Robust Statistics. International Conference on Robust Statistics 2001*, Eds. Dutter, R. and Filzmoser, P. and Rousseeuw, P. and Gather, U., Physica Verlag.

**Examples**

```
data(lh)
wle.ar(x=lh, order=c(3,0), group=30)
```

---

wle.binomial

*Robust Estimation in the Binomial Model*


---

**Description**

wle.binomial is used to robust estimate the proportion parameters via Weighted Likelihood.

**Usage**

```
wle.binomial(x, size, boot=30, group, num.sol=1, raf="HD",
             tol=10-6, equal=10-3, max.iter=500,
             verbose=FALSE)
```

**Arguments**

x	a vector contain the number of success in each size trials.
size	number of trials.
boot	the number of starting points based on bootstrap subsamples to use in the search of the roots.
group	the dimension of the bootstrap subsamples. The default value is $\max(\text{round}(\text{length}(x)/4), 2)$ .
num.sol	maximum number of roots to be searched.
raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.
equal	the absolute value for which two roots are considered the same. (This parameter must be greater than tol).
max.iter	maximum number of iterations.
verbose	if TRUE warnings are printed.

**Value**

wle.binomial returns an object of class "wle.binomial".

Only print method is implemented for this class.

The object returned by wle.binomial are:

p	the estimator of the proportion parameter, one value for each root found.
---	---

tot.weights	the sum of the weights divide by the number of observations, one value for each root found.
weights	the weights associated to each observation, one column vector for each root found.
f.density	the non-parametric density estimation.
m.density	the smoothed model.
delta	the Pearson residuals.
call	the match.call().
tot.sol	the number of solutions found.
not.conv	the number of starting points that does not converge after the max.iter iteration are reached.

**Author(s)**

Claudio Agostinelli

**References**

Markatou, M., Basu, A., and Lindsay, B.G., (1997) Weighted likelihood estimating equations: The discrete case with applications to logistic regression, *Journal of Statistical Planning and Inference*, 57, 215-232.

Agostinelli, C., (1998) Inferenza statistica robusta basata sulla funzione di verosimiglianza pesata: alcuni sviluppi, *Ph.D Thesis*, Department of Statistics, University of Padova.

**Examples**

```
library(wle)

set.seed(1234)

x <- rbinom(20,p=0.2,size=10)
wle.binomial(x,size=10)

x <- c(rbinom(20,p=0.2,size=10),rbinom(10,p=0.9,size=10))
wle.binomial(x,size=10)
```

---

wle.cp

*Weighted Mallows Cp*


---

**Description**

The Weighted Mallows Cp is evaluated for each submodel.

**Usage**

```
wle.cp(formula, data=list(), model=TRUE, x=FALSE,
       y=FALSE, boot=30, group, var.full=0, num.sol=1,
       raf="HD", smooth=0.031, tol=10-6,
       equal=10-3, max.iter=500, min.weight=0.5,
       method="full", alpha=2, contrasts=NULL, verbose=FALSE)
```

**Arguments**

formula	a symbolic description of the model to be fit. The details of model specification are given below.
data	an optional data frame containing the variables in the model. By default the variables are taken from the environment which <code>wle.cp</code> is called from.
model, x, y	logicals. If TRUE the corresponding components of the fit (the model frame, the model matrix, the response.)
boot	the number of starting points based on bootstrap subsamples to use in the search of the roots.
group	the dimension of the bootstrap subsamples. The default value is $\max(\text{round}(\text{size}/4), \text{var})$ where <i>size</i> is the number of observations and <i>var</i> is the number of variables.
var.full	the value of variance to be used in the denominator of the WCP, if 0 the variance estimated from the full model is used.
num.sol	maximum number of roots to be searched.
raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
smooth	the value of the smoothing parameter.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.c
equal	the absolute value for which two roots are considered the same. (This parameter must be greater than <code>tol</code> ).
max.iter	maximum number of iterations.
min.weight	see details.
method	see details.
alpha	penalty value.
contrasts	an optional list. See the <code>contrasts.arg</code> of <code>model.matrix.default</code> .
verbose	if TRUE warnings are printed.

**Details**

Models for `wle.cp` are specified symbolically. A typical model has the form `response ~ terms` where `response` is the (numeric) response vector and `terms` is a series of terms which specifies a linear predictor for response. A terms specification of the form `first+second` indicates all the terms in `first` together with all the terms in `second` with duplicates removed. A specification of



the form `first:second` indicates the the set of terms obtained by taking the interactions of all terms in `first` with all terms in `second`. The specification `first*second` indicates the *cross* of `first` and `second`. This is the same as `first+second+first:second`.

`min.weight`: the weighted likelihood equation could have more than one solution. These roots appear for particular situation depending on contamination level and type. The presence of multiple roots in the full model can create some problem in the set of weights we should use. Actually, the selection of the root is done by the minimum scale error provided. Since this choice is not always the one would choose, we introduce the `min.weight` parameter in order to choose only between roots that do not down weight everything. This is not still the optimal solution, and perhaps, in the new release, this part will be change.

`method`: this parameter, when set to "reduced", allows to use weights based on the reduced model. This is strongly discourage since the robust and asymptotic property of this kind of weighted Cp are not as good as the one based on `method="full"`.

## Value

`wle.cp` returns an object of `class "wle.cp"`.

The function `summary` is used to obtain and print a summary of the results. The generic accessor functions `coefficients` and `residuals` extract coefficients and residuals returned by `wle.cp`. The object returned by `wle.cp` are:

<code>wcp</code>	Weighted Mallows Cp for each submodels
<code>coefficients</code>	the parameters estimator, one row vector for each root found and each submodel.
<code>scale</code>	an estimation of the error scale, one value for each root found and each submodel.
<code>residuals</code>	the unweighted residuals from the estimated model, one column vector for each root found and each submodel.
<code>tot.weights</code>	the sum of the weights divide by the number of observations, one value for each root found and each submodel.
<code>weights</code>	the weights associated to each observation, one column vector for each root found and each submodel.
<code>freq</code>	the number of starting points converging to the roots.
<code>call</code>	the <code>match.call()</code> .
<code>contrasts</code>	
<code>xlevels</code>	
<code>terms</code>	the model frame.
<code>model</code>	if <code>model=TRUE</code> a matrix with first column the dependent variable and the remain column the explanatory variables for the full model.
<code>x</code>	if <code>x=TRUE</code> a matrix with the explanatory variables for the full model.
<code>y</code>	if <code>y=TRUE</code> a vector with the dependent variable.
<code>info</code>	not well working yet, if 0 no error occurred.

## Author(s)

Claudio Agostinelli

## References

Agostinelli, C., (1999). Robust model selection in regression via weighted likelihood methodology, *Working Paper n. 1999.4*, Department of Statistics, University of Padova.

Agostinelli, C., (2002). Robust model selection in regression via weighted likelihood methodology, *Statistics & Probability Letters*, 56, 289-300.

Agostinelli, C., (1998). Inferenza statistica robusta basata sulla funzione di verosimiglianza pesata: alcuni sviluppi, *Ph.D Thesis*, Department of Statistics, University of Padova.

Agostinelli, C., Markatou, M., (1998). A one-step robust estimator for regression based on the weighted likelihood reweighting scheme, *Statistics & Probability Letters*, Vol. 37, n. 4, 341-350.

Agostinelli, C., (1998). Verosimiglianza pesata nel modello di regressione lineare, *XXXIX Riunione scientifica della Societ a Italiana di Statistica*, Sorrento 1998.

## See Also

[wle.smooth](#) an algorithm to choose the smoothing parameter for normal distribution and normal kernel, [wle.lm](#) a function for estimating linear models with normal distribution error and normal kernel.

## Examples

```
library(wle)

x.data <- c(runif(60,20,80),runif(5,73,78))
e.data <- rnorm(65,0,0.6)
y.data <- 8*log(x.data+1)+e.data
y.data[61:65] <- y.data[61:65]-4
z.data <- c(rep(0,60),rep(1,5))

plot(x.data,y.data,xlab="X",ylab="Y")

xx.data <- cbind(x.data,x.data^2,x.data^3,log(x.data+1))
colnames(xx.data) <- c("X","X^2","X^3","log(X+1)")

result <- wle.cp(y.data~xx.data,boot=10,group=10,num.sol=2)

summary(result)

plot(result,num.max=15)

result <- wle.cp(y.data~xx.data+z.data,boot=10,group=10,num.sol=2)

summary(result)

plot(result,num.max=15)
```

---

wle.cp.summaries      *Summaries and methods for wle.cp*

---

## Description

All these functions are [methods](#) for class `wle.cp` or `summary.wle.cp`.

## Usage

```
## S3 method for class 'wle.cp'
summary(object, num.max=20, verbose=FALSE, ...)

## S3 method for class 'wle.cp'
print(x, digits = max(3, getOption("digits") - 3),
      num.max=max(1, nrow(x$wcp)), ...)

## S3 method for class 'summary.wle.cp'
print(x, digits = max(3, getOption("digits") - 3), ...)
```

## Arguments

<code>object</code>	an object of class <code>wle.cp</code> .
<code>x</code>	an object of class <code>wle.cp</code> or <code>summary.wle.cp</code> .
<code>digits</code>	number of digits to be used for most numbers.
<code>num.max</code>	the max number of models should be reported.
<code>verbose</code>	if TRUE warnings are printed.
<code>...</code>	additional arguments affecting the summary produced (in <code>summary.wle.cp</code> ) or further arguments passed to or from other methods (in <code>print.wle.cp</code> and <code>print.summary.wle.cp</code> ).

## Value

`summary.wle.cp` returns a list:

<code>wcp</code>	the first <code>num.max</code> best models with their Weighted Mallows Cp.
<code>num.max</code>	the number of models reported.
<code>call</code>	

## Author(s)

Claudio Agostinelli

## See Also

[wle.cp](#) a function for evaluate the Weighted Mallows Cp in the linear models.

**Description**

The Weighted Cross-Validation methods is used to choose the best linear model.

**Usage**

```
wle.cv(formula, data=list(), model=TRUE, x=FALSE,
       y=FALSE, monte.carlo=500, split, boot=30,
       group, num.sol=1, raf="HD", smooth=0.031,
       tol=10-6, equal=10-3, max.iter=500,
       min.weight=0.5, contrasts=NULL,
       type=c('fast', 'slow'), verbose=FALSE)
```

**Arguments**

formula	a symbolic description of the model to be fit. The details of model specification are given below.
data	an optional data frame containing the variables in the model. By default the variables are taken from the environment which <code>wle.cv</code> is called from.
model, x, y	logicals. If TRUE the corresponding components of the fit (the model frame, the model matrix, the response.)
monte.carlo	the number of Monte Carlo replication we use to estimate the average prediction error.
split	the size of the construction sample. When the suggested value is outside the possible range, the split size is let equal to $\max(\text{round}(\text{size}^{3/4}), \text{var} + 2)$ where <i>size</i> is the number of observations and <i>var</i> is the number of variables.
boot	the number of starting points based on bootstrap subsamples to use in the search of the roots.
group	the dimension of the bootstrap subsamples. The default value is $\max(\text{round}(\text{size}/4), \text{var})$ where <i>size</i> is the number of observations and <i>var</i> is the number of variables.
num.sol	maximum number of roots to be searched.
raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
smooth	the value of the smoothing parameter.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.
equal	the absolute value for which two roots are considered the same. (This parameter must be greater than tol).
max.iter	maximum number of iterations.

min.weight	see details.
contrasts	an optional list. See the contrasts.arg of model.matrix.default.
type	when 'fast' a weighted least squares is used to evaluate the parameters in the submodels, while if 'slow' a weighted likelihood is used.
verbose	if TRUE warnings are printed.

## Details

Models for `wle.cv` are specified symbolically. A typical model has the form `response ~ terms` where `response` is the (numeric) response vector and `terms` is a series of terms which specifies a linear predictor for `response`. A terms specification of the form `first+second` indicates all the terms in `first` together with all the terms in `second` with duplicates removed. A specification of the form `first:second` indicates the the set of terms obtained by taking the interactions of all terms in `first` with all terms in `second`. The specification `first*second` indicates the *cross* of `first` and `second`. This is the same as `first+second+first:second`.

`min.weight`: the weighted likelihood equation could have more than one solution. These roots appear for particular situation depending on contamination level and type. The presence of multiple roots in the full model can create some problem in the set of weights we should use. Actually, the selection of the root is done by the minimum scale error provided. Since this choice is not always the one would choose, we introduce the `min.weight` parameter in order to choose only between roots that do not down weight everything. This is not still the optimal solution, and perhaps, in the new release, this part will be change.

## Value

`wle.cv` returns an object of class `"wle.cv"`.

The function `summary` is used to obtain and print a summary of the results. The generic accessor functions `coefficients` and `residuals` extract coefficients and residuals returned by `wle.cv`. The object returned by `wle.cv` are:

<code>wcv</code>	Weighted Cross-Validation for each submodels
<code>coefficients</code>	the parameters estimator, one row vector for each root found in the full model.
<code>scale</code>	an estimation of the error scale, one value for each root found in the full model.
<code>residuals</code>	the unweighted residuals from the estimated model, one column vector for each root found in the full model.
<code>tot.weights</code>	the sum of the weights divide by the number of observations, one value for each root found in the full model.
<code>weights</code>	the weights associated to each observation, one column vector for each root found in the full model.
<code>freq</code>	the number of starting points converging to the roots.
<code>index</code>	position of the root used for the weights.
<code>call</code>	the <code>match.call()</code> .
<code>contrasts</code>	
<code>xlevels</code>	

terms	the model frame.
model	if model=TRUE a matrix with first column the dependent variable and the remain column the explanatory variables for the full model.
x	if x=TRUE a matrix with the explanatory variables for the full model.
y	if y=TRUE a vector with the dependent variable.
info	not well working yet, if 0 no error occurred.

### Author(s)

Claudio Agostinelli

### References

Agostinelli, C., (1999). Robust model selection by Cross-Validation via weighted likelihood methodology, *Working Paper n. 1999.37*, Department of Statistics, University of Padova.

Agostinelli, C., (1998). Inferenza statistica robusta basata sulla funzione di verosimiglianza pesata: alcuni sviluppi, *Ph.D Thesis*, Department of Statistics, University of Padova.

Agostinelli, C., Markatou, M., (1998). A one-step robust estimator for regression based on the weighted likelihood reweighting scheme, *Statistics & Probability Letters*, Vol. 37, n. 4, 341-350.

Agostinelli, C., (1998). Verosimiglianza pesata nel modello di regressione lineare, *XXXIX Riunione scientifica della Societ a Italiana di Statistica*, Sorrento 1998.

### See Also

[wle.smooth](#) an algorithm to choose the smoothing parameter for normal distribution and normal kernel, [wle.lm](#) a function for estimating linear models with normal distribution error and normal kernel.

### Examples

```
library(wle)

set.seed(1234)

x.data <- c(runif(60,20,80),runif(5,73,78))
e.data <- rnorm(65,0,0.6)
y.data <- 8*log(x.data+1)+e.data
y.data[61:65] <- y.data[61:65]-4
z.data <- c(rep(0,60),rep(1,5))

plot(x.data,y.data,xlab="X",ylab="Y")

xx.data <- cbind(x.data,x.data^2,x.data^3,log(x.data+1))
colnames(xx.data) <- c("X","X^2","X^3","log(X+1)")

result <- wle.cv(y.data~xx.data,boot=20,num.sol=2)

summary(result)
```

```

result <- wle.cv(y.data~xx.data+z.data,boot=20,num.sol=2,
               monte.carlo=1000,split=50)

summary(result)

```

---

wle.cv.summaries      *Summaries and methods for wle.cv*

---

## Description

All these functions are [methods](#) for class `wle.cv` or `summary.wle.cv`.

## Usage

```

## S3 method for class 'wle.cv'
summary(object, num.max=20, verbose=FALSE, ...)

## S3 method for class 'wle.cv'
print(x, digits = max(3, getOption("digits") - 3),
      num.max=max(1, nrow(x$wcv)), ...)

## S3 method for class 'summary.wle.cv'
print(x, digits = max(3, getOption("digits") - 3), ...)

```

## Arguments

<code>object</code>	an object of class <code>wle.cv</code> .
<code>x</code>	an object of class <code>wle.cv</code> or <code>summary.wle.cv</code> .
<code>digits</code>	number of digits to be used for most numbers.
<code>num.max</code>	the max number of models should be reported.
<code>verbose</code>	if TRUE warnings are printed.
<code>...</code>	additional arguments affecting the summary produced (in <code>summary.wle.cv</code> ) or further arguments passed to or from other methods (in <code>print.wle.cv</code> and <code>print.summary.wle.cv</code> ).

## Value

`summary.wle.cv` returns a list:

<code>wcv</code>	the first <code>num.max</code> best models with their estimated prediction error using WCV.
<code>num.max</code>	the number of models reported.
<code>call</code>	

## Author(s)

Claudio Agostinelli

**See Also**

[wle.cv](#) a function for evaluate the Weighted Cross Validation criterion in the linear models.

---

wle.fracdiff

*Fit Fractional Models to Time Series - Preliminary Version*


---

**Description**

This is a preliminary version of functions for the estimation of the fractional parameter via Weighted Likelihood Estimating Equations and a classification algorithm. The main function is `wle.fracdiff`, the remain functions are for internal use and they should not call by the users. They are not documented here.

**Usage**

```
wle.fracdiff(x, lower, upper, M, group, na.action=na.fail,
  tol=10-6, equal=10-3, raf="HD", smooth=0.0031, smooth.ao=smooth,
  boot=10, num.sol=1, x.init=rep(0,M), use.uniroot=FALSE, max.iter.out=20,
  max.iter.in=100, max.iter.step=5000, max.iter.start=max.iter.step,
  verbose=FALSE, w.level=0.4, min.weights=0.5, init.values=NULL,
  num.max=length(x), include.mean=FALSE, ao.list=NULL, elitist=5,
  size.generation=5, size.population=10, type.selection="roulette",
  prob.crossover=0.8, prob.mutation=0.02, type.scale="none", scale.c=2)
```

**Arguments**

<code>x</code>	a univariate time series.
<code>lower</code>	the lower end point of the interval to be searched.
<code>upper</code>	the upper end point of the interval to be searched.
<code>M</code>	the order of the finite memory process used to estimate the $d$ parameter.
<code>group</code>	the dimension of the bootstrap subsamples.
<code>na.action</code>	function to be applied to remove missing values.
<code>tol</code>	the absolute accuracy to be used to achieve convergence of the algorithm.
<code>equal</code>	the absolute value for which two roots are considered the same. (This parameter must be greater than <code>tol</code> ).
<code>raf</code>	type of Residual adjustment function to be use: <code>raf="HD"</code> : Hellinger Distance RAF, <code>raf="NED"</code> : Negative Exponential Disparity RAF, <code>raf="SCHI2"</code> : Symmetric Chi-Squared Disparity RAF.
<code>smooth</code>	the value of the smoothing parameter.
<code>smooth.ao</code>	the value of the smoothing parameter used in the outliers classificaton, default equal to <code>smooth</code> .
<code>boot</code>	the number of starting points based on bootstrap subsamples to use in the search of the roots.



num.sol	maximum number of roots to be searched.
x.init	initial values, a vector with the same length of the M parameter, or a number, default is 0.
use.uniroot	default: FALSE, if TRUE in each step the weighted likelihood estimating equations is solved, otherwise, a maximization is performed on a weighted log-likelihood function with fixed weights. The estimators obtain with the two methods is the same.
max.iter.out	maximum number of iterations in the outer loop.
max.iter.in	maximum number of iterations in the inner loop.
max.iter.step	maximum number of iterations in a step.
max.iter.start	maximum number of iterations in the starting process.
verbose	if TRUE warnings are printed.
w.level	the threshold used to decide if an observation could be an additive outlier.
init.values	a vector with initial values for the d and the innovations variance.
num.max	maximum number of observations can be considered as possible additive outliers.
include.mean	Should the model include a mean term? The default is TRUE.
ao.list	possible list of pattern of additive outliers.
min.weights	see details.
size.population	see details.
size.generation	see details.
prob.crossover	see details.
prob.mutation	see details.
type.scale	see details.
type.selection	see details.
elitist	see details.
scale.c	see details.

## Details

min.weight: the weighted likelihood equation could have more than one solution. These roots appear for particular situation depending on contamination level and type. We introduce the min.weight parameter in order to choose only between roots that do not down weight everything. This is not still the optimal solution, and perhaps, in the new release, this part will be change.

The algorithm used to classify the observations as additive outliers is a simple genetic algorithm as described in Goldberg (1989). The size.population, size.generation, type.selection, prob.crossover, prob.mutation, type.scale, type.selection, elitist and scale.c are parameters related to this algorithm.

**Value**

<code>d</code>	the WLE of the fractional parameter.
<code>sigma2</code>	the WLE of the innovations variance.
<code>x.mean</code>	the WLE of the mean.
<code>resid</code>	the residuals.
<code>resid.without.ao</code>	the residuals with the additive outliers effects.
<code>resid.with.ao</code>	the residuals without the additive outliers effects.
<code>x.ao</code>	the time series without the additive outliers effects.
<code>call</code>	the matched call.
<code>weights</code>	the weights.
<code>weights.with.ao</code>	the weights with the additive outliers effects.
<code>weights.without.ao</code>	the weights without the additive outliers effects
<code>tot.sol</code>	the number of solutions found.
<code>not.conv</code>	the number of starting points that does not converge after the <code>max.iter.out</code> iteration are reached.
<code>ao.position</code>	the position of the additive outliers.

**Author(s)**

Claudio Agostinelli

**References**

Agostinelli C., Bisaglia L., (2002) Robust estimation of ARFIMA processes, manuscript.  
 Goldberg, David E., (1989) Genetic Algorithms in Search, Optimization and Machine Learning.  
 Addison-Wesley Pub. Co. ISBN: 0201157675

**Examples**

```
## Not run:
set.seed(1234)
resw <- wle.fracdiff(Nile, M=100, include.mean=TRUE, lower=0.01, upper=0.96, group=20)
resw$d
resw$sigma2
resw$x.mean

x <- Nile
x[50] <- x[50]+4*sd(x)

set.seed(1234)
resw <- wle.fracdiff(x, M=100, include.mean=TRUE, lower=0.01, upper=0.96, group=40)
resw$d
resw$sigma2
```

```

    resw$x.mean
    resw$a0.position

## End(Not run)

```

---

wle.gamma

*Robust Estimation in the Gamma model*


---

## Description

wle.gamma is used to robust estimate the shape and the scale parameters via Weighted Likelihood, when the majority of the data are from a gamma distribution.

## Usage

```

wle.gamma(x, boot=30, group, num.sol=1, raf="HD", smooth=0.008,
          tol=10^(-6), equal=10^(-3), max.iter=500,
          shape.int=c(0.01, 100), use.smooth=TRUE, tol.int,
          verbose=FALSE, maxiter=1000)

```

## Arguments

x	a vector contain the observations.
boot	the number of starting points based on bootstrap subsamples to use in the search of the roots.
group	the dimension of the bootstrap subsamples. The default value is $\max(\text{round}(\text{size}/4), \text{var})$ where <i>size</i> is the number of observations and <i>var</i> is the number of variables.
num.sol	maximum number of roots to be searched.
raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
smooth	the value of the smoothing parameter.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.
equal	the absolute value for which two roots are considered the same. (This parameter must be greater than tol).
max.iter	maximum number of iterations for the main function.
shape.int	a 2 dimension vector for the interval search of the shape parameter.
use.smooth	if FALSE the unsmoothed model is used. This is usefull when the integration routine does not work well.
tol.int	the absolute accuracy to be used in the integration routine. The default value is $\text{tol} * 10^{-4}$ .
verbose	if TRUE warnings are printed.
maxiter	maximum number of iterations. This value is passed to <code>uniroot</code> function.

## Details

The gamma is parametrized as follows ( $\alpha = scale, \omega = shape$ ):

$$f(x) = 1/(\alpha^\omega \Gamma(\omega)) x^{\omega-1} e^{-x/\alpha}$$

for  $x > 0, \alpha > 0$  and  $\omega > 0$ .

The function use `uniroot` to solve the estimating equation for *shape*, errors from `uniroot` are handled by `try`. If errors occurs then the function returns NA.

You can use `shape.int` to avoid them. It also use a fortran routine (`dqagp`) to calculate the smoothed model, i.e., evaluate the integral. Sometime the accuracy is not satisfactory, you can use `use.smooth=FALSE` to have an approximate estimation using the model instead of the smoothed model.

The Folded Normal distribution is use as kernel. The bandwidth is  $smooth * shape / scale^2$ .

## Value

`wle.gamma` returns an object of class "wle.gamma".

Only print method is implemented for this class.

The object returned by `wle.gamma` are:

<code>shape</code>	the estimator of the shape parameter, one value for each root found.
<code>scale</code>	the estimator of the scale parameter, one value for each root found.
<code>rate</code>	the estimator of the rate parameter ( $1/scale$ ), one value for each root found.
<code>tot.weights</code>	the sum of the weights divide by the number of observations, one value for each root found.
<code>weights</code>	the weights associated to each observation, one column vector for each root found.
<code>f.density</code>	the non-parametric density estimation.
<code>m.density</code>	the smoothed model.
<code>delta</code>	the Pearson residuals.
<code>call</code>	the <code>match.call()</code> .
<code>tot.sol</code>	the number of solutions found.
<code>not.conv</code>	the number of starting points that does not converge after the <code>max.iter</code> iteration are reached.

## Author(s)

Claudio Agostinelli

## References

Markatou, M., Basu, A. and Lindsay, B.G., (1998). Weighted likelihood estimating equations with a bootstrap root search, *Journal of the American Statistical Association*, 93, 740-750.

Agostinelli, C., (1998). Inferenza statistica robusta basata sulla funzione di verosimiglianza pesata: alcuni sviluppi, *Ph.D Thesis*, Department of Statistics, University of Padova.

**Examples**

```
library(wle)

x <- rgamma(n=100, shape=2, scale=2)

wle.gamma(x)

x <- c(rgamma(n=30, shape=2, scale=2), rgamma(n=100, shape=20, scale=20))

wle.gamma(x, boot=10, group=10, num.sol=2) # depending on the sample, one or two roots.
```

---

wle.glm

*Robust Fitting Generalized Linear Models using Weighted Likelihood*


---

**Description**

wle.glm is used to robustly fit generalized linear models, specified by giving a symbolic description of the linear predictor and a description of the error distribution.

**Usage**

```
wle.glm(formula, family = binomial, data, weights, subset,
        na.action, start = NULL, etastart, mustart, offset,
        control = list(glm = glm.control(...), wle = wle.glm.control()),
        model = TRUE, method = "wle.glm.fit", x = FALSE, y = TRUE,
        contrasts = NULL, dist.method = "euclidean", ...)

wle.glm.fit(x, y, weights = NULL, wle.weights = rep(1, NROW(y)),
           start = NULL, etastart = NULL, mustart = NULL, offset = rep(0, NROW(y)),
           family = gaussian(), control = list(glm=glm.control(),
           wle=wle.glm.control()), dist.method='euclidean',
           intercept = TRUE, dispersion=NULL)

## S3 method for class 'wle.glm'
weights(object, type = c("prior", "working", "wle"), root="all", ...)
```

**Arguments**

formula	an object of class " <a href="#">formula</a> " (or one that can be coerced to that class): a symbolic description of the model to be fitted. The details of model specification are given under 'Details'.
family	a description of the error distribution and link function to be used in the model. This can be a character string naming a family function, a family function or the result of a call to a family function. (See <a href="#">family</a> for details of family functions.)

data	an optional data frame, list or environment (or object coercible by <code>as.data.frame</code> to a 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>wle.glm</code> is called.
weights	an optional vector of ‘prior weights’ to be used in the fitting process. Should be NULL or a numeric vector.
subset	an optional vector specifying a subset of observations to be used in the fitting process.
na.action	a function which indicates what should happen when the data contain NAs. The default is set by the <code>na.action</code> setting of <code>options</code> , and is <code>na.fail</code> if that is unset. The ‘factory-fresh’ default is <code>na.omit</code> . Another possible value is NULL, no action. Value <code>na.exclude</code> can be useful.
start	starting values for the parameters in the linear predictor.
etastart	starting values for the linear predictor.
mustart	starting values for the vector of means.
offset	this can be used to specify an <i>a priori</i> known component to be included in the linear predictor during fitting. This should be NULL or a numeric vector of length equal to the number of cases. One or more <code>offset</code> terms can be included in the formula instead or as well, and if more than one is specified their sum is used. See <code>model.offset</code> .
control	a list with two components of parameters for controlling the fitting process. The first component ( <code>glm</code> ) is set using the function <code>glm.control</code> while the second component ( <code>wle</code> ) is set using the function <code>wle.glm.control</code> and it is used to set the parameters regarding the behaviour of the robust method. See the documentation of these functions for details.
model	a logical value indicating whether <i>model frame</i> should be included as a component of the returned value.
method	the method to be used in fitting the model. The default method <code>"wle.glm.fit"</code> uses iteratively reweighted least squares (IWLS). The only current alternative is <code>"model.frame"</code> which returns the model frame and does no fitting.
x, y	For <code>wle.glm</code> : logical values indicating whether the response vector and model matrix used in the fitting process should be returned as components of the returned value. For <code>wle.glm.fit</code> : <code>x</code> is a design matrix of dimension $n * p$ , and <code>y</code> is a vector of observations of length <code>n</code> .
contrasts	an optional list. See the <code>contrasts.arg</code> of <code>model.matrix.default</code> .
dist.method	distance method passed to <code>dist</code> to measure the distance between predictor rows.
intercept	logical. Should an intercept be included in the <i>null</i> model?
dispersion	numeric or NULL. If provided used as starting value.
object	an object inheriting from class <code>"wle.glm"</code> .
type	character, partial matching allowed. Type of weights to extract from the fitted model object.
root	character ( <code>"all"</code> ) or a number. For which solutions the weights are reported.

wle.weights	For wle.glm.fit these are weights used in the iterative algorithm evaluated at each step by wle.glm.weights.
...	For wle.glm: arguments to be passed by default to <code>glm.control</code> : see argument control. For weights: further arguments passed to or from other methods.

## Details

A typical predictor has the form `response ~ terms` where `response` is the (numeric) response vector and `terms` is a series of terms which specifies a linear predictor for response. For binomial and quasibinomial families the response can also be specified as a `factor` (when the first level denotes failure and all others success) or as a two-column matrix with the columns giving the numbers of successes and failures. A terms specification of the form `first + second` indicates all the terms in `first` together with all the terms in `second` with any duplicates removed.

A specification of the form `first:second` indicates the the set of terms obtained by taking the interactions of all terms in `first` with all terms in `second`. The specification `first*second` indicates the *cross* of `first` and `second`. This is the same as `first + second + first:second`.

The terms in the formula will be re-ordered so that main effects come first, followed by the interactions, all second-order, all third-order and so on: to avoid this pass a `terms` object as the formula.

Non-NULL weights can be used to indicate that different observations have different dispersions (with the values in `weights` being inversely proportional to the dispersions); or equivalently, when the elements of `weights` are positive integers  $w_i$ , that each response  $y_i$  is the mean of  $w_i$  unit-weight observations. In case of binomial GLM prior weights CAN NOT be used to give the number of trials when the response is the proportion of successes; in this situation please submit the response variable as two columns (first column successes, second column unsuccesses). They would rarely be used for a Poisson GLM.

`wle.glm.fit` is the workhorse function: it is not normally called directly but can be more efficient where the response vector and design matrix have already been calculated. However, this function needs starting values and does not look for possible multiple roots in the system of equations.

If more than one of `etastart`, `start` and `mustart` is specified, the first in the list will be used. It is often advisable to supply starting values for a `quasi` family, and also for families with unusual links such as `gaussian("log")`.

All of `weights`, `subset`, `offset`, `etastart` and `mustart` are evaluated in the same way as variables in `formula`, that is first in data and then in the environment of `formula`.

For the background to warning messages about ‘fitted probabilities numerically 0 or 1 occurred’ for binomial GLMs, see Venables & Ripley (2002, pp. 197–8).

Multiple roots may occur if the asymptotic weights are used or in the case of continuous models. The function implements the bootstrap root serach approach described in Markatou, Basu and Lindsay (1998) in order to find these roots.

## Value

`wle.glm` returns an object of class inheriting from `"wle.glm"`.

The function `summary` (i.e., `summary.wle.glm`) can be used to obtain or print a summary of the results and the function `anova` (i.e., `anova.wle.glm.root`) to produce an analysis of variance table.

The generic accessor functions `coefficients`, `effects`, `fitted.values` and `residuals` can be used to extract various useful features of the value returned by `wle.glm`.

`weights` extracts a vector of weights, one for each case/root in the fit (after subsetting and `na.action`).

An object of class "`wle.glm`" is a (variable length) list containing at least the following components:

`root1` which is a list with the following components:

<code>coefficients</code>	a named vector of coefficients
<code>residuals</code>	the <i>working</i> residuals, that is the residuals in the final iteration of the IWLS fit. Since cases with zero weights are omitted, their working residuals are NA.
<code>fitted.values</code>	the fitted mean values, obtained by transforming the linear predictors by the inverse of the link function.
<code>rank</code>	the numeric rank of the fitted linear model.
<code>family</code>	the <code>family</code> object used.
<code>linear.predictors</code>	the linear fit on link scale.
<code>deviance</code>	up to a constant, minus twice the maximized log-likelihood. Where sensible, the constant is chosen so that a saturated model has deviance zero.
<code>aic</code>	Akaike's <i>An Information Criterion</i> , minus twice the maximized log-likelihood plus twice the number of coefficients (so assuming that the dispersion is known).
<code>null.deviance</code>	The deviance for the null model, comparable with <code>deviance</code> . The null model will include the offset, and an intercept if there is one in the model. Note that this will be incorrect if the link function depends on the data other than through the fitted mean: specify a zero offset to force a correct calculation.
<code>iter</code>	the number of iterations of IWLS used.
<code>weights</code>	the <i>working</i> weights, that is the weights in the final iteration of the IWLS fit.
<code>prior.weights</code>	the weights initially supplied, a vector of 1s if none were.
<code>df.residual</code>	the residual degrees of freedom.
<code>df.null</code>	the residual degrees of freedom for the null model.
<code>y</code>	if requested (the default) the <code>y</code> vector used. (It is a vector even for a binomial model.)
<code>x</code>	if requested, the model matrix.
<code>model</code>	if requested (the default), the model frame.
<code>converged</code>	logical. Was the IWLS algorithm judged to have converged?
<code>boundary</code>	logical. Is the fitted value on the boundary of the attainable values?
<code>wle.weights</code>	final (robust) weights based on the WLE approach.
<code>wle.asymptotic</code>	logicals. If TRUE asymptotic weight based on Anscombe residual is used for the corresponding observation.

In addition, non-empty fits will have components `qr`, `R`, `qraux`, `pivot` and `effects` relating to the final weighted linear fit.

and the following components:



family	the <code>family</code> object used.
call	the matched call.
formula	the formula supplied.
terms	the <code>terms</code> object used.
data	the <code>data</code> argument.
offset	the offset vector used.
control	the value of the <code>control</code> argument used.
method	the name of the fitter function used, currently always <code>"wle.glm.fit"</code> .
contrasts	(where relevant) the contrasts used.
xlevels	(where relevant) a record of the levels of the factors used in fitting.
tot.sol	the number of solutions found.
not.conv	the number of starting points that does not converge after the <code>max.iter</code> (defined using <code>wle.glm.control</code> ) iterations are reached.
na.action	(where relevant) information returned by <code>model.frame</code> on the special handling of NAs.

Objects of class `"wle.glm"` are normally of class `"wle.glm"`.

If a `binomial` `wle.glm` model was specified by giving a two-column response, the weights returned by `prior.weights` are the total numbers of cases (factored by the supplied case weights) and the component `y` of the result is the proportion of successes.

In case of multiple roots (i.e. `tot.sol > 1`) then objects of the same form as `root1` are reported with names `root2`, `root3` and so on until `tot.sol`.

## Warnings

Since in a model selection procedure and/or on an ANOVA table the weights of the WLE procedure must be that of the FULL model (and not that of the actual model) statistics on degrees of freedom, deviance and AIC are valid only if this is the FULL model.

## Author(s)

Claudio Agostinelli and Fatemah Al-quallaf

## References

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- Agostinelli, C. and Markatou, M., (1998) A one-step robust estimator for regression based on the weighted likelihood reweighting scheme, *Statistics & Probability Letters*, Vol. 37, n. 4, 341-350.
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- Agostinelli, C. and Al-quallaf, F. (2009) Robust inference in Generalized Linear Models. Manuscript in preparation.
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Hastie, T. J. and Pregibon, D. (1992) *Generalized linear models*. Chapter 6 of *Statistical Models in S* eds J. M. Chambers and T. J. Hastie, Wadsworth & Brooks/Cole.

Markatou, M., Basu, A. and Lindsay, B.G. (1998) Weighted likelihood estimating equations with a bootstrap root search. *Journal of the American Statistical Association*, 93:740-750.

McCullagh P. and Nelder, J. A. (1989) *Generalized Linear Models*. London: Chapman and Hall.

Venables, W. N. and Ripley, B. D. (2002) *Modern Applied Statistics with S*. New York: Springer.

### See Also

[anova.wle.glm.root](#), [summary.wle.glm](#), etc. for wle.glm methods, and the generic functions [anova](#), [summary](#), [effects](#), [fitted.values](#), and [residuals](#).

[wle.lm](#) for robust non-generalized *linear* models for ‘general’ linear models.

### Examples

```
## Dobson (1990) Page 93: Randomized Controlled Trial :
counts <- c(18,17,15,20,10,20,25,13,12)
outcome <- gl(3,1,9)
treatment <- gl(3,3)
print(d.AD <- data.frame(treatment, outcome, counts))
wle.glm.D93 <- wle.glm(counts ~ outcome + treatment, family=poisson(), x=TRUE, y=TRUE)
wle.glm.D93
anova(extractRoot(wle.glm.D93))
summary(wle.glm.D93)

## Not run:
## Support for gaussian family not provided yet!
## an example with offsets from Venables & Ripley (2002, p.189)
utils::data(anorexia, package="MASS")

anorex.2 <- wle.glm(Postwt ~ Prewt + Treat + offset(Prewt),
                  family = gaussian, data = anorexia)
anorex.2
summary(anorex.2)

## End(Not run)

## Not run:
# Gamma family is not yet implemented!
# A Gamma example, from McCullagh & Nelder (1989, pp. 300-2)
clotting <- data.frame(
  u = c(5,10,15,20,30,40,60,80,100),
  lot1 = c(118,58,42,35,27,25,21,19,18),
  lot2 = c(69,35,26,21,18,16,13,12,12))
wlot1 <- wle.glm(lot1 ~ log(u), data=clotting, family=Gamma,
control=list(glm=glm.control(), wle=wle.glm.control(use.asymptotic=1)))
wlot2 <- wle.glm(lot2 ~ log(u), data=clotting, family=Gamma,
control=list(glm=glm.control(), wle=wle.glm.control(use.asymptotic=1)))
wlot1
wlot2
summary(wlot1)
```

```
summary(wlot2)

## End(Not run)
```

---

wle.glm.control

*Auxiliary for Controlling GLM Robust Fitting*


---

## Description

Auxiliary function as user interface for glm robust fitting. Typically only used when calling `wle.glm` or `wle.glm.fit`.

## Usage

```
wle.glm.control(boot = 30, group = NULL, num.sol = 1,
  raf = c("GKL", "PWD", "HD", "NED", "SCHI2"), tau = 0.1,
  cutpoint = 0, powerdown = 1, delta = NULL, smooth = NULL,
  asy.smooth=0.031, tol = 10^(-6), equal = 10^(-3),
  max.iter = 500, window.size = NULL, use.asymptotic = NULL,
  use.smooth=TRUE, mle.dispersion = FALSE, verbose = FALSE)
```

## Arguments

<code>boot</code>	integer. Number of starting points based on bootstrap subsamples to use in the search of the roots.
<code>group</code>	integer. Dimension of the bootstrap subsamples. The default value is $\max(\text{round}(\text{size}/2), \text{var} + 1)$ where <i>size</i> is the number of observations and <i>var</i> is the number of predictors.
<code>num.sol</code>	integer. Maximum number of roots to be searched.
<code>raf</code>	type of Residual adjustment function to be used: <code>raf="GKL"</code> : Generalized Kullback-Leibler family RAF (see details), <code>raf="PWD"</code> : Power Divergence family RAF (see details), <code>raf="HD"</code> : Hellinger Distance RAF, <code>raf="NED"</code> : Negative Exponential Disparity RAF, <code>raf="SCHI2"</code> : Symmetric Chi-Squared Disparity RAF.
<code>tau</code>	positive real. Used in selecting the member of the RAF family in the case of GKL or PWD.
<code>cutpoint</code>	a value in the interval [0,1].
<code>powerdown</code>	a non negative number.
<code>delta</code>	between (0,1). Used in the construction of the weights for the Binomial family.
<code>smooth</code>	the value of the smoothing parameter; used in the evaluation of weights in the case of continuous models.
<code>asy.smooth</code>	the value of the smoothing parameter; used in the evaluation of asymptotic weights. or in the case of continuous models.

tol	the absolute accuracy to be used to achieve convergence of the algorithm.
equal	the absolute value for which two roots are considered the same. Two roots are compared using the corresponding final weights.
max.iter	maximum number of iterations.
window.size	positive real or NULL. The observations with a distance, in the predictors space, less than this threshold are used to estimate the conditional distribution for a given level of the predictor.
use.asymptotic	integer or NULL. The minimum number of observations for the level of the predictors under which asymptotic weights are used.
use.smooth	if TRUE the smoothed model is used in the computation of the Pearson Residuals. For now, the option is used only for the Gamma family.
mle.dispersion	if TRUE the weighted likelihood estimator for dispersion is used otherwise the weighted chi-squared statistics is used.
verbose	if TRUE warnings are printed.

### Details

The Generalized Kullback-Leibler family RAF is defined as:

$$\ln(\tau * x + 1) / \tau$$

for  $\tau > 0$ .

The Power Divergence family RAF is defined as:

$$\tau * ((x + 1)^{1/\tau} - 1)$$

for  $0 < \tau < Inf$  while

$$\ln(x + 1)$$

for  $\tau = Inf$ .

### Value

A list with the arguments as components.

### Author(s)

Claudio Agostinelli and Fatemah Alqallaf

### References

Agostinelli, C. and Alqallaf, F. (2009) Robust inference in Generalized Linear Models. Manuscript in preparation.

### See Also

[wle.glm](#)

**Examples**

```
### A variation on example(wle.glm) :

## Annette Dobson's example ...
counts <- c(18,17,15,20,10,20,25,13,12)
outcome <- gl(3,1,9)
treatment <- gl(3,3)
oo <- options(digits = 12) # to see more when tracing :
wle.glm.D93X <- wle.glm(counts ~ outcome + treatment, family=poisson(),
                      control=list(glm=glm.control(trace = TRUE),
                                    wle=wle.glm.control(raf='GKL', tau=0.15)))
options(oo)
coef(wle.glm.D93X)
```

---

wle.glm.summaries      *Accessing Generalized Linear Model Robust Fits*

---

**Description**

These functions are all [methods](#) for class `glm` or `summary.glm` objects.

**Usage**

```
## S3 method for class 'wle.glm'
family(object, ...)

## S3 method for class 'wle.glm'
residuals(object, type = c("deviance", "pearson", "working",
                          "response", "partial"), root="all", ...)
```

**Arguments**

<code>object</code>	an object of class <code>glm</code> , typically the result of a call to <a href="#">glm</a> .
<code>type</code>	the type of residuals which should be returned. The alternatives are: "deviance" (default), "pearson", "working", "response", and "partial".
<code>root</code>	vector of integer or characters. If "all" the residuals of all roots are reported, otherwise the position of the root should be supplied.
<code>...</code>	further arguments passed to or from other methods.

**Details**

The references define the types of residuals: Davison & Snell is a good reference for the usages of each.

The partial residuals are a matrix of working residuals, with each column formed by omitting a term from the model.

How residuals treats cases with missing values in the original fit is determined by the `na.action` argument of that fit. If `na.action = na.omit` omitted cases will not appear in the residuals, whereas if `na.action = na.exclude` they will appear, with residual value NA. See also [naresid](#).

For fits done with `y = FALSE` the response values are computed from other components.

### See Also

[wle.glm](#), [anova.wle.glm.root](#); the corresponding *generic* functions, [summary.wle.glm](#), [coef](#), [deviance](#), [df.residual](#), [effects](#), [fitted](#), [residuals](#).

---

<code>wle.glm.weights</code>	<i>Weights based on Weighted Likelihood for the GLM model</i>
------------------------------	---

---

### Description

Evaluate the weights for a given GLM model

### Usage

```
wle.glm.weights(y, x, fitted.values, family = gaussian(),
dispersion = 1, raf = "GKL", tau = 0.1, smooth = NULL,
asy.smooth=0.031, window.size = NULL, use.asymptotic = NULL,
use.smooth=TRUE, tol=10^(-6), dist.method = "euclidean",
cutpoint = 0, powerdown = 1)
```

### Arguments

<code>y</code>	<code>y</code> is a vector of observations of length <code>n</code> .
<code>x</code>	<code>x</code> is a design matrix of dimension <code>n * p</code> .
<code>fitted.values</code>	the fitted mean values, obtained by transforming the linear predictors by the inverse of the link function. Often obtain as a result of <a href="#">wle.glm.fit</a> call.
<code>family</code>	a description of the error distribution and link function to be used in the model. This can be a character string naming a family function, a family function or the result of a call to a family function. (See <a href="#">family</a> for details of family functions.)
<code>dispersion</code>	value of the dispersion parameter. Used only in the Gamma family for now.
<code>raf</code>	type of Residual adjustment function to be used: <code>raf="GKL"</code> : Generalized Kullback-Leibler family RAF (see details), <code>raf="PWD"</code> : Power Divergence family RAF (see details), <code>raf="HD"</code> : Hellinger Distance RAF, <code>raf="NED"</code> : Negative Exponential Disparity RAF, <code>raf="SCHI2"</code> : Symmetric Chi-Squared Disparity RAF.
<code>tau</code>	positive real. Used in selecting the member of the RAF family in the case of GKL or PWD.
<code>smooth</code>	the value of the smoothing parameter; used in the case of continuous models.

asy.smooth	the value of the smoothing parameter; used in the evaluation of asymptotic weights.
window.size	positive real or NULL. The observations with a distance, in the predictors space, less than this threshold are used to estimate the conditional distribution for a given level of the predictor.
use.asymptotic	integer or NULL. The minimum number of observations for the level of the predictors under which asymptotic weights are used.
use.smooth	if TRUE the smoothed model is used in the computation of the Pearson Residuals. For now, the option is used only for the Gamma family.
tol	the tolerance used in the numerical calculations. For now, the option is used only for the Gamma family.
dist.method	distance method passed to <a href="#">dist</a> to measure the distance between predictor rows.
cutpoint	a value in the interval [0,1].
powerdown	a non negative number.

**Value**

A list with two components

weights	the weights associated to the observations.
asy	logical. If TRUE the corresponding weights is evaluated using asymptotic considerations based on Anscombe residuals.

**Author(s)**

Claudio Agostinelli and Fatemah Al-quallaf

**References**

Agostinelli, C. and Al-quallaf, F. (2009) Robust inference in Generalized Linear Models. Manuscript in preparation.

**See Also**

[wle.glm](#)

**Examples**

```
# tau=0.1
wgr.D93 <- extractRoot(wle.glm.D93)
# tau=0.2
w1wgr.D93 <- wle.glm.weights(y = wgr.D93$y, x = wgr.D93$x,
  fitted.values = wgr.D93$fitted.values, family = wgr.D93$family,
  raf = "GKL", tau = 0.2, smooth = 0.031, window.size = NULL,
  use.asymptotic = NULL, dist.method = "euclidean")
# tau=0.3
w2wgr.D93 <- wle.glm.weights(y = wgr.D93$y, x = wgr.D93$x,
```

```
fitted.values = wgr.D93$fitted.values, family = wgr.D93$family,
raf = "GKL", tau = 0.3, smooth = 0.031, window.size = NULL,
use.asymptotic = NULL, dist.method = "euclidean")

plot(wgr.D93$wle.weights, ylim=c(0,1), ylab='Weights')
points(w1wgr.D93$weights, col=2)
points(w2wgr.D93$weights, col=3)
legend('bottomright', legend=expression(tau==0.1, tau==0.2, tau==0.3),
pch=rep(1,3), col=1:3, inset=0.05)
```

wle.lm

*Fitting Linear Models using Weighted Likelihood***Description**

wle.lm is used to fit linear models via Weighted Likelihood, when the errors are iid from a normal distribution with null mean and unknown variance. The carriers are considered fixed. Note that this estimator is robust against the presence of bad leverage points too.

**Usage**

```
wle.lm(formula, data=list(), model=TRUE, x=FALSE,
        y=FALSE, boot=30, group, num.sol=1, raf="HD",
        smooth=0.031, tol=10-6, equal=10-3,
        max.iter=500, contrasts=NULL, verbose=FALSE)
```

**Arguments**

formula	a symbolic description of the model to be fit. The details of model specification are given below.
data	an optional data frame containing the variables in the model. By default the variables are taken from the environment which wle.lm is called from.
model, x, y	logicals. If TRUE the corresponding components of the fit (the model frame, the model matrix, the response.)
boot	the number of starting points based on bootstrap subsamples to use in the search of the roots.
group	the dimension of the bootstrap subsamples. The default value is $\max(\text{round}(\text{size}/4), \text{var})$ where <i>size</i> is the number of observations and <i>var</i> is the number of variables.
num.sol	maximum number of roots to be searched.
raf	type of Residual adjustment function to be used: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
smooth	the value of the smoothing parameter.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.



equal	the absolute value for which two roots are considered the same. (This parameter must be greater than tol).
max.iter	maximum number of iterations.
contrasts	an optional list. See the contrasts.arg of model.matrix.default.
verbose	if TRUE warnings are printed.

### Details

Models for `wle.lm` are specified symbolically. A typical model has the form `response ~ terms` where `response` is the (numeric) response vector and `terms` is a series of terms which specifies a linear predictor for response. A terms specification of the form `first+second` indicates all the terms in `first` together with all the terms in `second` with duplicates removed. A specification of the form `first:second` indicates the the set of terms obtained by taking the interactions of all terms in `first` with all terms in `second`. The specification `first*second` indicates the *cross* of `first` and `second`. This is the same as `first+second+first:second`.

### Value

`wle.lm` returns an object of class `"wle.lm"`.

The function `summary` is used to obtain and print a summary of the results. The generic accessor functions `coefficients`, `residuals` and `fitted.values` extract coefficients, residuals and fitted values returned by `wle.lm`.

The object returned by `wle.lm` are:

<code>coefficients</code>	the parameters estimator, one row vector for each root found.
<code>standard.error</code>	an estimation of the standard error of the parameters estimator, one row vector for each root found.
<code>scale</code>	an estimation of the error scale, one value for each root found.
<code>residuals</code>	the unweighted residuals from the estimated model, one column vector for each root found.
<code>fitted.values</code>	the fitted values from the estimated model, one column vector for each root found.
<code>tot.weights</code>	the sum of the weights divide by the number of observations, one value for each root found.
<code>weights</code>	the weights associated to each observation, one column vector for each root found.
<code>f.density</code>	the non-parametric density estimation.
<code>m.density</code>	the smoothed model.
<code>delta</code>	the Pearson residuals.
<code>freq</code>	the number of starting points converging to the roots.
<code>tot.sol</code>	the number of solutions found.
<code>not.conv</code>	the number of starting points that does not converge after the <code>max.iter</code> iterations are reached.
<code>call</code>	the <code>match.call()</code> .

contrasts	
xlevels	
terms	the model frame.
model	if model=TRUE a matrix with first column the dependent variable and the remain column the explanatory variables for the full model.
x	if x=TRUE a matrix with the explanatory variables for the full model.
y	if y=TRUE a vector with the dependent variable.
info	not well working yet, if 0 no error occurred.

**Author(s)**

Claudio Agostinelli

**References**

Agostinelli, C., (1998) Inferenza statistica robusta basata sulla funzione di verosimiglianza pesata: alcuni sviluppi, *Ph.D Thesis*, Department of Statistics, University of Padova.

Agostinelli, C., Markatou, M., (1998) A one-step robust estimator for regression based on the weighted likelihood reweighting scheme, *Statistics & Probability Letters*, Vol. 37, n. 4, 341-350.

Agostinelli, C., (1998) Verosimiglianza pesata nel modello di regressione lineare, *XXXIX Riunione scientifica della Societa' Italiana di Statistica*, Sorrento 1998.

**See Also**

[wle.smooth](#) an algorithm to choose the smoothing parameter for normal distribution and normal kernel.

**Examples**

```
library(wle)
# You can find this data set in:
# Hawkins, D.M., Bradu, D., and Kass, G.V. (1984).
# Location of several outliers in multiple regression data using
# elemental sets. Technometrics, 26, 197-208.
#
data(artificial)

result <- wle.lm(y.artificial~x.artificial,boot=40,num.sol=3)

summary(result)

plot(result)
```

---

wle.lm.control

*Auxiliary for Controlling LM Robust Fitting*


---

## Description

Auxiliary function as user interface for lm robust fitting.

## Usage

```
wle.lm.control(nstart = 30, group = NULL, num.sol = 1,
  raf = c("HD", "NED", "SCHI2", "GKL", "PWD"), tau = 0.1,
  cutpoint = 0, powerdown = 1, smooth = 0.031, tol = 10^(-6),
  equal = 10^(-3), max.iter = 500, verbose = FALSE)
```

## Arguments

nstart	interger. Number of starting points based on bootstrap subsamples to use in the search of the roots.
group	integer. Dimension of the bootstrap subsamples. The default value is $\max(\text{round}(\text{size}/2), \text{var} + 1)$ where <i>size</i> is the number of observations and <i>var</i> is the number of predictors.
num.sol	interger. Maximum number of roots to be searched.
raf	type of Residual adjustment function to be used: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF, raf="GKL": Generalized Kullback-Leibler family RAF (see details), raf="PWD": Power Divergence family RAF (see details).
tau	positive real. Used in selecting the member of the RAF family in the case of GKL or PWD.
cutpoint	a value in the interval [0,1].
powerdown	a non negative number.
smooth	the value of the smoothing parameter; used in the evaluation of weights in the case of continuous models.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.
equal	the absolute value for which two roots are considered the same. Two roots are compared using the corresponding final weights.
max.iter	maximum number of iterations.
verbose	if TRUE warnings are printed.

**Details**

The Generalized Kullback-Leibler family RAF is defined as:

$$\ln(\tau * x + 1) / \tau$$

for  $\tau > 0$ .

The Power Divergence family RAF is defined as:

$$\tau * ((x + 1)^{1/\tau} - 1)$$

for  $0 < \tau < \text{Inf}$  while

$$\ln(x + 1)$$

for  $\tau = \text{Inf}$ .

**Value**

A list with the arguments as components.

**Author(s)**

Claudio Agostinelli

**References**

Agostinelli, C., (1998) Inferenza statistica robusta basata sulla funzione di verosimiglianza pesata: alcuni sviluppi, *Ph.D Thesis*, Department of Statistics, University of Padova.

Agostinelli, C., Markatou, M., (1998) A one-step robust estimator for regression based on the weighted likelihood reweighting scheme, *Statistics & Probability Letters*, Vol. 37, n. 4, 341-350.

Agostinelli, C., (1998) Verosimiglianza pesata nel modello di regressione lineare, *XXXIX Riunione scientifica della Societ a Italiana di Statistica*, Sorrento 1998.

---

wle.lm.summaries

*Accessing Linear Model Fits for wle.lm*

---

**Description**

All these functions are [methods](#) for class `wle.lm` or `summary.wle.lm`.

**Usage**

```
## S3 method for class 'wle.lm'
coef(object, ...)
## S3 method for class 'wle.lm'
formula(x, ...)
## S3 method for class 'wle.lm'
fitted(object, ...)
## S3 method for class 'wle.lm'
```

```

model.frame(formula, data, na.action, ...)
## S3 method for class 'wle.lm'
summary(object, root="ALL", ...)
## S3 method for class 'wle.lm.root'
summary(object, root=1, ...)

## S3 method for class 'wle.lm'
print(x, digits = max(3, getOption("digits") - 3), ...)

## S3 method for class 'summary.wle.lm'
print(x, digits = max(3, getOption("digits") - 3),
      signif.stars= getOption("show.signif.stars"), ...)

## S3 method for class 'summary.wle.lm.root'
print(x, digits = max(3, getOption("digits") - 3),
      signif.stars= getOption("show.signif.stars"), ...)

```

### Arguments

<code>object</code>	an object of class <code>wle.lm</code> .
<code>x</code>	an object of class <code>wle.lm</code> or <code>summary.wle.lm</code> .
<code>formula</code>	a model formula
<code>data</code>	<code>data.frame</code> , list, environment or object coercible to <code>data.frame</code> containing the variables in formula.
<code>na.action</code>	how NAs are treated. The default is first, any <code>na.action</code> attribute of <code>data</code> , second a <code>na.action</code> setting of <code>options</code> , and third <code>na.fail</code> if that is unset. The “factory-fresh” default is <code>na.omit</code> .
<code>root</code>	the root to be printed, in <code>summary.wle.lm</code> it could be "ALL", all the roots are printed, or a vector of integers.
<code>digits</code>	number of digits to be used for most numbers.
<code>signif.stars</code>	logical; if TRUE, P-values are additionally encoded visually as “significance stars” in order to help scanning of long coefficient tables. It defaults to the <code>show.signif.stars</code> slot of <code>options</code> .
<code>...</code>	additional arguments.

### Details

`print.summary.wle.lm` and `print.summary.wle.lm.root` tries formatting for each root the coefficients, standard errors, etc. and additionally gives “significance stars” if `signif.stars` is TRUE. The generic accessor functions `coefficients`, `fitted.values`, `residuals` and `weights` can be used to extract various useful features of the value returned by `wle.lm`.

### Value

The function `summary.wle.lm` (the `summary.wle.lm.root` do the same for just one selected root) computes and returns, for each selected root, a list of summary statistics of the fitted linear model given in `object`, using the components (list elements) `"call"` and `"terms"` from its argument, plus

residuals	the <i>weighted</i> residuals, the usual residuals rescaled by the square root of the weights given by <code>wle.lm</code> .
coefficients	a $p \times 4$ matrix with columns for the estimated coefficient, its standard error, weighted-t-statistic and corresponding (two-sided) p-value.
sigma	the square root of the estimated variance of the random error.
df	degrees of freedom, a 3-vector $(p, \sum weights - p, p^*)$ .
fstatistic	a 3-vector with the value of the weighted-F-statistic with its numerator and denominator degrees of freedom.
r.squared	$R^2$ , the “fraction of variance explained by the model”.
adj.r.squared	the above $R^2$ statistic “ <i>adjusted</i> ”, penalizing for higher $p$ .
root	the label of the root reported.

**Author(s)**

Claudio Agostinelli

**See Also**

[wle.lm](#) a function for estimating linear models with normal distribution error and normal kernel, [plot.wle.lm](#) for plot method.

**Examples**

```
library(wle)
# You can find this data set in:
# Hawkins, D.M., Bradu, D., and Kass, G.V. (1984).
# Location of several outliers in multiple regression data using
# elemental sets. Technometrics, 26, 197-208.
#
data(artificial)

result <- wle.lm(y.artificial~x.artificial,boot=40,group=6,num.sol=3)

#summary only for the first root
summary(result,root=1)
#summary for all the roots
summary(result,root="ALL")
```

---

wle.negativebinomial    *Robust Estimation in the Negative Binomial Model*

---

**Description**

`wle.negativebinomial` is used to robust estimate the proportion parameters via Weighted Likelihood.

**Usage**

```
wle.negativebinomial(x, size, boot=30, group, num.sol=1,
  raf="HD", tol=10^(-6), equal=10^(-3),
  max.iter=500, verbose=FALSE)
```

**Arguments**

x	a vector contain the number of failures which occur in a sequence of Bernoulli trials before a target number of successes size is reached.
size	target number of successes.
boot	the number of starting points based on bootstrap subsamples to use in the search of the roots.
group	the dimension of the bootstrap subsamples. The default value is $\max(\text{round}(\text{length}(x)/4), 2)$ .
num.sol	maximum number of roots to be searched.
raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.
equal	the absolute value for which two roots are considered the same. (This parameter must be greater than tol).
max.iter	maximum number of iterations.
verbose	if TRUE warnings are printed.

**Value**

wle.negativebinomial returns an object of class "wle.negativebinomial".

Only print method is implemented for this class.

The object returned by wle.negativebinomial are:

p	the estimator of the proportion parameter, one value for each root found.
tot.weights	the sum of the weights divide by the number of observations, one value for each root found.
weights	the weights associated to each observation, one column vector for each root found.
f.density	the non-parametric density estimation.
m.density	the smoothed model.
delta	the Pearson residuals.
call	the match.call().
tot.sol	the number of solutions found.
not.conv	the number of starting points that does not converge after the max.iter iteration are reached.

**Author(s)**

Claudio Agostinelli

**References**

Markatou, M., Basu, A., and Lindsay, B.G., (1997) Weighted likelihood estimating equations: The discrete case with applications to logistic regression, *Journal of Statistical Planning and Inference*, 57, 215-232.

Agostinelli, C., (1998) Inferenza statistica robusta basata sulla funzione di verosimiglianza pesata: alcuni sviluppi, *Ph.D Thesis*, Department of Statistics, University of Padova.

**Examples**

```
library(wle)

set.seed(1234)

x <- rbinom(20, size=10, prob=0.2)
wle.negativebinomial(x, size=10)

x <- c(rnbinom(20, size=10, prob=0.2),rnbinom(10, size=10, p=0.9))
result <- wle.negativebinomial(x, size=10)
print(result)
plot(result$weights)
```

---

wle.normal

---

*Robust Estimation in the Normal Model*


---

**Description**

wle.normal is used to robust estimate the location and the scale parameters via Weighted Likelihood, when the sample is iid from a normal distribution with unknown mean and variance.

**Usage**

```
wle.normal(x, boot=30, group, num.sol=1, raf="HD",
           smooth=0.003, tol=10-6, equal=10-3,
           max.iter=500, verbose=FALSE)
```

**Arguments**

x	a vector contain the observations.
boot	the number of starting points based on bootstrap subsamples to use in the search of the roots.
group	the dimension of the bootstrap subsamples. The default value is $\max(\text{round}(\text{size}/4), 2)$ where <i>size</i> is the number of observations.
num.sol	maximum number of roots to be searched.



raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
smooth	the value of the smoothing parameter.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.
equal	the absolute value for which two roots are considered the same. (This parameter must be greater than tol).
max.iter	maximum number of iterations.
verbose	if TRUE warnings are printed.

### Value

wle.normal returns an object of class "wle.normal".

Only print method is implemented for this class.

The object returned by wle.normal are:

location	the estimator of the location parameter, one value for each root found.
scale	the estimator of the scale parameter, one value for each root found.
residuals	the residuals associated to each observation, one column vector for each root found.
tot.weights	the sum of the weights divide by the number of observations, one value for each root found.
weights	the weights associated to each observation, one column vector for each root found.
f.density	the non-parametric density estimation.
m.density	the smoothed model.
delta	the Pearson residuals.
freq	the number of starting points converging to the roots.
call	the match.call().
tot.sol	the number of solutions found.
not.conv	the number of starting points that does not converge after the max.iter iteration are reached.

### Author(s)

Claudio Agostinelli

### References

Markatou, M., Basu, A. and Lindsay, B.G., (1998) Weighted likelihood estimating equations with a bootstrap root search, *Journal of the American Statistical Association*, 93, 740-750.

Agostinelli, C., (1998) Inferenza statistica robusta basata sulla funzione di verosimiglianza pesata: alcuni sviluppi, *Ph.D Thesis*, Department of Statistics, University of Padova.



**Arguments**

x	a vector contain the observations.
m	numbers of components.
boot	the number of starting points based on bootstrap subsamples to use in the search of the roots.
group	the dimension of the bootstrap subsamples. The default value is $\max(\text{round}(\text{size}/4), 2)$ where <i>size</i> is the number of observations.
num.sol	maximum number of roots to be searched.
raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
smooth	the value of the smoothing parameter.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.
equal	the absolute value for which two roots are considered the same. (This parameter must be greater than tol).
max.iter	maximum number of iterations.
all.comp	try to find all the components.
min.size	see details
min.weights	see details
boot.start	the number of starting points for the starting process.
group.start	the dimension of the bootstrap subsamples in the starting process. The default value is $\max(\text{round}(\text{group}/4), 2)$ .
tol.start	the absolute accuracy to be used to achieve convergence of the algorithm in the starting process.
equal.start	the absolute value for which two roots are considered the same in the starting process. (This parameter must be greater than tol.start).
smooth.start	the value of the smoothing parameter in the starting process.
max.iter.start	maximum number of iterations in the starting process.
max.iter.boot	maximum number of iterations of the starting process.
verbose	if TRUE warnings are printed.

**Details**

this function use an iterative procedure to evaluate starting points. First, using `wle.normal` we try to find the biggest components, then we discard each observation with weight greater than `min.weights`. The `wle.normal` is run on the remain observations if the ratio between the number of observations and the original sample size is greater than `min.size`. The convergence of the algorithm is determined by the difference between two iterations. This stopping rule could have some problems, as soon as possible it will replace with the one proposed in Markatou (2000) pag. 485 (5).

**Value**

wle.normal.mixture returns an object of class "wle.normal.mixture".

Only print method is implemented for this class.

The objects returned by wle.normal.mixture are:

location	the estimator of the location parameters, one vector for each root found.
scale	the estimator of the scale parameters, one vector for each root found.
pi	the estimator of the proportion parameters, one vector for each root found.
tot.weights	the sum of the weights, divide by the number of observations, one value for each root found.
weights	the weights associated to each observation, one column vector for each root found.
f.density	the non-parametric density estimation.
m.density	the smoothed model.
delta	the Pearson residuals.
freq	the number of starting points converging to the roots.
tot.sol	the number of solutions found.
not.conv	the number of starting points that does not converge after the max.iter iteration are reached.
call	the match.call().

**Author(s)**

Claudio Agostinelli

**References**

Markatou, M., (2000) Mixture models, robustness and the weighted likelihood methodology, *Biometrics*, 56, 483-486.

Markatou, M., (2001) A closer look at the weighted likelihood in the context of mixtures, *Probability and Statistical Models with Applications*, Charalambides, C.A., Koutras, M.V. and Balakrishnan, N. (eds.), Chapman and Hall/CRC, 447-467.

**Examples**

```
library(wle)
set.seed(1234)
x <- c(rnorm(150,0,1),rnorm(50,15,2))
wle.normal.mixture(x,m=2,group=50,group.start=2,boot=5,num.sol=3)
wle.normal(x,group=2,boot=10,num.sol=3)
```

**Description**

wle.normal.multi is used to robust estimate the location and the covariance matrix via Weighted Likelihood, when the sample is iid from a normal multivariate distribution with unknown means and variance matrix.

**Usage**

```
wle.normal.multi(x, boot=30, group, num.sol=1,
                 raf="HD", smooth, tol=10-6,
                 equal=10-3, max.iter=500,
                 verbose=FALSE)
```

**Arguments**

x	a matrix contain the observations.
boot	the number of starting points based on bootstrap subsamples to use in the search of the roots.
group	the dimension of the bootstrap subsamples. The default value is $\max(\text{round}(\text{size}/4), (\text{var} * (\text{var} + 1)/2 + \text{var}))$ where <i>size</i> is the number of observations and <i>var</i> is the number of variables.
num.sol	maximum number of roots to be searched.
raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
smooth	the value of the smoothing parameter.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.
equal	the absolute value for which two roots are considered the same. (This parameter must be greater than tol).
max.iter	maximum number of iterations.
verbose	if TRUE warnings are printed.

**Value**

wle.normal.multi returns an object of class "wle.normal.multi".

Only print method is implemented for this class.

The object returned by wle.normal.multi are:

location	the estimator of the location parameters, one vector for each root found.
----------	---



```
barplot(result$weights,col=2,xlab="Observations",
        ylab="Weights",ylim=c(0,1),
        names.arg=seq(1:length(result$weights)))
```

---

wle.normal.multi.summaries

*Summaries and methods for wle.normal.multi*

---

### Description

Until now, only print [methods](#) is available for class `wle.normal.multi`. `print.wle.normal.multi` print nicely the output.

### Usage

```
## S3 method for class 'wle.normal.multi'
print(x, digits = max(3, getOption("digits") - 3), ...)
```

### Arguments

<code>x</code>	an object of class <code>wle.normal.multi</code> .
<code>digits</code>	number of digits to be used for most numbers.
<code>...</code>	further arguments passed to or from other methods.

### Author(s)

Claudio Agostinelli

### See Also

[wle.normal.multi](#) a function for estimating normal multivariate location and scale models.

---

wle.normal.summaries

*Summaries and methods for wle.normal*

---

### Description

Until now, only print [methods](#) is available for class `wle.normal`. `print.wle.normal` print nicely the output.

### Usage

```
## S3 method for class 'wle.normal'
print(x, digits = max(3, getOption("digits") - 3), ...)
```

**Arguments**

x	an object of class <code>wle.normal</code> .
digits	number of digits to be used for most numbers.
...	further arguments passed to or from other methods.

**Author(s)**

Claudio Agostinelli

**See Also**

[wle.normal](#) a function for estimating normal location and scale models.

---

wle.onestep

*A One-Step Weighted Likelihood Estimator for Linear model*

---

**Description**

This function evaluate the One-step weighted likelihood estimator for the regression and scale parameters.

**Usage**

```
wle.onestep(formula, data=list(), model=TRUE, x=FALSE,
             y=FALSE, ini.param, ini.scale, raf="HD",
             smooth=0.031, num.step=1,
             contrasts=NULL, verbose=FALSE)
```

**Arguments**

formula	a symbolic description of the model to be fit. The details of model specification are given below.
data	an optional data frame containing the variables in the model. By default the variables are taken from the environment which <code>wle.stepwise</code> is called from.
model, x, y	logicals. If TRUE the corresponding components of the fit (the model frame, the model matrix, the response.)
ini.param	starting values for the coefficients.
ini.scale	starting values for the scale parameters.
raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
smooth	the value of the smoothing parameter.
num.step	number of the steps.
contrasts	an optional list. See the <code>contrasts.arg</code> of <code>model.matrix.default</code> .
verbose	if TRUE warnings are printed.



**Value**

wle.onestep returns an object of class "wle.onestep".

Only print method is implemented for this class.

The object returned by wle.onestep are:

coefficients	the parameters estimator.
standard.error	an estimation of the standard error of the parameters estimator.
scale	an estimation of the error scale.
residuals	the unweighted residuals from the estimated model.
fitted.values	the fitted values from the estimated model.
tot.weights	the sum of the weights divide by the number of observations.
weights	the weights associated to each observation.
f.density	the non-parametric density estimation.
m.density	the smoothed model.
delta	the Pearson residuals.
call	the match.call().
contrasts	
xlevels	
terms	the model frame.
model	if model=TRUE a matrix with first column the dependent variable and the remain column the explanatory variables for the full model.
x	if x=TRUE a matrix with the explanatory variables for the full model.
y	if y=TRUE a vector with the dependent variable.

**Author(s)**

Claudio Agostinelli

**References**

Agostinelli, C., (1997) A one-step robust estimator based on the weighted likelihood methodology, *Working Paper n. 1997.16*, Department of Statistics, University of Padova.

Agostinelli, C., (1998) Inferenza statistica robusta basata sulla funzione di verosimiglianza pesata: alcuni sviluppi, *Ph.D Thesis*, Department of Statistics, University of Padova.

Agostinelli, C., Markatou, M., (1998) A one-step robust estimator for regression based on the weighted likelihood reweighting scheme, *Statistics & Probability Letters*, Vol. 37, n. 4, 341-350.

Agostinelli, C., (1998) Verosimiglianza pesata nel modello di regressione lineare, *XXXIX Riunione scientifica della Societ a Italiana di Statistica*, Sorrento 1998.

**See Also**

[wle.smooth](#) an algorithm to choose the smoothing parameter for normal distribution and normal kernel, [wle.lm](#) a function for estimating linear models with normal distribution error and normal kernel.

## Examples

```
#library(wle)
#library(lqs)

#data(artificial)

#result.lts <- lqs(y.artificial~x.artificial,
#                 method = "lts")

#result.wle <- wle.onestep(y.artificial~x.artificial,
#                          ini.param=result.lts$coefficients,
#                          ini.scale=result.lts$scale[1])

#result.wle
```

---

wle.onestep.summaries *Summaries and methods for wle.onestep*

---

## Description

Until now, only print [methods](#) is available for class `wle.onestep`. `print.wle.onestep` print nicely the output.

## Usage

```
## S3 method for class 'wle.onestep'
print(x, digits = max(3, getOption("digits") - 3), ...)
```

## Arguments

<code>x</code>	an object of class <code>wle.onestep</code> .
<code>digits</code>	number of digits to be used for most numbers.
<code>...</code>	further arguments passed to or from other methods.

## Author(s)

Claudio Agostinelli

## See Also

[wle.onestep](#) a function for one-step estimation in linear models.

**Description**

wle.poisson is used to robust estimate the lambda parameters in the poisson model via Weighted Likelihood.

**Usage**

```
wle.poisson(x, boot=30, group, num.sol=1,
  raf=c("HD", "NED", "GKL", "PWD", "SCHI2"),
  tau=NULL, tol=10-6, equal=10-3,
  max.iter=500, verbose=FALSE)
```

**Arguments**

x	a vector contain the number of success.
boot	the number of starting points based on bootstrap subsamples to use in the search of the roots.
group	the dimension of the bootstrap subsamples. The default value is $\max(\text{round}(\text{length}(x)/4), 2)$ .
num.sol	maximum number of roots to be searched.
raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="GKL": Generalized Kullback-Leibler RAF family with parameter tau. raf="PWD": Power Divergence Measure RAF family with parameter tau. raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
tau	this is to set the member inside the GKL and PWD family. It must be in [0,1] for GKL and in [-1, Inf] for PWD.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.
equal	the absolute value for which two roots are considered the same. (This parameter must be greater than tol).
max.iter	maximum number of iterations.
verbose	if TRUE warnings are printed.

**Value**

wle.poisson returns an object of `class` "wle.poisson".

Only print method is implemented for this class.

The object returned by wle.poisson are:

lambda	the estimator of the lambda parameter, one value for each root found.
--------	---

tot.weights	the sum of the weights divide by the number of observations, one value for each root found.
weights	the weights associated to each observation, one column vector for each root found.
f.density	the non-parametric density estimation.
m.density	the smoothed model.
delta	the Pearson residuals.
call	the match.call().
tot.sol	the number of solutions found.
not.conv	the number of starting points that does not converge after the max.iter iteration are reached.

### Author(s)

Claudio Agostinelli

### References

Markatou, M., Basu, A., and Lindsay, B.G., (1997) Weighted likelihood estimating equations: The discrete case with applications to logistic regression, *Journal of Statistical Planning and Inference*, 57, 215-232.

Agostinelli, C., (1998) Inferenza statistica robusta basata sulla funzione di verosimiglianza pesata: alcuni sviluppi, *Ph.D Thesis*, Department of Statistics, University of Padova.

### Examples

```
library(wle)

set.seed(1234)

x <- rpois(40,5)
wle.poisson(x)

x <- c(rpois(40,5),rpois(10,20))
wle.poisson(x)
```

---

wle.smooth

*Bandwidth selection for the normal kernel and normal model.*

---

### Description

The bandwidth of the kernel is choose for normal model and normal kernel in such a way a contaminated point costant times away from the mean of the distribution in scale units and mass level has a weight no bigger than weight.

## Usage

```
wle.smooth(weight=0.31, costant=3, level=0.2,  
           dimension=1, raf="HD", interval=c(0.00001, 0.5),  
           tol=10^-6, max.iter=1000)
```

## Arguments

weight	weights associated to an observation that is constant scale units away from the mean of the distribution.
constant	times the contaminated point mass is away from the mean of the distribution in scale units.
level	mass of the contaminated point.
dimension	dimension of the normal distribution.
raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
interval	interval from which to search the root.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.
max.iter	maximum number of iterations.

## Details

The `wle.smooth` use `uniroot` function to solve the non linear equation. No handling error is provided yet. For the Symmetric Chi-Squared Disparity RAF you should use `weight=0.2` and `interval=c(0.1, 1)` to have a solution.

## Value

`wle.smooth` returns an object of `class` "wle.smooth".

Only `print` method is implemented for this class.

The object returned by `wle.smooth` is a list with four components: `root` and `f.root` give the location of the root and the value of the function evaluated at that point. `iter` and `estim.prec` give the number of iterations used and an approximate estimated precision for root.

`root` is the value of the bandwidth.

## Author(s)

Claudio Agostinelli

## References

- Agostinelli, C., (1998) Inferenza statistica robusta basata sulla funzione di verosimiglianza pesata: alcuni sviluppi, *Ph.D. thesis*, Department of Statistics, University of Padova.
- Markatou, M., Basu, A. and Lindsay, B.G. (1998) Weighted likelihood estimating equations with a bootstrap root search. *Journal of the American Statistical Association*, 93, 740-750.
- Agostinelli, C., and Markatou, M., (2001) Test of hypotheses based on the Weighted Likelihood Methodology, *Statistica Sinica*, vol. 11, n. 2, 499-514.

## See Also

[uniroot](#), uniroot: one dimensional root finding.

## Examples

```
library(wle)

wle.smooth()
```

---

wle.stepwise

*Weighted Stepwise, Backward and Forward selection methods*

---

## Description

This function performs Weighted Stepwise, Forward and Backward model selection.

## Usage

```
wle.stepwise(formula, data=list(), model=TRUE, x=FALSE,
             y=FALSE, boot=30, group, num.sol=1, raf="HD",
             smooth=0.031, tol=10(-6), equal=10(-3),
             max.iter=500, min.weight=0.5, type="Forward",
             f.in=4.0, f.out=4.0, method="WLE",
             contrasts=NULL, verbose=FALSE)
```

## Arguments

- |             |  |
|-------------|--|
| formula     | a symbolic description of the model to be fit. The details of model specification are given below.   |
| data        | an optional data frame containing the variables in the model. By default the variables are taken from the environment which <code>wle.stepwise</code> is called from.                                      |
| model, x, y | logicals. If TRUE the corresponding components of the fit (the model frame, the model matrix, the response.)   |
| boot        | the number of starting points based on bootstrap subsamples to use in the search of the roots.   |
| group       | the dimension of the bootstrap subsamples. The default value is $\max(\text{round}(\text{size}/4), \text{var})$ where <i>size</i> is the number of observations and <i>var</i> is the number of variables. |

num.sol	maximum number of roots to be searched.
raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
smooth	the value of the smoothing parameter.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.
equal	the absolute value for which two roots are considered the same. (This parameter must be greater than tol).
max.iter	maximum number of iterations.
min.weight	see details.
type	type="Stepwise": the weighted stepwise methods is used, type="Forward": the weighted forward methods is used, type="Backward": the weighted backward method is used.
f.in	the in value
f.out	the out value
method	method="WLS": the submodel parameters are estimated by weighted least square with weights from the weighted likelihood estimator on the full model. method="WLE": the submodel parameters are estimated by weighted likelihood estimators.
contrasts	an optional list. See the contrasts.arg of model.matrix.default.
verbose	if TRUE warnings are printed.

## Details

Models for `wle.stepwise` are specified symbolically. A typical model has the form `response ~ terms` where `response` is the (numeric) response vector and `terms` is a series of terms which specifies a linear predictor for response. A terms specification of the form `first+second` indicates all the terms in `first` together with all the terms in `second` with duplicates removed. A specification of the form `first:second` indicates the the set of terms obtained by taking the interactions of all terms in `first` with all terms in `second`. The specification `first*second` indicates the *cross* of `first` and `second`. This is the same as `first+second+first:second`.

`min.weight`: the weighted likelihood equation could have more than one solution. These roots appear for particular situation depending on contamination level and type. The presence of multiple roots in the full model can create some problem in the set of weights we should use. Actually, the selection of the root is done by the minimum scale error provided. Since this choice is not always the one would choose, we introduce the `min.weight` parameter in order to choose only between roots that do not down weight everything. This is not still the optimal solution, and perhaps, in the new release, this part will be change.

**Value**

wle.stepwise returns an object of class "wle.stepwise".

The function summary is used to obtain and print a summary of the results. The generic accessor functions coefficients and residuals extract coefficients and residuals returned by wle.stepwise.

The object returned by wle.stepwise are:

wstep	the iterations with the model selected.
coefficients	the parameters estimator, one row vector for each root found in the full model.
scale	an estimation of the error scale, one value for each root found in the full model.
residuals	the unweighted residuals from the estimated model, one column vector for each root found in the full model.
tot.weights	the sum of the weights divide by the number of observations, one value for each root found in the full model.
weights	the weights associated to each observation, one column vector for each root found in the full model.
freq	the number of starting points converging to the roots.
index	position of the root used for the weights.
call	the match.call().
contrasts	
xlevels	
terms	the model frame.
model	if model=TRUE a matrix with first column the dependent variable and the remain column the explanatory variables for the full model.
x	if x=TRUE a matrix with the explanatory variables for the full model.
y	if y=TRUE a vector with the dependent variable.
info	not well working yet, if 0 no error occurred.
type	"Stepwise": the weighted stepwise methods is used, "Forward": the weighted forward methods is used, "Backward": the weighted backward method is used.
f.in	the in value.
f.out	the out value.
method	if "WLS" the submodel parameters are estimated by weighted least square with weights from the weighted likelihood estimator on the full model else if "WLE" the submodel parameters are estimated by weighted likelihood estimators.

**Author(s)**

Claudio Agostinelli



## References

Agostinelli, C., (2000) Robust stepwise regression, Working Paper n. 2000.10 del Dipartimento di Scienze Statistiche, Università di Padova, Padova.

Agostinelli, C., (2002) Robust stepwise regression, *Journal of Applied Statistics* 29, 6, 825-840.

Agostinelli, C., (1998) Inferenza statistica robusta basata sulla funzione di verosimiglianza pesata: alcuni sviluppi, *Ph.D Thesis*, Department of Statistics, University of Padova.

Agostinelli, C., (1998) Verosimiglianza pesata nel modello di regressione lineare, *XXXIX Riunione scientifica della Società Italiana di Statistica*, Sorrento 1998.

## See Also

[wle.smooth](#) an algorithm to choose the smoothing parameter for normal distribution and normal kernel, [wle.lm](#) a function for estimating linear models with normal distribution error and normal kernel.

## Examples

```
library(wle)

# You can find this dataset in:
# Agostinelli, C., (2002). Robust model selection in regression
# via weighted likelihood methodology, Statistics &
# Probability Letters, 56, 289-300.

data(selection)

result <- wle.stepwise(ydata~xdata, boot=100, group=6, num.sol=3,
min.weight=0.8, type="Stepwise", method="WLS")

summary(result)
```

---

```
wle.stepwise.summaries
```

*Accessing summaries for wle.stepwise*

---

## Description

All these functions are [methods](#) for class `wle.stepwise` or `summary.wle.stepwise`.

## Usage

```
## S3 method for class 'wle.stepwise'
summary(object, num.max=20, verbose=FALSE, ...)

## S3 method for class 'wle.stepwise'
print(x, digits = max(3, getOption("digits")) -
```

```
3), num.max=max(1, nrow(x$wstep)), ...)

## S3 method for class 'summary.wle.stepwise'
print(x, digits = max(3, getOption("digits") - 3), ...)
```

### Arguments

<code>object</code>	an object of class <code>wle.stepwise</code> .
<code>x</code>	an object of class <code>wle.stepwise</code> or <code>summary.wle.stepwise</code> .
<code>digits</code>	number of digits to be used for most numbers.
<code>num.max</code>	the number of the last iterations reported.
<code>verbose</code>	if TRUE warnings are printed.
<code>...</code>	additional arguments affecting the summary produced (in <code>summary.wle.stepwise</code> ) or further arguments passed to or from other methods (in <code>print.wle.stepwise</code> and <code>print.summary.wle.stepwise</code> ).

### Details

The generic accessor functions `coefficients`, `fitted.values`, `residuals` and `weights` can be used to extract various useful features of the value returned by `wle.stepwise`.

### Value

The function `summary.wle.stepwise` returns the last `num.max` iterations, call plus:

<code>wstep</code>	the model for each iteration reported.
<code>num.max</code>	the number of iterations reported.
<code>type</code>	the type of selection procedure used.
<code>f.in</code>	the in value.
<code>f.out</code>	the out value.

### Author(s)

Claudio Agostinelli

---

wle.t.test

*Weighted Likelihood Student's t-Test*

---

### Description

`wle.t.test` performs one and two sample Weighted Likelihood t-tests on vectors of data. This is a robust version of the classical t-test. It should be used when the majority of the data follows a normal distribution.

**Usage**

```
wle.t.test(x, y = NULL, alternative = c("two.sided", "less", "greater"),
  mu = 0, paired = FALSE, var.equal = FALSE, conf.level = 0.95,
  boot=30, group, num.sol=1, raf="HD", smooth=0.003,
  tol=10-6, equal=10-3, max.iter=500)
```

**Arguments**

x	a numeric vector of data values.
y	an optional numeric vector data values.
alternative	character specifying the alternative hypothesis, must be one of "two.sided" (default), "greater" or "less". You can specify just the initial letter.
mu	a number indicating the true value of the mean (or difference in means if you are performing a two sample test).
paired	a logical indicating whether you want a paired weighted t-test.
var.equal	a logical variable indicating whether to treat the two variances as being equal. If TRUE then the pooled variance is used to estimate the variance otherwise the Welch approximation to the degrees of freedom is used.
conf.level	confidence level of the interval.
boot	the number of starting points based on bootstrap subsamples to use in the search of the roots.
group	the dimension of the bootstrap subsamples. The default value is $\max(\text{round}(\text{size}/4), 2)$ where <i>size</i> is the number of observations.
num.sol	maximum number of roots to be searched.
raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
smooth	the value of the smoothing parameter.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.
equal	the absolute value for which two roots are considered the same. (This parameter must be greater than tol).
max.iter	maximum number of iterations.

**Details**

If paired is TRUE then both x and y must be specified and they must be the same length. Missing values are removed (in pairs if paired is TRUE). If var.equal is TRUE then the pooled estimate of the variance is used. By default, if var.equal is FALSE then the variance is estimated separately for both groups and the Welch modification to the degrees of freedom is used.

**Value**

The function return a list of class "wle. t. test" with the following components:

test	A list with two dimensions. Each cell is related with a combination of 'x', 'y' roots. In each cell a list of class "htest" containing the following components: statistic the value of the t-statistic. parameters the degrees of freedom for the t-statistic. p.value the p-value for the test. conf.int a confidence interval for the mean appropriate to the specified alternative hypothesis. estimate the estimated mean or difference in means depending on whether it was a one-sample test or a two-sample test. null.value the specified hypothesized value of the mean or mean difference depending on whether it was a one-sample test or a two-sample test. alternative a character string describing the alternative hypothesis. method a character string indicating what type of t-test was performed. data.name a character string giving the name(s) of the data. x.weights the weights related to the 'x' data. y.weights the weights related to the 'y' data. x.root the number of the 'x' root. y.root the number of the 'y' root.
x.tot.sol	the number of solutions for the dataset 'x'.
y.tot.sol	the number of solutions for the dataset 'y' or 1.
call	the match.call().
paired	a logical indicating whether is a paired weighted t-test.
x	'x' data.
y	'y' data or NULL.

**Author(s)**

Claudio Agostinelli

**References**

- Agostinelli, C., (1998) Inferenza statistica robusta basata sulla funzione di verosimiglianza pesata: alcuni sviluppi, *Ph.D Thesis*, Department of Statistics, University of Padova (in italian).
- Agostinelli, C., (2002) Un approccio alla verifica d'ipotesi robusta basato sulla funzione di verosimiglianza pesata - Robust Testing Hypotheses via Weighted Likelihood function, *Statistica*, Anno LXII, 1, 87-110.
- Agostinelli, C., and Markatou, M., (2001) Test of hypotheses based on the Weighted Likelihood Methodology, *Statistica Sinica*, vol. 11, n. 2, 499-514.

**Examples**

```

library(wle)

set.seed(1234)

x <- rnorm(20,0,1)
y <- rnorm(20,6,1)

t.test(x,y)           # P < 2.2e-16
wle.t.test(x,y,group=5) # P < 2.2e-16

t.test(x,y=c(y,250))   # P = 0.1419 -- NOT significant anymore
wle.t.test(x,y=c(y,250),group=5) # P < 2.2e-16 -- still significant
set.seed(1234)

# three roots for 'x' and three roots for 'y'
# with nine t-test value
res <- wle.t.test(x=c(rnorm(40,0,1),rnorm(40,10,1)),
                  y=c(rnorm(40,0,1),rnorm(40,10,1)),
                  group=4,num.sol=3,boot=100)

print(res) # print ALL the t-test
print(res,x.root=1,y.root=1) # print the test associated to the
                              # x.root=1,y.root=1

root.1.1 <- res$test[[1]][[1]] # access to the object associated
                              # to the x.root=1,y.root=1

names(root.1.1)

set.seed(1234)

# one root and NOT significant t-test
wle.t.test(x=c(rnorm(40,0,1),rnorm(40,10,1)),
           y=c(rnorm(40,0,1),rnorm(40,10,1)),
           group=4,num.sol=3,boot=100,paired=TRUE)

```

---

wle.var.test

*Weighted F Test to Compare Two Variances*


---

**Description**

Performs an Weighted F test to compare the variances of two samples from normal populations. The WF-test is based on weighted likelihood.

**Usage**

```

wle.var.test(x, y, ratio = 1, alternative = c("two.sided", "less", "greater"),
             conf.level = 0.95, x.root=1, y.root=1)

```

**Arguments**

x, y	fitted linear model objects (inheriting from class "wle.lm") or fitted normal model objects (inheriting from class "wle.normal").
ratio	the hypothesized ratio of the population variances of x and y.
alternative	the alternative hypothesis; must be one of "two.sided" (default), "greater" or "less". You can specify just the initial letter.
conf.level	confidence level for the returned confidence interval.
x.root	the 'x' root used.
y.root	the 'y' root used.

**Details**

The null hypothesis is that the ratio of the variances in the data to which the normal model ([wle.normal](#)) or linear models ([wle.lm](#)) x and y were fitted, is equal to ratio.

**Value**

A list with class "hctest" containing the following components:

statistic	the value of the WF test statistic.
parameter	the degrees of the freedom of the WF distribution of the test statistic.
p.value	the p-value of the test.
conf.int	a confidence interval for the ratio of the population variances.
estimate	the ratio of the sample variances from x and y.
null.value	the ratio of population variances under the null.
alternative	a character string describing the alternative hypothesis.
method	the string "WF test to compare two variances".
data.name	a character string giving the names of the data.

**Author(s)**

Claudio Agostinelli

**References**

- Agostinelli, C., (1998). Inferenza statistica robusta basata sulla funzione di verosimiglianza pesata: alcuni sviluppi, *Ph.D Thesis*, Department of Statistics, University of Padova (in italian).
- Agostinelli, C., (2001) Un approccio robusto alla verifica d'ipotesi basato sulla funzione di verosimiglianza pesata - Robust Testing Hypotheses via Weighted Likelihood function, in press *Statistica*, (in italian).
- Agostinelli, C., and Markatou, M., (2001) Test of hypotheses based on the Weighted Likelihood Methodology, *Statistica Sinica*, vol. 11, n. 2, 499-514.

**Examples**

```

set.seed(2345)

x <- rnorm(50,0,1)
y <- rnorm(50,10,1)

res.x <- wle.normal(x,group=5)
res.y <- wle.normal(y,group=5)

wle.var.test(res.x, res.y) # Do x and y have the same variance?

set.seed(2345)

x <- c(rnorm(50,0,1),rnorm(20,10,1))
y <- c(rnorm(50,10,1),rnorm(10,0,5))

res.x <- wle.normal(x,group=5,num.sol=2)
res.y <- wle.normal(y,group=5)

res.x
wle.var.test(res.x, res.y, x.root=1)
if (res.x$tot.sol>1) wle.var.test(res.x, res.y, x.root=2)

```

---

wle.vonmises

*von Mises Weighted Likelihood Estimates*


---

**Description**

Computes the weighted likelihood estimates for the parameters of a von Mises distribution: the mean direction and the concentration parameter.

**Usage**

```

wle.vonmises(x, boot = 30, group, num.sol = 1, raf = "HD", smooth, tol =
10-6, equal = 10-3, max.iter = 500, bias = FALSE, mle.bias =
FALSE, max.kappa = 500, min.kappa = 0.01, use.smooth = TRUE, alpha =
NULL, p = 2, verbose = FALSE, control.circular = list())
## S3 method for class 'wle.vonmises'
print(x, digits = max(3, getOption("digits") - 3), ...)

```

**Arguments**

x	a vector. The object is coerced to class <a href="#">circular</a> .
boot	the number of starting points based on bootstrap subsamples to use in the search of the roots.
group	the dimension of the bootstrap subsamples.

<code>num.sol</code>	maximum number of roots to be searched.
<code>raf</code>	type of Residual adjustment function to be use: <code>raf="HD"</code> : Hellinger Distance RAF, <code>raf="NED"</code> : Negative Exponential Disparity RAF, <code>raf="SCHI2"</code> : Symmetric Chi-Squared Disparity RAF.
<code>smooth</code>	the value of the smoothing parameter.
<code>tol</code>	the absolute accuracy to be used to achieve convergence of the algorithm.
<code>equal</code>	the absolute value for which two roots are considered the same. (This parameter must be greater than <code>tol</code> ).
<code>max.iter</code>	maximum number of iterations.
<code>bias</code>	logical, if TRUE, the estimate for kappa is computed with a bias corrected method. Default is FALSE, i.e. no bias correction.
<code>mle.bias</code>	logical, if TRUE a bias corrected method is used to estimate the concentration parameter for the initial values.
<code>max.kappa</code>	maximum value for the concentration parameter.
<code>min.kappa</code>	minimum value for the concentration parameter.
<code>use.smooth</code>	logical, if TRUE a smoothed model is used, default is TRUE.
<code>alpha</code>	if not NULL overrides the value of <code>p</code> . See the next argument <code>p</code> . This is a different parameterization, <code>alpha=-1/2</code> provides Hellinger Distance RAF, <code>alpha=-1</code> provides Kullback-Leibler RAF and <code>alpha=-2</code> provides Neyman's Chi-Square RAF.
<code>p</code>	this parameter works only when <code>raf="HD"</code> . <code>p=2</code> provides Hellinger Distance RAF, <code>p=-1</code> provides Kullback-Leibler RAF and <code>p=Inf</code> provides Neyman's Chi-Square RAF.
<code>verbose</code>	logical, if TRUE warnings are printed.
<code>control.circular</code>	the attribute of the resulting object ( <code>mu</code> )
<code>digits</code>	integer indicating the precision to be used.
<code>...</code>	further parameters in <code>print.wle.vonmises</code> .

### Details

Parameters `p` and `raf` will be change in the future. See the reference below for the definition of all the RAF.

### Value

Returns a list with the following components:

<code>call</code>	the <code>match.call()</code> .
<code>mu</code>	the estimate of the mean direction or the value supplied. If <code>num.sol &gt; 1</code> then <code>mu</code> may have length greater than 1, i.e. one value for each root found.
<code>kappa</code>	the estimate of the concentration parameter or the value supplied. If <code>num.sol &gt; 1</code> then <code>kappa</code> may have length greater than 1, i.e. one value for each root found.



tot.weights	the sum of the weights divide by the number of observations, one value for each root found.
weights	the weights associated to each observation, one column vector for each root found.
f.density	the non-parametric density estimation.
m.density	the smoothed model.
delta	the Pearson residuals.
tot.sol	the number of solutions found.
not.conv	the number of starting points that does not converge after the max.iter iteration are reached.

**Author(s)**

Claudio Agostinelli

**References**

C. Agostinelli. Robust estimation for circular data. *Computational Statistics & Data Analysis*, 51(12):5867-5875, 2007.

**See Also**

[circular](#), [mle.vonmises](#).

**Examples**

```
x <- c(rvonmises(n=50, mu=circular(0), kappa=10), rvonmises(n=5, mu=circular(pi/2), kappa=20))
wle.vonmises(x, smooth=20, group=5)
```

---

wle.weights

*Weights based on Weighted Likelihood for the normal model*


---

**Description**

This function evaluated the weights for the vector 'x' using the vector 'y' in the estimation of the density by the kernel density estimator.

**Usage**

```
wle.weights(x, y=NULL, smooth=0.0031, sigma2, raf=1, location=FALSE,
max.iter=1000, tol=10^(-6))
```

**Arguments**

x	the data set for which the weights would be calculate.
y	the data set used to calculate the weights.
smooth	the value of the smoothing parameter.
sigma2	an estimate of the variance.
raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
location	if TRUE the location is estimated. Only available when y=NULL.
max.iter	maximum number of iterations.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.

**Value**

weights	the weights associated to the x vector.
location	the location.
conv	TRUE if the convergence is achived.

**Author(s)**

Claudio Agostinelli

---

wle.wrappednormal      *Wrapped Normal Weighted Likelihood Estimates*

---

**Description**

Computes the weighted likelihood estimates for the parameters of a Wrapped Normal distribution: the mean direction and the concentration parameter (and the scale parameter).

**Usage**

```
wle.wrappednormal(x, mu, rho, sd, K, boot = 30, group, num.sol = 1, raf = "HD",
  smooth = 0.0031, tol = 10^(-6), equal = 10^(-3), min.sd = 0.001,
  min.k = 10, max.iter = 100, use.smooth = TRUE, alpha=NULL, p = 2,
  verbose = FALSE, control.circular=list())
```

```
## S3 method for class 'wle.wrappednormal'
print(x, digits = max(3, getOption("digits") - 3), ...)
```

**Arguments**

x	a vector. The object is coerced to class <code>circular</code> .
mu	if a values if provided the parameter is considered known.
rho	if a values if provided the parameter (and sd) is considered known.
sd	if a values if provided the parameter (and rho) is considered known.
K	number of elements used to approximate the density of the wrapped normal.
boot	the number of starting points based on bootstrap subsamples to use in the search of the roots.
group	the dimension of the bootstrap subsamples.
num.sol	maximum number of roots to be searched.
raf	type of Residual adjustment function to be use: raf="HD": Hellinger Distance RAF, raf="NED": Negative Exponential Disparity RAF, raf="SCHI2": Symmetric Chi-Squared Disparity RAF.
smooth	the value of the smoothing parameter.
tol	the absolute accuracy to be used to achieve convergence of the algorithm.
equal	the absolute value for which two roots are considered the same. (This parameter must be greater than tol).
min.sd	minimum value for the sd parameter.
min.k	minimum number of elements used to approximate the density of the wrapped normal.
max.iter	maximum number of iterations.
use.smooth	logical, if TRUE a smoothed model is used, default is TRUE.
alpha	if not NULL overrides the value of p. See the next argument p. This is a different parameterization, alpha=-1/2 provides Hellinger Distance RAF, alpha=-1 provides Kullback-Leibler RAF and alpha=-2 provides Neyman's Chi-Square RAF.
p	this parameter works only when raf="HD". p=2 provide Hellinger Distance RAF, p=-1 provide Kullback-Leibler RAF and p=Inf provide Neyman's Chi-Square RAF.
verbose	logical, if TRUE warnings are printed.
control.circular	the attribute of the resulting objects (mu)
digits	integer indicating the precision to be used.
...	further parameters in <code>print.wle.vonmises</code> .

**Details**

Parameters p and raf will be change in the future. See the reference below for the definition of all the RAF.

**Value**

Returns a list with the following components:

<code>call</code>	the <code>match.call()</code> .
<code>mu</code>	the estimate of the mean direction or the value supplied. If <code>num.sol &gt; 1</code> then <code>mu</code> may have length greater than 1, i.e, one value for each root found.
<code>rho</code>	the estimate of the concentration parameter or the value supplied. If <code>num.sol &gt; 1</code> then <code>rho</code> may have length greater than 1, i.e, one value for each root found.
<code>sd</code>	the estimate of the standard deviation parameter or the value supplied. If <code>num.sol &gt; 1</code> then <code>sd</code> may have length greater than 1, i.e, one value for each root found.
<code>tot.weights</code>	the sum of the weights divide by the number of observations, one value for each root found.
<code>weights</code>	the weights associated to each observation, one column vector for each root found.
<code>f.density</code>	the non-parametric density estimation.
<code>m.density</code>	the smoothed model.
<code>delta</code>	the Pearson residuals.
<code>tot.sol</code>	the number of solutions found.
<code>not.conv</code>	the number of starting points that does not converge after the <code>max.iter</code> iteration are reached.

**Author(s)**

Claudio Agostinelli

**References**

C. Agostinelli. Robust estimation for circular data. *Computational Statistics & Data Analysis*, 51(12):5867-5875, 2007.

**See Also**

[circular](#), [mle.wrappednormal](#).

**Examples**

```
x <- c(rwrappednormal(n=50, mu=circular(0), sd=1), rwrappednormal(n=5, mu=circular(pi/2), sd=0.5))
wle.wrappednormal(x, smooth=1/20, group=5)
```

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