

# Package ‘CARMAgeddon’

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**LinkingTo** Rcpp, RcppArmadillo (>= 0.6.000.0.0), RcppParallel

**Suggests** mvtnorm

**SystemRequirements** GNU make

## Description

Functions for fitting and working with multivariate overlapping noisy observation continuous-time autoregression (MONOCAR) models and other CARMA models.

**LazyLoad** true

**LazyData** true

**License** GPL-3

**URL** <http://monocar.tahk.us/>

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create.ctdata                      *Create and combine MONOCAR data sets*

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### Description

Create and combine data sets for estimation of MONOCAR models by `monocar.estimate` and `monocar.hist`

### Usage

```
create.ctdata(x, v, t1, t2 = NULL,
              data = NULL,
              series.name = NULL,
              house.name = NULL,
              process.name = NULL,
              transform = NULL,
              transform.var = NULL,
              time.origin = NULL,
              time.unit = NULL,
              exo.data = NULL,
              inclusive.end.date = FALSE)
```

```
join.ctdata(...)
```

### Arguments

<code>x</code>	A numeric vector giving the value of each observation.
<code>v</code>	A numeric vector giving the (untransformed) variance of each observation.
<code>t1</code>	A numeric, date-time, or date vector giving the start time for each observation.
<code>t2</code>	An optional numeric, date-time, or date vector giving the end time for each observation. Must be of the same class as <code>t1</code> . The default is to assume <code>t2</code> equals <code>t1</code> , meaning instantaneous observations.
<code>data</code>	An optional data frame in which to search for variables.
<code>series.name</code>	An optional character, factor, or integer vector identifying the series for each observation. The default value of <code>NULL</code> uses the quoted <code>x</code> argument.
<code>house.name</code>	An optional character, factor, or integer vector identifying the house for each observation. The default is to use the series.
<code>process.name</code>	An optional character, factor, or integer vector identifying the process for each observation. The default is to assume all observations belong to the same process.
<code>transform</code>	An optional function specifying a transformation to apply to each observation. The function must take two inputs corresponding to vectors of means and variances. The default is leave the observations untransformed. Alternatively, a character object giving the name of a common variance-stabilizing transformation. The two implemented transformations are "arcsin" for binomial-distributed data (which uses $\arcsin\left(\left(1 - \frac{1}{5000}\right)\left(2\frac{x}{v} - 1\right)\right)$ and implied variance $\frac{1}{v}$ ; this assumes <code>x</code> gives the number of successes and <code>v</code> gives the sample sizes for each observation) and "sqrt" for Poisson-distributed data (which uses $\sqrt{x + \frac{3}{8}}$ and implied variance $\frac{1}{4}$ ; <code>v</code> is ignored).

<code>transform.var</code>	An optional function specifying a transformation to apply to the variance of each observation. The function must take two inputs corresponding to vectors of means and variances. The default is leave the variances untransformed or, if <code>transform</code> is a character object, the corresponding transformation of the variances for the named variance-stabilizing transformation.
<code>time.origin</code>	An optional numeric, date-time, or date object of length one specifying origin time to be subtracted from <code>t1</code> and <code>t2</code> .
<code>time.unit</code>	An optional numeric (if <code>t1</code> is a numeric vector) or <code>difftime</code> (if <code>t1</code> is a date-time or date vector) object specifying the units for <code>t1</code> and <code>t2</code> (i.e., the amount of time that should be considered to be length one).
<code>exo.data</code>	A vector, data frame, matrix, or formula of possible exogenous variables to be included in the data.
<code>inclusive.end.date</code>	If <code>t1</code> and <code>t2</code> are date vectors, whether the end date means the start or end of the day. Thus, for an observation which starts and ends on January 1, a <code>FALSE</code> value for <code>inclusive.end.date</code> would imply a length of zero hours while a <code>TRUE</code> value would imply a length of 24 hours.
<code>...</code>	Multiple arguments giving objects of class <code>ct.data.frame</code> , usually each created by <code>create.ctdata</code> , to be joined together. Alternatively, a single list containing multiple objects of class <code>ct.data.frame</code> .

## Details

See [monocar.estimate](#) for more details about the MONOCAR model.

Observations are grouped into series, houses, and processes. Series correspond to different elements of the latent process, houses allow for house-specific effects, and processes allow for multiple, independent latent processes that share common parameters (specifically,  $\Theta$  and  $\Sigma$  as well as, optionally,  $\mu$  and/or  $\delta$ ; whether the latter are included is controlled by `mu.by.process` and `delta.by.process`). Series are the most fundamental of the three. Houses are usually nested within series or identical to series. In most cases, all observations belong to the same process.

If series are not specified, all observations will be assumed to belong to a single series. Usually this will not be the case. However, even when it is not the case, it may be easier to create separate `ct.data.frame` objects with `create.ctdata` for each series and combine them with `joint.ctdata`.

If houses are not specified, they will be assumed to equal the series (implying series are not further divided into houses). If processes are not specified, all observations will be assumed to be based on a single (usually multivariate) latent process, as is typically the case.

## Value

An object of class `ct.data.frame` containing the data.

## See Also

[monocar.estimate](#) and [monocar.hist](#).

## Examples

```
## Create a simple data set with instantaneous observations every
## day for the first month of 2000
obs <- c(-0.25, 1.91, 0.06, 2.24, -0.48, -1.84, -1, -0.07,
```

```

-0.5, -1.17, 0.32, 0.67, 0.99, 0.39, 1.2, 0.12, -0.7, -0.65,
-1.52, -2.28, 0.98, 1.62, -0.36, 0.94, -1.61, -0.25, -0.76,
-1.81, -1.09, 0.9, -0.79)
obs.time <- seq(as.Date("2000-1-1"), by=1, length.out=length(obs))
obs.var <- rep(1, length(obs))
new.data <- create.ctdata(obs, obs.var, obs.time)

```

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monocar.estimate      *Estimate MONOCAR model*

---

## Description

Estimate a Multivariate overlapping noisy observation continuous-time autoregressive (MONOCAR) model

## Usage

```

monocar.estimate(data, init=list(), restrict=list(),
  subset=NULL, exovars=NULL, byexo="series",
  enforce.bounds=TRUE, lower.bounds=NULL, upper.bounds=NULL,
  delta.group=TRUE, verbose=1, mu.by.process=FALSE,
  delta.by.process=FALSE, var.by.house=FALSE, remove.unused=TRUE,
  ...)

```

## Arguments

data	A data file of class <code>ct.data.frame</code> (see <a href="#">create.ctdata</a> )
init	A list containing initial values for parameters. Omitted items will be guessed. The default is for <code>theta</code> , <code>sigma</code> , and <code>alpha</code> to be identity matrices, one of <code>mu</code> or <code>delta</code> to be the mean of each series or house (depending on whether <code>delta.group</code> is used), and variance-transformation parameters to be 0 for <code>var.add</code> and 1 for <code>var.mult</code> and <code>var.pow</code> (i.e., no transformation).
restrict	A list of which parameters to be restricted to their initial values. TRUE means restricted to initial value, FALSE means unrestricted, NA means restricted to equal the previous parameter (useful, for example, to set the variance-transformation parameters to be equal for multiple series. The parameter <code>sigma</code> can also be specified as one of the following strings: "diagonal" (meaning only the off-diagonal elements of <code>sigma</code> are restricted), "restricted" (meaning all elements of <code>sigma</code> are restricted), or "unrestricted" (meaning all elements of <code>sigma</code> are unrestricted). The default is for <code>theta</code> to be unrestricted, <code>sigma</code> to be restricted to be diagonal, <code>alpha</code> to be fully restricted, one of <code>mu</code> or <code>delta</code> to be unrestricted (depending on <code>delta.group</code> or whether initial values are explicitly specified), and variance-transformation parameters to be restricted unless initial values are explicitly specified.
subset	An optional vector specifying a subset of observations to be used in the fitting process.
exovars	An optional formula giving the exogenous variables to be included, a character vector giving the names of exogenous variables to be included from <code>data</code> , or an integer vector giving the location of the exogenous variables to be included from <code>data</code> . The default is to include no exogenous variables.

byexo	A string or null value specifying whether exogenous variables are to have a single coefficient, separate coefficients for each series, or separate coefficients for each house. Possible values are <code>series</code> , <code>house</code> or <code>NULL</code> . Default is <code>NULL</code> , meaning a single coefficient, common to all series and houses, for each exogenous variable.
enforce.bounds	A boolean or integer specifying whether to enforce bounds on the variance-transformation parameters during optimization. If equal to <code>FALSE</code> or <code>0</code> , no bounds are imposed. If equal to <code>TRUE</code> or <code>1</code> , <code>var.add</code> and <code>var.mult</code> are restricted to be non-negative, avoiding negative transformed variances. If equal to <code>2</code> , <code>var.pow</code> is also restricted to be non-negative, ensuring the variance transformation is non-decreasing. The default is <code>TRUE</code> .
lower.bounds	An optional vector of lower bounds for parameters if desired. The default is <code>NULL</code> , meaning no lower bounds other than those implied by <code>enforce.bounds</code> .
upper.bounds	An optional vector of lower bounds for parameters if desired. The default is <code>NULL</code> , meaning no lower bounds.
delta.group	An optional vector of integers specifying which houses are associated with which series. Alternatively, the logical value <code>TRUE</code> , in which case the data will be used to determine which series corresponds to each house, or <code>NULL</code> , in which case no attempt to recenter $\delta$ will be taken. <code>delta.group</code> is used to recenter $\delta$ such that the mean of <i>delta</i> within each series is zero with $\mu$ adjusted to compensate (thus providing an equivalent parameterization). This only makes sense if houses are nested within series. The value <code>TRUE</code> will be treated as <code>NULL</code> if houses are not nested within series. The default is <code>TRUE</code> .
verbose	An integer specifying the amount of output. <code>0</code> means no output, <code>1</code> means a dot ( <code>.</code> ) for each gradient computation and a star ( <code>*</code> ) for each Hessian evaluation, and a <code>2</code> means a dot for each likelihood evaluation, a colon ( <code>:</code> ) for each gradient evaluation, and a star for each Hessian evaluation. A <code>+</code> indicates an optimization algorithm converged while <code>-</code> means it failed to do so. The default is <code>1</code> .
mu.by.process	If multiple processes are specified, a boolean indicating whether the series means should be identical across processes. The default is <code>FALSE</code> .
delta.by.process	If multiple processes are specified, a boolean indicating whether the house means should be identical across processes. The default is <code>FALSE</code> .
var.by.house	Whether variance-transformation parameters should be applied by house or by series. The default is to apply by series.
remove.unused	Whether to remove unused series and houses, thereby reducing the number of parameters. This can be particularly useful when using a subset of the data that does not contain all houses. Note that the size of the parameters specified in <code>init</code> and <code>restrict</code> may change as a result.
...	Additional parameters to be passed to the <code>monocar.cal</code> , which carries out the optimization.

## Details

Observations are grouped into series, houses, and processes. Series correspond to different elements of the latent process, houses allow for house-specific effects, and processes allow for multiple, independent latent processes that share common parameters (specifically,  $\Theta$  and  $\Sigma$  as well as, optionally,  $\mu$  and/or  $\delta$ ; whether the latter are included is controlled by `mu.by.process` and `delta.by.process`). Series are the most fundamental of the three. Houses are usually nested within series or identical to series. In most cases, all observations belong to a single process.

The latent process,  $x_t$  is an Ornstein-Uhlenbeck process defined by the stochastic differential equation

$$x_t = \Theta(\mu - x_t) dt + \Sigma^{\frac{1}{2}} dW_t$$

where  $\Theta$  describes how each element of the latent process affects itself and other elements of the latent process,  $\Sigma$  is the instantaneous variance-covariance matrix for the error process,  $\mu$  is the long-run mean of  $\mu$ , and  $W_t$  is a multivariate Wiener process.

The observation equations are defined by

$$y_i = \int_{r_i}^{s_i} x_{s_i,t} dt + \delta_{h_i} + \epsilon_i$$

when  $q_i < r_i$  or

$$y_i = x_{s_i,q_i} + \delta_{h_i} + \epsilon_i$$

when  $q_i = r_i$  with

$$\epsilon_i \sim \mathcal{N}\left(0, \gamma_{s_i} + \nu_{s_i} \left(\frac{v_i}{\bar{v}_{s_i}}\right)^{\xi_{s_i}}\right)$$

and independent across observations. Here,  $y_i$  is the value of observation  $i$ ,  $v_i$  is the untransformed (or naive) variance of the error term for the observation,  $\delta_j$  is a house-specific effect for house  $j$ ,  $h_i$  is the house of equation  $i$ ,  $q_i$  is the start time for observation  $i$ ,  $r_i$  is the end time for observation  $i$ ,  $\bar{v}_j$  is the variance center for house  $j$ , and  $s_i$  and  $h_i$  index the series and house of observation  $i$ .  $y_i$ ,  $v_i$ ,  $q_i$ ,  $r_i$ ,  $s_i$ , and  $h_i$  are all known (entered as `x`, `v`, `t1`, and `t2` in `create.ctdata`). If `var.by.house` is TRUE, then  $\gamma$ ,  $\nu$ ,  $\xi$ , and  $\bar{v}$  are subscripted by  $h_i$  rather than  $s_i$ .

The variance centers ( $\bar{v}$ ) are used to improve parameter estimation and interpretation. By default, they are the mean of the untransformed variances for observations in each house. These default to the mean variance of each series or house, but other values can be specified (see `monocar.cal`). Aside from possible computational issues, a change of variance centers does not affect the model once  $\nu$  and  $\xi$  are adjusted appropriately.

For multiple processes, the latent process is replaced by multiple latent processes with common parameters  $\Theta$  and  $\Sigma$  defined by

$$x_{k,t} = \Theta(\mu_k - x_{k,t}) dt + \Sigma^{\frac{1}{2}} dW_{k,t}$$

for all  $k$ , where  $k$  indexes the latent process. The observation equations also likewise adjusted by replacing  $x_t$  by  $x_{k_i,t}$ , where  $k_i$  is the latent process observed by observation  $i$ .

For a structural MONOCAR model, the latent process becomes

$$Ax_t = \Theta(\mu - x_t) dt + \Sigma^{\frac{1}{2}} dW_t$$

where  $A$  is a matrix describing instantaneous, non-gradual effects and is assumed to be the identity matrix by default.

In specifying `init` and `restrict`, the possible parameters to list are `theta` ( $\Theta$ , an  $N \times N$  matrix), `sigma` ( $\Sigma$ , an  $n \times n$  matrix), `mu` ( $\mu$ , an  $N$ -element vector), `delta` ( $\delta$ , an  $H$ -element vector), `alpha` ( $A$ , an  $N \times N$  matrix), `var.add` ( $\gamma$ , an  $H$ -element vector), `var.mult` ( $\nu$ , an  $H$ -element vector), and `var.pow` ( $\xi$ , an  $H$ -element vector), where  $N$  is the number of series and  $H$  is the number of houses. Parameters can always be omitted from `init` and/or `restrict`, in which case reasonable guesses are taken. If `mu.by.process` is TRUE, then `mu` is a  $K \times N$  matrix, where  $K$  is the number of processes. If `delta.by.process` is TRUE, then `delta` should be a  $K \times H$  matrix.

The actual estimation is carried out by the internal function `monocar.cal`. See `monocar.cal` for additional arguments which control the estimation process. The arguments `optimizers`, `maxiters`, `tolerances`, `epsilon`, `lower.bounds`, `upper.bounds`, and `var.centers` can be included as arguments to `monocar.estimate`, which passes them to `monocar.cal`.

**Value**

An object of class `monocar`.

**References**

Tahk, A. M. (2015) “A Continuous-Time, Latent-Variable Model of Time-Series Data,” *Political Analysis* 23 (2): 278–298.

**See Also**

[summary.monocar](#), [monocar.estimate](#), and [create.ctdata](#).

**Examples**

```
## Run a simple MONOCAR model
car.model <- monocar.estimate(data = PresUnemploy)
summary(car.model)

## Run a MONOCAR model with a Bush administration
## dummy variable added
car.model.bush <-
  monocar.estimate(data = PresUnemploy, exo="Bush")
summary(car.model.bush)

## Run a MONOCAR model on Gallup polls only
car.model.gallup <-
  monocar.estimate(data = PresUnemploy,
                  subset = house == "Gallup" |
                  house == "unemployment")
summary(car.model.gallup)

## Run a MONOCAR model with correlated error processes
car.model.cor <-
  monocar.estimate(data = PresUnemploy,
                  restrict=list(sigma="unrestricted"))
summary(car.model.cor)

## Run a MONOCAR model with correlated error processes and transformed variances
car.model.cor.var <-
  monocar.estimate(data = PresUnemploy,
                  restrict=list(sigma="unrestricted",
                                var.add=c(TRUE,FALSE),
                                var.mult=c(FALSE,TRUE),
                                var.pow=c(FALSE,TRUE)))
summary(car.model.cor.var)
```

---

`monocar.hist`

*Estimate a series history from a MONOCAR model*

---

**Description**

Estimate the history of a latent process from a Multivariate overlapping noisy observation continuous-time autoregressive (MONOCAR) model

**Usage**

```
monocar.hist(data, model, estimates=NULL, exovars=NULL, byexo="series",
             delta.group=TRUE, verbose=1, mu.by.process=FALSE,
             delta.by.process=FALSE, var.by.house=FALSE, ...)
```

**Arguments**

<code>data</code>	A data file of class <code>ct.data.frame</code> (see <a href="#">create.ctdata</a> )
<code>model</code>	A model of class <code>monocar</code> generated by <a href="#">monocar.estimate</a> . Can be <code>NULL</code> , in which case the specification is taken from <code>estimates</code> and <code>exovars</code> .
<code>estimates</code>	If <code>model</code> is <code>NULL</code> , a list of parameters.
<code>exovars</code>	If <code>model</code> is <code>NULL</code> , an optional formula giving the exogenous variables to be included, a character vector giving the names of exogenous variables to be included from <code>data</code> , or an integer vector giving the location of the exogenous variables to be included from <code>data</code> . The default is to include no exogenous variables.
<code>byexo</code>	If <code>model</code> is <code>NULL</code> , a string or null value specifying whether exogenous variables are to have a single coefficient, separate coefficients for each series, or separate coefficients for each house. Possible values are <code>series</code> , <code>house</code> or <code>NULL</code> . Default is <code>NULL</code> , meaning a single coefficient, common to all series and houses, for each exogenous variable.
<code>delta.group</code>	An optional vector of integers specifying which houses are associated with which series.
<code>verbose</code>	An integer specifying the amount of output. 0 means no output, 1 means a dot (.) for each gradient computation and a star (*) for each hessian evaluation, and a 2 means a dot for each likelihood evaluation, a colon (:) for each gradient evaluation, and a star for each hessian evaluation. The default is 1.
<code>mu.by.process</code>	If multiple processes are specified, a boolean indicating whether the series means should be identical across processes. The default is <code>FALSE</code> .
<code>delta.by.process</code>	If multiple processes are specified, a boolean indicating whether the house means should be identical across processes. The default is <code>FALSE</code> .
<code>var.by.house</code>	Whether variance-transformation parameters should be applied by house or by series. The default is to apply by series.
<code>...</code>	Additional parameters to be passed to the <code>monocar.cal</code> , which carries out the estimation.

**References**

Tahk, A. M. (2015) "A Continuous-Time, Latent-Variable Model of Time-Series Data," *Political Analysis* 23 (2): 278–298.

**See Also**

[monocar.estimate](#) and [create.ctdata](#).

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PresUnemploy	<i>Presidential Approval and Unemployment in the Reagan and Bush administrations</i>
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**Description**

This data set gives the monthly unemployment rate and polls of presidential approval during the Reagan and Bush administrations.

**Usage**

```
PresUnemploy
```

**Format**

A `ct.data.frame` object containing 677 observations.

**Source**

Roper Center's iPoll DataBank and Federal Reserve Economic Data (FRED)

**References**

Tahk, A. M. (2015) "A Continuous-Time, Latent-Variable Model of Time-Series Data," *Political Analysis* 23 (2): 278–298.

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<code>simulate.monocar</code>	<i>Simulate continuous-time data sets</i>
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**Description**

Generate simulated continuous-time data sets

**Usage**

```
## S3 method for class 'monocar'
simulate(object, nsim=1, seed=NULL,
  var=NULL, t1=seq(0,100,1.0), t2=NULL, data=NULL,
  series.name = NULL, house.name = NULL, process.name = NULL,
  transform = function(x, v) x, transform.var = function(x, v) v,
  exo.data=NULL, exovars=NULL, byexo=NULL, var.centers=NULL, ...)
```

**Arguments**

<code>object</code>	A an object of class <code>monocar</code> or a list specifying the parameters of the model.
<code>nsim</code>	Number of data sets to simulate. Defaults to 1.
<code>seed</code>	An object specifying if and how the random number generator should be initialized. Either <code>NULL</code> or an integer that will be used in a call to <code>set.seed</code> before simulating the response vectors. If set, the value is saved as the <code>seed</code> attribute of the returned value. The default, <code>NULL</code> will not change the random generator state, and return <code>.Random.seed</code> as the <code>seed</code> attribute.

var	A numeric vector giving the variance of the measurement error for each observation. The default value of NULL means zero measurement error for each observation.
t1	A numeric vector giving the start time for each observation.
t2	An optional numeric vector giving the end time for each observation. By default, the end time of each observation equals the start time.
data	An optional data frame or object of class <code>ct.data.frame</code> .
series.name	An optional character, factor, or integer vector identifying the series for each observation. The default value of NULL uses the quoted <code>x</code> argument.
house.name	An optional character, factor, or integer vector identifying the house for each observation. The default is to use the series.
process.name	An optional character, factor, or integer vector identifying the process for each observation. The default is to assume all observations belong to the same process.
transform	An optional function specifying a transformation to apply to each observation. See <a href="#">monocar.data</a> for more details.
transform.var	An optional function specifying a transformation to apply to the variance of each observation. See <a href="#">monocar.data</a> for more details.
exo.data	A vector, data frame, matrix, or formula of possible exogenous variables to be included in the data.
exovars	An optional formula giving the exogenous variables to be included, a character vector giving the names of exogenous variables to be included from <code>data</code> , or an integer vector giving the location of the exogenous variables to be included from <code>data</code> . The default is to include no exogenous variables.
byexo	A string or null value specifying whether exogenous variables are to have a single coefficient, separate coefficients for each series, or separate coefficients for each house. Possible values are <code>series</code> , <code>house</code> or <code>NULL</code> . Default is <code>NULL</code> , meaning a single coefficient, common to all series and houses, for each exogenous variable.
var.centers	For each house, the value around which to recenter the variances. See <a href="#">monocar.cal</a> for more details.
...	Additional optional arguments.

### Value

If `nsim==1`, an object of class `ct.data.frame` containing the data. Otherwise, a list containing `nsim` objects of class `ct.data.frame`, each containing one simulated data set.

### See Also

[create.ctdata](#), [monocar.estimate](#), and [monocar.hist](#).

---

summary.monocar      *Print and summarize MONOCAR models*

---

## Description

Print and summary methods for MONOCAR models from `monocar.estimate`.

## Usage

```
## S3 method for class 'monocar'
print(x, reverse.offdiag = TRUE,
      digits = max(3L, getOption("digits") - 3L), ...)

## S3 method for class 'monocar'
summary(object, reverse.offdiag = TRUE, ...)

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

## Arguments

<code>x</code>	For <code>print.monocar</code> , an object of class <code>monocar</code> . For <code>print.summary.monocar</code> , an object of class <code>summary.monocar</code> .
<code>object</code>	An object of class <code>monocar</code> .
<code>reverse.offdiag</code>	A logical value indicating whether to reverse the sign of off-diagonal elements of Theta. In a MONOCAR model, a positive value of an off-diagonal element of Theta indicates that an increase in one latent process tends to <i>decrease</i> the level of another latent process. Reversing the sign of these parameters means that a positive value indicates one process tends to increase another. This can simplify the interpretation of the parameters. The default is to reverse the sign of off-diagonal elements of Theta. For <code>print.summary.monocar</code> , whether these elements are reversed is set by <code>summary.monocar</code> and stored in the <code>summary.monocar</code> object.
<code>digits</code>	The minimal number of significant digits, see <code>print.default</code> .
<code>signif.stars</code>	A logical value indicating whether stars should be printed based on p-values, see <code>printCoefmat</code> .
<code>...</code>	Additional arguments to pass to <code>print.default</code> or <code>printCoefmat</code> .

## Value

`summary.monocar` returns an object of class `summary.monocar` containing a summary information about the model.

## See Also

[monocar.estimate](#) and [monocar.hist](#).

**Examples**

```
## Run a simple MONOCAR model
car.model <- monocar.estimate(data = PresUnemploy)
## Print model estimates
car.model
## Print model estimates with diagonal not reversed
print(car.model, reverse.offdiag=FALSE)
## Print summary information about the model
summary(car.model)
```

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