

# Package ‘hmmTMB’

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**Type** Package

**Title** Fit Hidden Markov Models using Template Model Builder

**Version** 1.0.1

**Maintainer** Theo Michelot <theo.michelot@dal.ca>

**Description** Fitting hidden Markov models using automatic differentiation and Laplace approximation, allowing for fast inference and flexible covariate effects (including random effects and smoothing splines) on model parameters. The package is described by Michelot (2022) <[arXiv:2211.14139](#)>.

**URL** <https://github.com/TheoMichelot/hmmTMB>

**License** GPL-3

**Depends** R6, mgcv, TMB, ggplot2

**Imports** Matrix, stringr, optimx, CircStats, MASS, tmbstan, methods

**LinkingTo** TMB, RcppEigen

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**Author** Theo Michelot [aut, cre],  
Richard Glennie [aut, ctb]

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---

as\_character\_formula *Read formula with as.character without splitting*

---

## Description

Read formula with as.character without splitting

## Usage

```
as_character_formula(x, ...)
```

## Arguments

x	R formula
...	Unused

## Details

Citation: this function was taken from the R package formula.tools: Christopher Brown (2018). formula.tools: Programmatic Utilities for Manipulating Formulas, Expressions, Calls, Assignments and Other R Objects. R package version 1.7.1. <https://CRAN.R-project.org/package=formula.tools>

---

as_sparse	<i>Transforms matrix to dgTMatrix</i>
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**Description**

Transforms matrix to dgTMatrix

**Usage**

```
as_sparse(x)
```

**Arguments**

x                    Matrix or vector. If this is a vector, it is formatted into a single-column matrix.

**Value**

Sparse matrix of class dgTMatrix

---

bdiag_check	<i>Create block diagonal matrix (safe version)</i>
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**Description**

This version of bdiag checks whether the matrices passed as arguments are NULL. This avoids errors that would arise if using bdiag directly.

**Usage**

```
bdiag_check(...)
```

**Arguments**

...                    Matrix or list of matrices (only the first argument is used)

**Value**

Block diagonal matrix

---

cov_grid	<i>Grid of covariates</i>
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**Description**

Grid of covariates

**Usage**

```
cov_grid(var, data = NULL, obj = NULL, covs = NULL, formulas, n_grid = 1000)
```

**Arguments**

var	Name of variable
data	Data frame containing the covariates. If not provided, data are extracted from obj
obj	HMM model object containing data and formulas
covs	Optional named list for values of covariates (other than 'var') that should be used in the plot (or dataframe with single row). If this is not specified, the mean value is used for numeric variables, and the first level for factor variables.
formulas	List of formulas used in the model
n_grid	Grid size (number of points). Default: 1000.

**Value**

Data frame of covariates, with 'var' defined over a grid, and other covariates fixed to their mean (numeric) or first level (factor).

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Dist	<i>R6 class for probability distribution</i>
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**Description**

R6 class for probability distribution

R6 class for probability distribution

**Details**

Contains the probability density/mass function, and the link and inverse link functions for a probability distribution.

**Methods****Public methods:**

- `Dist$new()`
- `Dist$name()`
- `Dist$pdf()`
- `Dist$rng()`
- `Dist$link()`
- `Dist$invlink()`
- `Dist$npar()`
- `Dist$parnames()`
- `Dist$parapprox()`
- `Dist$fixed()`
- `Dist$code()`
- `Dist$name_long()`
- `Dist$set_npar()`
- `Dist$set_parnames()`
- `Dist$set_code()`
- `Dist$pdf_apply()`
- `Dist$rng_apply()`
- `Dist$n2w()`
- `Dist$w2n()`
- `Dist$clone()`

**Method** `new()`: Create a Dist object

*Usage:*

```
Dist$new(  
  name,  
  pdf,  
  rng,  
  link,  
  invlink,  
  npar,  
  parnames,  
  parapprox = NULL,  
  fixed = NULL,  
  name_long = name  
)
```

*Arguments:*

`name` Name of distribution

`pdf` Probability density/mass function of the distribution (e.g. `dnorm` for normal distribution).

`rng` Random generator function of the distribution (e.g. `rnorm` for normal distribution).

`link` Named list of link functions for distribution parameters

`invlink` Named list of inverse link functions for distribution parameters

`npar` Number of parameters of the distribution  
`parnames` Character vector with name of each parameter  
`parapprox` Function that takes a sample and produces approximate values for the unknown parameters  
`fixed` Vector with element for each parameter which is TRUE if parameter is fixed  
`name_long` Long version of the name of the distribution, possibly more user-readable than name.

*Returns:* A new Dist object

**Method** `name()`: Return name of Dist object

*Usage:*

`Dist$name()`

**Method** `pdf()`: Return pdf of Dist object

*Usage:*

`Dist$pdf()`

**Method** `rng()`: Return random generator function of Dist object

*Usage:*

`Dist$rng()`

**Method** `link()`: Return link function of Dist object

*Usage:*

`Dist$link()`

**Method** `invlink()`: Return inverse link function of Dist object

*Usage:*

`Dist$invlink()`

**Method** `npar()`: Return number of parameters of Dist object

*Usage:*

`Dist$npar()`

**Method** `parnames()`: Return names of parameters

*Usage:*

`Dist$parnames()`

**Method** `parapprox()`: Return function that approximates parameters

*Usage:*

`Dist$parapprox()`

**Method** `fixed()`: Return which parameters are fixed

*Usage:*

`Dist$fixed()`

**Method** `code()`: Return code of Dist object

*Usage:*

`Dist$code()`

**Method** `name_long()`: Human-readable name of Dist object

*Usage:*

`Dist$name_long()`

**Method** `set_npar()`: Set number of parameters this distribution has

*Usage:*

`Dist$set_npar(new_npar)`

*Arguments:*

`new_npar` Number of parameters

**Method** `set_parnames()`: Set parameter names

*Usage:*

`Dist$set_parnames(new_parnames)`

*Arguments:*

`new_parnames` Parameter names

**Method** `set_code()`: Set distribution code

*Usage:*

`Dist$set_code(new_code)`

*Arguments:*

`new_code` Distribution code

**Method** `pdf_apply()`: Evaluate probability density/mass function

This method is used in the `Dist$obs_probs()` method. It is a wrapper around `Dist$pdf()`, which prepares the parameters and passes them to the function.

*Usage:*

`Dist$pdf_apply(x, par, log = FALSE)`

*Arguments:*

`x` Value at which the function should be evaluated

`par` Vector of parameters. The entries should be named if they are not in the same order as expected by the R function. (E.g. `shape/scale` rather than `shape/rate` for gamma distribution.)

`log` Logical. If TRUE, the log-density is returned. Default: FALSE.

*Returns:* Probability density/mass function evaluated at `x` for parameters `par`

**Method** `rng_apply()`: Random number generator

This method is a wrapper around `Dist$rng()`, which prepares the parameters and passes them to the function.

*Usage:*

`Dist$rng_apply(n, par)`

*Arguments:*

n Number of realisations to generate

par Vector of parameters. The entries should be named if they are not in the same order as expected by the R function. (E.g. shape/scale rather than shape/rate for gamma distribution.)

*Returns:* Vector of n realisations of this distribution

**Method** n2w(): Natural to working parameter transformation

This method transforms parameters from the natural scale (i.e., their domain of definition) to the "working" or "linear predictor" scale (i.e., the real line). It is a wrapper for Dist\$link().

*Usage:*

```
Dist$n2w(par)
```

*Arguments:*

par List of parameters

*Returns:* Vector of parameters on the working scale

**Method** w2n(): Working to natural parameter transformation

This method transforms parameters from the "working" or "linear predictor" scale (i.e., the real line) to the natural scale (i.e., their domain of definition). It is a wrapper for Dist\$invlink().

*Usage:*

```
Dist$w2n(wpar, as_matrix = FALSE)
```

*Arguments:*

wpar Vector of working parameters

as\_matrix Logical. If TRUE, the natural parameters are returned as a matrix with one row for each state and one column for each parameter. If FALSE, the natural parameters are returned as a list (default).

*Returns:* List or matrix of parameters on natural scale

**Method** clone(): The objects of this class are cloneable with this method.*Usage:*

```
Dist$clone(deep = FALSE)
```

*Arguments:*

deep Whether to make a deep clone.

---

 find\_re

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*Find s(, bs = "re") terms in formula*


---

**Description**

This function is used to identify the variables "x" which are included as s(x, bs = "re") in the formula, in particular to check that they are factors.



**Usage**

```
find_re(form)
```

**Arguments**

form            Model formula

**Value**

Vector of names of variables for which a random effect term is included in the model.

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HMM

*R6 class for hidden Markov model*

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**Description**

R6 class for hidden Markov model

R6 class for hidden Markov model

**Details**

Encapsulates the observation and hidden state models for a hidden Markov model.

**Methods****Public methods:**

- [HMM\\$new\(\)](#)
- [HMM\\$obs\(\)](#)
- [HMM\\$hid\(\)](#)
- [HMM\\$out\(\)](#)
- [HMM\\$tm\\_b\\_obj\(\)](#)
- [HMM\\$tm\\_b\\_obj\\_joint\(\)](#)
- [HMM\\$tm\\_b\\_rep\(\)](#)
- [HMM\\$states\(\)](#)
- [HMM\\$coeff\\_fe\(\)](#)
- [HMM\\$coeff\\_re\(\)](#)
- [HMM\\$coeff\\_list\(\)](#)
- [HMM\\$fixpar\(\)](#)
- [HMM\\$coeff\\_array\(\)](#)
- [HMM\\$lambda\(\)](#)
- [HMM\\$update\\_par\(\)](#)
- [HMM\\$sd\\_re\(\)](#)
- [HMM\\$par\(\)](#)
- [HMM\\$set\\_priors\(\)](#)

- `HMM$priors()`
- `HMM$iters()`
- `HMM$out_stan()`
- `HMM$llk()`
- `HMM$edf()`
- `HMM$setup()`
- `HMM$fit_stan()`
- `HMM$fit()`
- `HMM$mle()`
- `HMM$forward_backward()`
- `HMM$cond()`
- `HMM$pseudores()`
- `HMM$viterbi()`
- `HMM$sample_states()`
- `HMM$state_probs()`
- `HMM$post_coeff()`
- `HMM$post_linpred()`
- `HMM$post_fn()`
- `HMM$predict()`
- `HMM$confint()`
- `HMM$simulate()`
- `HMM$check()`
- `HMM$plot_ts()`
- `HMM$plot_dist()`
- `HMM$plot()`
- `HMM$AIC_marginal()`
- `HMM$AIC_conditional()`
- `HMM$print_obspar()`
- `HMM$print_tpm()`
- `HMM$formulation()`
- `HMM$print()`
- `HMM$clone()`

**Method** `new()`: Create new HMM object

*Usage:*

```
HMM$new(obs = NULL, hid = NULL, file = NULL, init = NULL, fixpar = NULL)
```

*Arguments:*

`obs` Observation object, created with `Observation$new()`. This contains the formulation for the observation model.

`hid` MarkovChain object, created with `MarkovChain$new()`. This contains the formulation for the state process model.

`file` Path to specification file for HMM. If this argument is used, then `obs` and `hid` are unnecessary.

`init` HMM object, used to initialise the parameters for this model. If `init` is passed, then all parameters that are included in `init` and in the present model are copied. This may be useful when fitting increasingly complex models: start from a simple model, then pass it as `init` to create a more complex model, and so on.

`fixpar` Named list, with optional elements: 'hid', 'obs', 'delta0', 'lambda\_obs', and 'lambda\_hid'. Each element is a named vector of parameters in `coeff_fe` that should either be fixed (if the corresponding element is set to NA) or estimated to a common value (using integers or factor levels). See examples in the vignettes, and check the TMB documentation to understand the inner workings (argument `map` of `TMB::MakeADFun()`).

*Returns:* A new HMM object

*Examples:*

```
# Load data set (included with R)
data(nottem)
data <- data.frame(temp = as.vector(t(nottem)))

# Create hidden state and observation models
hid <- MarkovChain$new(data = data, n_states = 2)
par0 <- list(temp = list(mean = c(40, 60), sd = c(5, 5)))
obs <- Observation$new(data = data, n_states = 2,
                      dists = list(temp = "norm"),
                      par = par0)

# Create HMM
hmm <- HMM$new(hid = hid, obs = obs)
```

**Method** `obs()`: Observation object for this model

*Usage:*

```
HMM$obs()
```

**Method** `hid()`: MarkovChain object for this model

*Usage:*

```
HMM$hid()
```

**Method** `out()`: Output of optimiser after model fitting

*Usage:*

```
HMM$out()
```

**Method** `tmb_obj()`: Model object created by TMB. This is the output of the TMB function `MakeADFun`, and it is a list including elements

- `fn` Objective function
- `gr` Gradient function of `fn`
- `par` Vector of initial parameters on working scale

*Usage:*

```
HMM$tmb_obj()
```

**Method** `tmb_obj_joint()`: Model object created by TMB for the joint likelihood of the fixed and random effects. This is the output of the TMB function `MakeADFun`, and it is a list including elements

- `fn` Objective function
- `gr` Gradient function of `fn`
- `par` Vector of initial parameters on working scale

*Usage:*

`HMM$tmb_obj_joint()`

**Method** `tmb_rep()`: Output of the TMB function `sdreport`, which includes estimates and standard errors for all model parameters.

*Usage:*

`HMM$tmb_rep()`

**Method** `states()`: Vector of estimated states, after `viterbi` has been run

*Usage:*

`HMM$states()`

**Method** `coeff_fe()`: Coefficients for fixed effect parameters

*Usage:*

`HMM$coeff_fe()`

**Method** `coeff_re()`: Coefficients for random effect parameters

*Usage:*

`HMM$coeff_re()`

**Method** `coeff_list()`: List of all model coefficients

These are the parameters estimated by the model, including fixed and random effect parameters for the observation parameters and the transition probabilities, (transformed) initial probabilities, and smoothness parameters.

*Usage:*

`HMM$coeff_list()`

**Method** `fixpar()`: Fixed parameters

*Usage:*

`HMM$fixpar(all = FALSE)`

*Arguments:*

`all` Logical. If `FALSE`, only user-specified fixed parameters are returned, but not parameters that are fixed by definition (e.g., size of binomial distribution).

**Method** `coeff_array()`: Array of working parameters

*Usage:*

`HMM$coeff_array()`

**Method** `lambda()`: Smoothness parameters

*Usage:*

HMM\$lambda()

**Method** update\_par(): Update parameters stored inside model object

*Usage:*

HMM\$update\_par(par\_list = NULL, iter = NULL)

*Arguments:*

par\_list List with elements for coeff\_fe\_obs, coeff\_fe\_hid, coeff\_re\_obs, coeff\_re\_hid, log\_delta0, log\_lambda\_hid, and log\_lambda\_obs

iter Optional argument to update model parameters based on MCMC iterations (if using rstan). Either the index of the iteration to use, or "mean" if the posterior mean should be used.

**Method** sd\_re(): Standard deviation of smooth terms (or random effects)

This function transforms the smoothness parameter of each smooth term into a standard deviation, given by  $SD = 1/\sqrt{\lambda}$ . It is particularly helpful to get the standard deviations of independent normal random effects.

*Usage:*

HMM\$sd\_re()

*Returns:* List of standard deviations for observation model and hidden state model.

**Method** par(): Model parameters

*Usage:*

HMM\$par(t = 1)

*Arguments:*

t returns parameters at time t, default is t = 1

*Returns:* A list with elements:

- obsparParameters of observation model
- tpmTransition probability matrix of hidden state model

**Method** set\_priors(): Set priors for coefficients

*Usage:*

HMM\$set\_priors(new\_priors = NULL)

*Arguments:*

new\_priors is a list of matrices for optionally coeff\_fe\_obs, coeff\_fe\_hid, log\_lambda\_obs, log\_lambda\_hid each matrix has two rows (first row = mean, second row = sd) specifying parameters for Normal priors

**Method** priors(): Extract stored priors

*Usage:*

HMM\$priors()

**Method** iters(): Iterations from stan MCMC fit

*Usage:*

```
HMM$iters(type = "response")
```

*Arguments:*

*type* Either "response" for parameters on the response (natural) scale, or "raw" for parameters on the linear predictor scale.

*Returns:* see output of `as.matrix` in `stan`

**Method** `out_stan()`: fitted stan object from MCMC fit

*Usage:*

```
HMM$out_stan()
```

*Returns:* the stanfit object

**Method** `llk()`: Log-likelihood at current parameters

*Usage:*

```
HMM$llk()
```

*Returns:* Log-likelihood

**Method** `edf()`: Effective degrees of freedom

*Usage:*

```
HMM$edf()
```

*Returns:* Number of effective degrees of freedom (accounting for flexibility in non-parametric terms implied by smoothing)

**Method** `setup()`: TMB setup

This creates an attribute `tmb_obj`, which can be used to evaluate the negative log-likelihood function.

*Usage:*

```
HMM$setup(silent = TRUE)
```

*Arguments:*

*silent* Logical. If TRUE, all tracing outputs are hidden (default).

**Method** `fit_stan()`: Fit model using `tmbstan`

Consult documentation of the `tmbstan` package for more information. After this method has been called, the Stan output can be accessed using the method `out_stan()`. This Stan output can for example be visualised using functions from the `rstan` package. The parameters stored in this HMM object are automatically updated to the mean posterior estimate, although this can be changed using `update_par()`.

*Usage:*

```
HMM$fit_stan(..., silent = FALSE)
```

*Arguments:*

*...* Arguments passed to `tmbstan`

*silent* Logical. If FALSE, all tracing outputs are shown (default).

**Method** `fit()`: Model fitting

The negative log-likelihood of the model is minimised using the function `optimx`. TMB uses the Laplace approximation to integrate the random effects out of the likelihood.

After the model has been fitted, the output of `optim` can be accessed using the method `out`.

*Usage:*

```
HMM$fit(silent = FALSE, ...)
```

*Arguments:*

`silent` Logical. If `FALSE`, all tracing outputs are shown (default).

`...` Other arguments to `optimx` which is used to optimise likelihood, see `?optimx`

*Examples:*

```
# Load data set (included with R)
data(nottem)
data <- data.frame(temp = as.vector(t(nottem)))

# Create hidden state and observation models
hid <- MarkovChain$new(data = data, n_states = 2)
par0 <- list(temp = list(mean = c(40, 60), sd = c(5, 5)))
obs <- Observation$new(data = data, n_states = 2,
                       dists = list(temp = "norm"),
                       par = par0)

# Create HMM
hmm <- HMM$new(hid = hid, obs = obs)

# Fit HMM
hmm$fit(silent = TRUE)
```

**Method** `mle()`: Get maximum likelihood estimates once model fitted*Usage:*

```
HMM$mle()
```

*Returns:* list of maximum likelihood estimates as described as input for the function `update_par()`

**Method** `forward_backward()`: Forward-backward algorithm

The forward probability for time step  $t$  and state  $j$  is the joint pdf/pmf of observations up to time  $t$  and of being in state  $j$  at time  $t$ ,  $p(Z[1], Z[2], \dots, Z[t], S[t] = j)$ . The backward probability for time  $t$  and state  $j$  is the conditional pdf/pmf of observations between time  $t + 1$  and  $n$ , given state  $j$  at time  $t$ ,  $p(Z[t+1], Z[t+2], \dots, Z[n] \mid S[t] = j)$ . This function returns their logarithm, for use in other methods `state_probs`, and `sample_states`.

*Usage:*

```
HMM$forward_backward()
```

*Returns:* Log-forward and log-backward probabilities

**Method** `cond()`: Compute conditional cumulative distribution functions*Usage:*

```
HMM$cond(ngrid = 1000, silent = FALSE)
```

*Arguments:*

`ngrid` how many cells on the grid that CDF is computed on

`silent` if TRUE then no messages are printed

*Returns:* cdfs on grid for each variable

**Method** `pseudores()`: Pseudo-residuals

Compute pseudo-residuals for the fitted model. If the fitted model is the "true" model, the pseudo-residuals follow a standard normal distribution. Deviations from normality suggest lack of fit.

*Usage:*

```
HMM$pseudores()
```

*Returns:* Matrix of pseudo-residuals, with one row for each response variable and one column for each observation

**Method** `viterbi()`: Viterbi algorithm

*Usage:*

```
HMM$viterbi()
```

*Returns:* Most likely state sequence

**Method** `sample_states()`: Sample posterior state sequences using forward-filtering backward-sampling

The forward-filtering backward-sampling algorithm returns a sequence of states, similarly to the Viterbi algorithm, but it generates it from the posterior distribution of state sequences, i.e., accounting for uncertainty in the state classification. Multiple generated sequences will therefore generally not be the same.

*Usage:*

```
HMM$sample_states(nsamp = 1, full = FALSE)
```

*Arguments:*

`nsamp` Number of samples to produce

`full` If TRUE and model fit by `fit_stan` then parameter estimates are sampled from the posterior samples before simulating each sequence

*Returns:* Matrix where each column is a different sample of state sequences, and each row is a time of observation

**Method** `state_probs()`: Compute posterior probability of being in each state

*Usage:*

```
HMM$state_probs()
```

*Returns:* matrix with a row for each observation and a column for each state

**Method** `post_coeff()`: Posterior sampling for model coefficients

*Usage:*

```
HMM$post_coeff(n_post)
```

*Arguments:*



`n_post` Number of posterior samples

*Returns:* Matrix with one column for each coefficient and one row for each posterior draw

**Method** `post_linpred()`: Posterior sampling for linear predictor

*Usage:*

```
HMM$post_linpred(n_post)
```

*Arguments:*

`n_post` Number of posterior samples

*Returns:* List with elements `obs` and `hid`, where each is a matrix with one column for each predictor and one row for each posterior draw

**Method** `post_fn()`: Create posterior simulations of a function of a model component

*Usage:*

```
HMM$post_fn(fn, n_post, comp = NULL, ..., level = 0, return_post = FALSE)
```

*Arguments:*

`fn` Function which takes a vector of linear predictors as input and produces either a scalar or vector output

`n_post` Number of posterior simulations

`comp` Either "obs" for observation model linear predictor, or "hid" for hidden model linear predictor

`...` Arguments passed to `fn`

`level` Confidence interval level if required (e.g., 0.95 for 95 confidence intervals). Default is 0, i.e., confidence intervals are not returned.

`return_post` Logical indicating whether to return the posterior samples. If FALSE (default), only mean estimates and confidence intervals are returned

*Returns:* A list with elements:

- `post`If `return_post = TRUE`, this is a vector (for scalar outputs of `fn`) or matrix (for vector outputs) with a column for each simulation
- `mean`Mean over posterior samples
- `lcl`Lower confidence interval bound (if `level != 0`)
- `ucl`Upper confidence interval bound (if `level != 0`)

**Method** `predict()`: Predict estimates from a fitted model

By default, this returns point estimates of the HMM parameters for a new data frame of covariates. See the argument 'n\_post' to also get confidence intervals.

*Usage:*

```
HMM$predict(
  what,
  t = 1,
  newdata = NULL,
  n_post = 0,
  level = 0.95,
  return_post = FALSE
)
```

*Arguments:*

*what* Which estimates to predict? Options include transition probability matrices "tpm", stationary distributions "delta", or observation distribution parameters "obspar"

*t* Time points to predict at

*newdata* New dataframe to use for prediction

*n\_post* If greater than zero then *n\_post* posterior samples are produced, and used to create confidence intervals.

*level* Level of the confidence intervals, e.g. CI = 0.95 will produce 95% confidence intervals (default)

*return\_post* Logical. If TRUE, a list of posterior samples is returned.

... Other arguments to the respective functions for *hid\$tpm*, *hid\$delta*, *obs\$par*

*Returns:* Named array of predictions and confidence intervals, if requested

*Examples:*

```
# Load data set (included with R)
data(nottem)
data <- data.frame(temp = as.vector(t(nottem)))

# Create hidden state and observation models
hid <- MarkovChain$new(data = data, n_states = 2)
par0 <- list(temp = list(mean = c(40, 60), sd = c(5, 5)))
obs <- Observation$new(data = data, n_states = 2,
                       dists = list(temp = "norm"),
                       par = par0)

# Create HMM
hmm <- HMM$new(hid = hid, obs = obs)

# Fit HMM
hmm$fit(silent = TRUE)

# Get transition probability matrix with confidence intervals
hmm$predict(what = "tpm", n_post = 1000)
```

**Method** `confint()`: Confidence intervals for working parameters

This function computes standard errors for all fixed effect model parameters based on the diagonal of the inverse of the Hessian matrix, and then derives Wald-type confidence intervals.

*Usage:*

```
HMM$confint(level = 0.95)
```

*Arguments:*

*level* Level of confidence intervals. Defaults to 0.95, i.e., 95% confidence intervals.

*Returns:* List of matrices with three columns: mle (maximum likelihood estimate), lcl (lower confidence limit), and ucl (upper confidence limit). One such matrix is produced for the working parameters of the observation model, the working parameters of the hidden state model, the smoothness parameters of the observation model, and the smoothness parameters of the hidden state model.

**Method** `simulate()`: Simulate from hidden Markov model

*Usage:*

```
HMM$simulate(n, data = NULL, silent = FALSE)
```

*Arguments:*

`n` Number of time steps to simulate  
`data` Optional data frame including covariates  
`silent` if TRUE then no messages are printed

*Returns:* Data frame including columns of data (if provided), and simulated data variables

**Method** `check()`: Compute goodness-of-fit statistics using simulation

Many time series are simulated from the fitted model, and the statistic(s) of interest are calculated for each. A histogram of those values can for example be used to compare to the observed value of the statistic. An observation far in the tails of the distribution of simulated statistics suggests lack of fit.

*Usage:*

```
HMM$check(check_fn, nsims = 100, full = FALSE, silent = FALSE)
```

*Arguments:*

`check_fn` Goodness-of-fit function which accepts "data" as input and returns a statistic (either a vector or a single number) to be compared between observed data and simulations.  
`nsims` Number of simulations to perform  
`full` If model fitted with 'fit\_stan', then `full = TRUE` will sample from posterior for each simulation  
`silent` Logical. If FALSE, simulation progress is shown. (Default: TRUE)

*Returns:* List with elements:

- `obs_stat`: Vector of values of goodness-of-fit statistics for the observed data
- `stats`: Matrix of values of goodness-of-fit statistics for the simulated data sets (one row for each statistic, and one column for each simulation)
- `plot`: ggplot object

**Method** `plot_ts()`: Time series plot coloured by states

Creates a plot of the data coloured by the most likely state sequence, as estimated by the Viterbi algorithm. If one variable name is passed as input, it is plotted against time. If two variables are passed, they are plotted against each other.

*Usage:*

```
HMM$plot_ts(var, var2 = NULL)
```

*Arguments:*

`var` Name of the variable to plot.  
`var2` Optional name of a second variable, for 2-d plot.

*Returns:* A ggplot object

**Method** `plot_dist()`: Plot observation distributions weighted by frequency in Viterbi

This is a wrapper around `Observation$plot_dist`, where the distribution for each state is weighted by the proportion of time spent in that state (according to the Viterbi state sequence).

*Usage:*

```
HMM$plot_dist(var)
```

*Arguments:*

var Name of data variable

*Returns:* Plot of distribution with data histogram

**Method plot():** Plot a model component

*Usage:*

```
HMM$plot(
  what,
  var = NULL,
  covs = NULL,
  i = NULL,
  j = NULL,
  n_grid = 50,
  n_post = 1000
)
```

*Arguments:*

what Name of model component to plot: should be one of "tpm" (transition probabilities), "delta" (stationary state probabilities), or "obspar" (state-dependent observation parameters)

var Name of covariate to plot on x-axis

covs Optional named list for values of covariates (other than 'var') that should be used in the plot (or dataframe with single row). If this is not specified, the mean value is used for numeric variables, and the first level for factor variables.

i If plotting tpm then rows of tpm; if plotting delta then indices of states to plot; if plotting obspar then full names of parameters to plot (e.g., obsvar.mean)

j If plotting tpm then columns of tpm to plot; if plotting delta then this is ignored; if plotting obspar then indices of states to plot

n\_grid Number of points in grid over x-axis (default: 50)

n\_post Number of posterior simulations to use when computing confidence intervals; default: 1000. See predict function for more detail.

*Returns:* A ggplot object

**Method AIC\_marginal():** Marginal Akaike Information Criterion

The marginal AIC is for example defined by Wood (2017), as  $AIC = -2L + 2k$  where L is the maximum marginal log-likelihood (of fixed effects), and k is the number of degrees of freedom of the fixed effect component of the model

*Usage:*

```
HMM$AIC_marginal()
```

*Returns:* Marginal AIC

**Method AIC\_conditional():** Conditional Akaike Information Criterion

The conditional AIC is for example defined by Wood (2017), as  $AIC = -2L + 2k$  where L is the maximum joint log-likelihood (of fixed and random effects), and k is the number of effective degrees of freedom of the model (accounting for flexibility in non-parametric terms implied by smoothing)

*Usage:*

HMM\$AIC\_conditional()

*Returns:* Conditional AIC

**Method** print\_obspar(): Print observation parameters at t = 1

*Usage:*

HMM\$print\_obspar()

**Method** print\_tpm(): Print observation parameters at t = 1

*Usage:*

HMM\$print\_tpm()

**Method** formulation(): Print model formulation

*Usage:*

HMM\$formulation()

**Method** print(): Print HMM object

*Usage:*

HMM\$print()

**Method** clone(): The objects of this class are cloneable with this method.

*Usage:*

HMM\$clone(deep = FALSE)

*Arguments:*

deep Whether to make a deep clone.

## Examples

```
## -----
## Method `HMM$new`
## -----

# Load data set (included with R)
data(nottem)
data <- data.frame(temp = as.vector(t(nottem)))

# Create hidden state and observation models
hid <- MarkovChain$new(data = data, n_states = 2)
par0 <- list(temp = list(mean = c(40, 60), sd = c(5, 5)))
obs <- Observation$new(data = data, n_states = 2,
                      dists = list(temp = "norm"),
                      par = par0)

# Create HMM
hmm <- HMM$new(hid = hid, obs = obs)

## -----
```

```

## Method `HMM$fit`
## -----

# Load data set (included with R)
data(nottem)
data <- data.frame(temp = as.vector(t(nottem)))

# Create hidden state and observation models
hid <- MarkovChain$new(data = data, n_states = 2)
par0 <- list(temp = list(mean = c(40, 60), sd = c(5, 5)))
obs <- Observation$new(data = data, n_states = 2,
                      dists = list(temp = "norm"),
                      par = par0)

# Create HMM
hmm <- HMM$new(hid = hid, obs = obs)

# Fit HMM
hmm$fit(silent = TRUE)

## -----
## Method `HMM$predict`
## -----

# Load data set (included with R)
data(nottem)
data <- data.frame(temp = as.vector(t(nottem)))

# Create hidden state and observation models
hid <- MarkovChain$new(data = data, n_states = 2)
par0 <- list(temp = list(mean = c(40, 60), sd = c(5, 5)))
obs <- Observation$new(data = data, n_states = 2,
                      dists = list(temp = "norm"),
                      par = par0)

# Create HMM
hmm <- HMM$new(hid = hid, obs = obs)

# Fit HMM
hmm$fit(silent = TRUE)

# Get transition probability matrix with confidence intervals
hmm$predict(what = "tpm", n_post = 1000)

```

---

hmmTMB\_cols

*hmmTMB colour palette*


---

## Description

p hmmTMB colour palette

**Usage**

hmmTMB\_cols

**Format**

An object of class character of length 6.

---

invmlgit	<i>Multivariate inverse logit function</i>
----------	--

---

**Description**

Multivariate inverse logit function

**Usage**

invmlgit(x)

**Arguments**

x                    Numeric vector

---

is_whole_number	<i>Check if number of whole number</i>
-----------------	--

---

**Description**

Check if number of whole number

**Usage**

is\_whole\_number(x, tol = 1e-10)

**Arguments**

x                    number to check or vector of numbers  
tol                    how far away from whole number is ok?

**Value**

TRUE if it is a whole number within tolerance

---

logLik.HMM	<i>logLik function for SDE objects</i>
------------	--

---

**Description**

This function makes it possible to call generic R methods such as AIC and BIC on HMM objects. It is based on the number of degrees of freedom of the \*conditional\* AIC (rather than marginal AIC), i.e., including degrees of freedom from the smooth/random effect components of the model.

**Usage**

```
## S3 method for class 'HMM'
logLik(object, ...)
```

**Arguments**

object	SDE model object
...	For compatibility with S3 method

**Value**

Maximum log-likelihood value for the model, with attributes df (degrees of freedom) and nobs (number of observations)

---

logsumexp	<i>Log of sum of exponentials</i>
-----------	-----------------------------------

---

**Description**

Log of sum of exponentials

**Usage**

```
logsumexp(x)
```

**Arguments**

x	Numeric vector
---	----------------



---

make_formulas	<i>Process formulas and store in nested list</i>
---------------	--

---

## Description

Process formulas and store in nested list

## Usage

```
make_formulas(input_forms, var_names, par_names, n_states)
```

## Arguments

input_forms	Nested list of formulas, with two levels: observed variable, and parameter of the observation distribution. The formulas can contain state-specific terms, e.g. " $\sim$ state1(x1) + x2".
var_names	character vector name of each observation variable
par_names	list with element for each observation variable that contains character vector of name of each parameter in its distribution
n_states	Number of states

## Details

Formulas for the observation parameters can be different for the different states, using special functions of the form "state1", "state2", etc. This method processes the list of formulas passed by the user to extract the state-specific formulas. Missing formulas are assumed to be intercept-only  $\sim 1$ .

## Value

Nested list of formulas, with three levels: observed variable, parameter of the observation distribution, and state.

## Examples

```
input_forms <- list(step = list(shape = ~ state1(x1) + x2,
                               scale = ~ x1),
                  count = list(lambda = ~ state1(x1) + state2(s(x2, bs = "cs"))))

make_formulas(input_forms = input_forms,
              var_names = names(input_forms),
              par_names = lapply(input_forms, names),
              n_states = 2)
```

---

make_matrices	<i>Create model matrices</i>
---------------	------------------------------

---

**Description**

Create model matrices

**Usage**

```
make_matrices(formulas, data, new_data = NULL)
```

**Arguments**

formulas	List of formulas (possibly nested, e.g. for use within Observation)
data	Data frame including covariates
new_data	Optional new data set, including covariates for which the design matrices should be created. This needs to be passed in addition to the argument 'data', for cases where smooth terms or factor covariates are included, and the original data set is needed to determine the full range of covariate values.

**Value**

A list of

- X\_fe Design matrix for fixed effects
- X\_re Design matrix for random effects
- S Smoothness matrix
- ncol\_fe Number of columns of X\_fe for each parameter
- ncol\_re Number of columns of X\_re and S for each random effect

---

MarkovChain	<i>R6 class for HMM hidden process model</i>
-------------	--

---

**Description**

R6 class for HMM hidden process model

R6 class for HMM hidden process model

**Details**

Contains the parameters and model formulas for the hidden process model.

**Methods****Public methods:**

- `MarkovChain$new()`
- `MarkovChain$formula()`
- `MarkovChain$formulas()`
- `MarkovChain$tpm()`
- `MarkovChain$ref()`
- `MarkovChain$ref_mat()`
- `MarkovChain$ref_delta0()`
- `MarkovChain$coeff_fe()`
- `MarkovChain$delta()`
- `MarkovChain$delta0()`
- `MarkovChain$stationary()`
- `MarkovChain$fixpar()`
- `MarkovChain$coeff_re()`
- `MarkovChain$X_fe()`
- `MarkovChain$X_re()`
- `MarkovChain$lambda()`
- `MarkovChain$sd_re()`
- `MarkovChain$nstates()`
- `MarkovChain$terms()`
- `MarkovChain$unique_ID()`
- `MarkovChain$initial_state()`
- `MarkovChain$update_tpm()`
- `MarkovChain$update_coeff_fe()`
- `MarkovChain$update_coeff_re()`
- `MarkovChain$update_X_fe()`
- `MarkovChain$update_X_re()`
- `MarkovChain$update_delta0()`
- `MarkovChain$update_lambda()`
- `MarkovChain$update_fixpar()`
- `MarkovChain$make_mat()`
- `MarkovChain$make_mat_grid()`
- `MarkovChain$tpm2par()`
- `MarkovChain$par2tpm()`
- `MarkovChain$linpred()`
- `MarkovChain$simulate()`
- `MarkovChain$formulation()`
- `MarkovChain$print()`
- `MarkovChain$clone()`

**Method** `new()`: Create new MarkovChain object

*Usage:*

```
MarkovChain$new(
  data,
  formula = NULL,
  n_states,
  tpm = NULL,
  initial_state = "estimated",
  fixpar = NULL,
  ref = 1:n_states
)
```

*Arguments:*

**data** Data frame, needed to create model matrices, and to identify the number of time series (which each have a separate initial distribution)

**formula** Either (1) R formula, used for all transition probabilities, or (2) matrix of character strings giving the formula for each transition probability, with "." along the diagonal (or for reference elements; see **ref** argument). (Default: no covariate dependence.)

**n\_states** Number of states. If not specified, then **formula** needs to be provided as a matrix, and **n\_states** is deduced from its dimensions.

**tpm** Optional transition probability matrix, to initialise the model parameters (intercepts in model with covariates). If not provided, the default is a matrix with 0.9 on the diagonal.

**initial\_state** Specify model for initial state distribution. There are five different options:

- "estimated": a separate initial distribution is estimated for each ID (default)
- "stationary": the initial distribution is fixed to the stationary distribution of the transition probability matrix for the first time point of each ID
- "shared": a common initial distribution is estimated for all IDs
- integer value between 1 and **n\_states**: used as the known initial state for all IDs
- vector of integers between 1 and **n\_states** (of length the number of IDs): each element is used as the known initial state for the corresponding ID

**fixpar** List with optional elements "hid" (fixed parameters for transition probabilities), "lambda\_hid" (fixed smoothness parameters), and "delta0" (fixed parameters for initial distribution). Each element is a named vector of coefficients that should either be fixed (if the corresponding element is set to NA) or estimated to a common value (using integers or factor levels).

**ref** Vector of indices for reference transition probabilities, of length **n\_states**. The *i*-th element is the index for the reference in the *i*-th row of the transition probability matrix. For example, **ref** = c(1, 1) means that the first element of the first row  $\text{Pr}(1>1)$  and the first element of the second row  $\text{Pr}(2>1)$  are used as reference elements and are not estimated. If this is not provided, the diagonal transition probabilities are used as references.

*Returns:* A new MarkovChain object

*Examples:*

```
# Load data set from MSwM package
data(energy, package = "MSwM")

# Create 2-state covariate-free model and initialise transition
# probability matrix
hid <- MarkovChain$new(data = energy, n_states = 2,
```

```

tpm = matrix(c(0.8, 0.3, 0.2, 0.7), 2, 2))

# Create 2-state model with non-linear effect of Oil on all transition
# probabilities
hid <- MarkovChain$new(data = energy, n_states = 2,
                      formula = ~ s(Oil, k = 5, bs = "cs"))

# Create 2-state model with quadratic effect of Oil on Pr(1 > 2)
structure <- matrix(c(".", "~poly(Oil, 2)",
                      "~1", "."),
                    ncol = 2, byrow = TRUE)
hid <- MarkovChain$new(data = energy, n_states = 2,
                      formula = structure)

```

**Method** formula(): Formula of MarkovChain model

*Usage:*

MarkovChain\$formula()

**Method** formulas(): List of formulas for MarkovChain model

*Usage:*

MarkovChain\$formulas()

**Method** tpm(): Get transition probability matrices

*Usage:*

MarkovChain\$tpm(t = 1, linpred = NULL)

*Arguments:*

t Time index or vector of time indices; default = 1. If t = "all" then all transition probability matrices are returned.

linpred Optional custom linear predictor

*Returns:* Array with one slice for each transition probability matrix

**Method** ref(): Indices of reference elements in transition probability matrix

*Usage:*

MarkovChain\$ref()

**Method** ref\_mat(): Matrix of reference elements in transition probability matrix

*Usage:*

MarkovChain\$ref\_mat()

**Method** ref\_delta0(): Indices of reference elements in initial distribution

*Usage:*

MarkovChain\$ref\_delta0()

**Method** coeff\_fe(): Current parameter estimates (fixed effects)

*Usage:*

MarkovChain\$coeff\_fe()

**Method** `delta()`: Stationary distribution

*Usage:*

```
MarkovChain$delta(t = NULL, linpred = NULL)
```

*Arguments:*

`t` Time point(s) for which stationary distribution should be returned. If `t = "all"`, all deltas are returned; else this should be a vector of time indices. If `NULL` (default), the stationary distribution for the first time step is returned.

`linpred` Optional custom linear predictor

*Returns:* Matrix of stationary distributions. Each row corresponds to a row of the design matrices, and each column corresponds to a state.

**Method** `delta0()`: Initial distribution

*Usage:*

```
MarkovChain$delta0(log = FALSE, as_matrix = TRUE)
```

*Arguments:*

`log` Logical indicating whether to return the log of the initial probabilities (default: `FALSE`). If `TRUE`, then the last element is excluded, as it is not estimated.

`as_matrix` Logical indicating whether the output should be formatted as a matrix (default). If `as_matrix` is `FALSE` and `log` is `TRUE`, the result is formatted as a column vector.

*Returns:* Matrix with one row for each time series ID, and one column for each state. For each ID, the *i*-th element of the corresponding row is the probability  $\Pr(S[1] = i)$

**Method** `stationary()`: Use stationary distribution as initial distribution?

*Usage:*

```
MarkovChain$stationary()
```

**Method** `fixpar()`: Fixed parameters

*Usage:*

```
MarkovChain$fixpar(all = FALSE)
```

*Arguments:*

`all` Logical. If `FALSE`, only user-specified fixed parameters are returned, but not parameters that are fixed for some other reason (e.g., from `'.'` in formula)

**Method** `coeff_re()`: Current parameter estimates (random effects)

*Usage:*

```
MarkovChain$coeff_re()
```

**Method** `X_fe()`: Fixed effect design matrix

*Usage:*

```
MarkovChain$X_fe()
```

**Method** `X_re()`: Random effect design matrix

*Usage:*

```
MarkovChain$X_re()
```

**Method** `lambda()`: Smoothness parameters

*Usage:*

`MarkovChain$lambda()`

**Method** `sd_re()`: Standard deviation of smooth terms

This function transforms the smoothness parameter of each smooth term into a standard deviation, given by  $SD = 1/\sqrt{\lambda}$ . It is particularly helpful to get the standard deviations of independent normal random effects.

*Usage:*

`MarkovChain$sd_re()`

**Method** `nstates()`: Number of states

*Usage:*

`MarkovChain$nstates()`

**Method** `terms()`: Terms of model formulas

*Usage:*

`MarkovChain$terms()`

**Method** `unique_ID()`: Number of time series

*Usage:*

`MarkovChain$unique_ID()`

**Method** `initial_state()`: Initial state (see constructor argument)

*Usage:*

`MarkovChain$initial_state()`

**Method** `update_tpm()`: Update transition probability matrix

*Usage:*

`MarkovChain$update_tpm(tpm)`

*Arguments:*

`tpm` New transition probability matrix

**Method** `update_coeff_fe()`: Update coefficients for fixed effect parameters

*Usage:*

`MarkovChain$update_coeff_fe(coeff_fe)`

*Arguments:*

`coeff_fe` Vector of coefficients for fixed effect parameters

**Method** `update_coeff_re()`: Update coefficients for random effect parameters

*Usage:*

`MarkovChain$update_coeff_re(coeff_re)`

*Arguments:*

`coeff_re` Vector of coefficients for random effect parameters

**Method** `update_X_fe()`: Update design matrix for fixed effects

*Usage:*

`MarkovChain$update_X_fe(X_fe)`

*Arguments:*

`X_fe` new design matrix for fixed effects

**Method** `update_X_re()`: Update design matrix for random effects

*Usage:*

`MarkovChain$update_X_re(X_re)`

*Arguments:*

`X_re` new design matrix for random effects

**Method** `update_delta0()`: Update initial distribution

*Usage:*

`MarkovChain$update_delta0(delta0)`

*Arguments:*

`delta0` Either a matrix where the *i*-th row is the initial distribution for the *i*-th time series in the data, or a vector which is then used for all time series. Entries of each row of `delta0` should sum to one.

**Method** `update_lambda()`: Update smoothness parameters

*Usage:*

`MarkovChain$update_lambda(lambda)`

*Arguments:*

`lambda` New smoothness parameter vector

**Method** `update_fixpar()`: Update information about fixed parameters

*Usage:*

`MarkovChain$update_fixpar(fixpar)`

*Arguments:*

`fixpar` New list of fixed parameters, in the same format expected by `MarkovChain$new()`

**Method** `make_mat()`: Make model matrices

*Usage:*

`MarkovChain$make_mat(data, new_data = NULL)`

*Arguments:*

`data` Data frame containing all needed covariates

`new_data` Optional new data set, including covariates for which the design matrices should be created. This needs to be passed in addition to the argument 'data', for cases where smooth terms or factor covariates are included, and the original data set is needed to determine the full range of covariate values.

*Returns:* A list with elements:



- `X_fe` Design matrix for fixed effects
- `X_re` Design matrix for random effects
- `SSmoothness` matrix for random effects
- `ncol_fe` Number of columns of `X_fe` for each parameter
- `ncol_re` Number of columns of `X_re` and `S` for each random effect

**Method** `make_mat_grid()`: Design matrices for grid of covariates

Used in plotting functions such as `HMM$plot_tpm` and `HMM$plot_stat_dist`

*Usage:*

```
MarkovChain$make_mat_grid(var, data, covs = NULL, n_grid = 1000)
```

*Arguments:*

`var` Name of variable

`data` Data frame containing the covariates

`covs` Optional named list for values of covariates (other than 'var') that should be used in the plot (or dataframe with single row). If this is not specified, the mean value is used for numeric variables, and the first level for factor variables.

`n_grid` Grid size (number of points). Default: 1000.

*Returns:* A list with the same elements as the output of `make_mat`, plus a data frame of covariates values.

**Method** `tpm2par()`: Transform transition probabilities to working scale

Apply the multinomial logit link function to get the corresponding parameters on the working scale (i.e., linear predictor scale).

*Usage:*

```
MarkovChain$tpm2par(tpm)
```

*Arguments:*

`tpm` Transition probability matrix

*Returns:* Vector of parameters on linear predictor scale

**Method** `par2tpm()`: Transform working parameters to transition probabilities

Apply the inverse multinomial logit link function to transform the parameters on the working scale (i.e., linear predictor scale) into the transition probabilities.

*Usage:*

```
MarkovChain$par2tpm(par)
```

*Arguments:*

`par` Vector of parameters on working scale

*Returns:* Transition probability matrix

**Method** `linpred()`: Linear predictor for transition probabilities

*Usage:*

```
MarkovChain$linpred()
```

**Method** `simulate()`: Simulate from Markov chain

*Usage:*

```
MarkovChain$simulate(n, data = NULL, new_data = NULL, silent = FALSE)
```

*Arguments:*

n Number of time steps to simulate

data Optional data frame containing all needed covariates

new\_data Optional new data set, including covariates for which the design matrices should be created. This needs to be passed in addition to the argument 'data', for cases where smooth terms or factor covariates are included, and the original data set is needed to determine the full range of covariate values.

silent if TRUE then no messages are printed

*Returns:* Sequence of states of simulated chain

**Method** formulation(): Print model formulation

*Usage:*

```
MarkovChain$formulation()
```

**Method** print(): Print MarkovChain object

*Usage:*

```
MarkovChain$print()
```

**Method** clone(): The objects of this class are cloneable with this method.

*Usage:*

```
MarkovChain$clone(deep = FALSE)
```

*Arguments:*

deep Whether to make a deep clone.

**Examples**

```
## -----
## Method `MarkovChain$new`
## -----

# Load data set from MSwM package
data(energy, package = "MSwM")

# Create 2-state covariate-free model and initialise transition
# probability matrix
hid <- MarkovChain$new(data = energy, n_states = 2,
                      tpm = matrix(c(0.8, 0.3, 0.2, 0.7), 2, 2))

# Create 2-state model with non-linear effect of Oil on all transition
# probabilities
hid <- MarkovChain$new(data = energy, n_states = 2,
                      formula = ~ s(Oil, k = 5, bs = "cs"))

# Create 2-state model with quadratic effect of Oil on Pr(1 > 2)
```

```
structure <- matrix(c(".", "~poly(Oil, 2)",
                    "~1", "."),
                  ncol = 2, byrow = TRUE)
hid <- MarkovChain$new(data = energy, n_states = 2,
                      formula = structure)
```

---

mlogit	<i>Multivariate logit function</i>
--------	------------------------------------

---

**Description**

Multivariate logit function

**Usage**

```
mlogit(x)
```

**Arguments**

x	Numeric vector
---	----------------

---

mvnorm_invlink	<i>Multivariate Normal inverse link function</i>
----------------	--

---

**Description**

Multivariate Normal inverse link function

**Usage**

```
mvnorm_invlink(x)
```

**Arguments**

x	Vector of parameters on linear predictor scale (in the order: means, SDs, correlations)
---	---

---

mvnorm_link	<i>Multivariate Normal link function</i>
-------------	--

---

**Description**

Multivariate Normal link function

**Usage**

```
mvnorm_link(x)
```

**Arguments**

x                      Vector of parameters on natural scale (in the order: means, SDs, correlations)

---

na_fill	<i>Fill in NAs</i>
---------	--------------------

---

**Description**

Replace NA entries in a vector by the last non-NA value. If the first entry of the vector is NA, it is replaced by the first non-NA value. If the vector passed as input doesn't contain NAs, it is returned as is.

**Usage**

```
na_fill(x)
```

**Arguments**

x                      Vector in which NAs should be removed

**Value**

Copy of x in which NAs have been replaced by nearest available value.

---

Observation

*R6 class for HMM observation model*

---

### Description

R6 class for HMM observation model

R6 class for HMM observation model

### Details

Contains the data, distributions, parameters, and formulas for the observation model from a hidden Markov model.

### Methods

#### Public methods:

- `Observation$new()`
- `Observation$data()`
- `Observation$dists()`
- `Observation$nstates()`
- `Observation$par()`
- `Observation$inipar()`
- `Observation$coeff_fe()`
- `Observation$coeff_re()`
- `Observation$X_fe()`
- `Observation$X_re()`
- `Observation$lambda()`
- `Observation$sd_re()`
- `Observation$formulas()`
- `Observation$terms()`
- `Observation$obs_var()`
- `Observation$known_states()`
- `Observation$fixpar()`
- `Observation$update_par()`
- `Observation$update_coeff_fe()`
- `Observation$update_coeff_re()`
- `Observation$update_X_fe()`
- `Observation$update_X_re()`
- `Observation$update_lambda()`
- `Observation$update_data()`
- `Observation$update_fixpar()`
- `Observation$make_mat()`

- `Observation$make_newdata_grid()`
- `Observation$n2w()`
- `Observation$w2n()`
- `Observation$linpred()`
- `Observation$obs_probs()`
- `Observation$suggest_initial()`
- `Observation$plot_dist()`
- `Observation$formulation()`
- `Observation$print()`
- `Observation$clone()`

**Method** `new()`: Create new Observation object

*Usage:*

```
Observation$new(data, dists, formulas = NULL, n_states, par, fixpar = NULL)
```

*Arguments:*

`data` Data frame containing response variables (named in `dists` and `par`) and covariates (named in `formulas`)

`dists` Named list of distribution names for each data stream, with the following options: `beta`, `binom`, `cat`, `dir`, `exp`, `foldednorm`, `gamma`, `gamma2`, `lnorm`, `mvnorm`, `nbinom`, `norm`, `pois`, `t`, `truncnorm`, `tweedie`, `vm`, `weibull`, `wrpcauchy`, `zibinom`, `zgamma`, `zgamma2`, `zinbinom`, `zipois`, `ztnbinom`, `ztpois`. See vignette about list of distributions for more detail, e.g., list of parameters for each distribution.

`formulas` List of formulas for observation parameters. This should be a nested list, where the outer list has one element for each observed variable, and the inner lists have one element for each parameter. Any parameter that is not included is assumed to have the formula `~1`. By default, all parameters have the formula `~1` (i.e., no covariate effects).

`n_states` Number of states (needed to construct model formulas)

`par` List of initial observation parameters. This should be a nested list, where the outer list has one element for each observed variable, and the inner lists have one element for each parameter. The choice of good initial values can be important, especially for complex models; the package vignettes discuss approaches to selecting them (e.g., see `Observation$suggest_initial()`).

`fixpar` List with optional elements `"obs"` (fixed coefficients for observation parameters), and `"lambda_obs"` (fixed smoothness parameters). Each element is a named vector of coefficients that should either be fixed (if the corresponding element is set to `NA`) or estimated to a common value (using integers or factor levels).

*Returns:* A new Observation object

*Examples:*

```
# Load data set from MSwM package
data(energy, package = "MSwM")

# Initial observation parameters
par0 <- list(Price = list(mean = c(3, 6), sd = c(2, 2)))

# Model "energy" with normal distributions
```

```

obs <- Observation$new(data = energy,
                      dists = list(Price = "norm"),
                      par = par0,
                      n_states = 2)

# Model "energy" with gamma distributions
obs <- Observation$new(data = energy,
                      dists = list(Price = "gamma2"),
                      par = par0,
                      n_states = 2)

# Model with non-linear effect of EurDol on mean price
f <- list(Price = list(mean = ~ s(EurDol, k = 5, bs = "cs")))
obs <- Observation$new(data = energy,
                      dists = list(Price = "norm"),
                      par = par0,
                      n_states = 2,
                      formula = f)

```

**Method** `data()`: Data frame

*Usage:*

```
Observation$data()
```

**Method** `dists()`: List of distributions

*Usage:*

```
Observation$dists()
```

**Method** `nstates()`: Number of states

*Usage:*

```
Observation$nstates()
```

**Method** `par()`: Parameters on natural scale

*Usage:*

```
Observation$par(t = 1, full_names = TRUE, linpred = NULL, as_list = FALSE)
```

*Arguments:*

`t` Time index or vector of time indices; default `t = 1`. If `t = "all"`, then return observation parameters for all time points.

`full_names` Logical. If `TRUE`, the rows of the output are named in the format "variable.parameter" (default). If `FALSE`, the rows are names in the format "parameter". The latter is used in various internal functions, when the parameters need to be passed on to an R function.

`linpred` Optional custom linear predictor.

`as_list` Logical. If `TRUE`, the output is a nested list with three levels: (1) time step, (2) observed variable, (3) observation parameter. If `FALSE` (default), the output is an array with one row for each observation parameter, one column for each state, and one slice for each time step.

*Returns:* Array of parameters with one row for each observation parameter, one column for each state, and one slice for each time step. (See `as_list` argument for alternative output format.)

*Examples:*

```
# Load data set from MSwM package
data(energy, package = "MSwM")

# Initial observation parameters
par0 <- list(Price = list(mean = c(3, 6), sd = c(2, 2)))

# Model with linear effect of EurDol on mean price
f <- list(Price = list(mean = ~ EurDol))
obs <- Observation$new(data = energy,
                      dists = list(Price = "norm"),
                      par = par0,
                      n_states = 2,
                      formula = f)

# Set slope coefficients
obs$update_coeff_fe(coeff_fe = c(3, 2, 6, -2, log(2), log(2)))

# Observation parameter values for given data rows
obs$par(t = c(1, 10, 20))
```

**Method** `inipar()`: Return initial parameter values supplied

*Usage:*

```
Observation$inipar()
```

**Method** `coeff_fe()`: Fixed effect parameters on working scale

*Usage:*

```
Observation$coeff_fe()
```

**Method** `coeff_re()`: Random effect parameters

*Usage:*

```
Observation$coeff_re()
```

**Method** `X_fe()`: Fixed effect design matrix

*Usage:*

```
Observation$X_fe()
```

**Method** `X_re()`: Random effect design matrix

*Usage:*

```
Observation$X_re()
```

**Method** `lambda()`: Smoothness parameters

*Usage:*

```
Observation$lambda()
```

**Method** `sd_re()`: Standard deviation of smooth terms

This function transforms the smoothness parameter of each smooth term into a standard deviation, given by  $SD = 1/\sqrt{\lambda}$ . It is particularly helpful to get the standard deviations of independent normal random effects.



*Usage:*

```
Observation$sd_re()
```

**Method** `formulas()`: List of model formulas for observation model

*Usage:*

```
Observation$formulas(raw = FALSE)
```

*Arguments:*

`raw` Logical. If FALSE, returns the nested list created by `make_formulas` (default). If TRUE, returns formulas passed as input.

**Method** `terms()`: Terms of model formulas

*Usage:*

```
Observation$terms()
```

**Method** `obs_var()`: Data frame of response variables

*Usage:*

```
Observation$obs_var(expand = FALSE)
```

*Arguments:*

`expand` If TRUE, then multivariate variables in observations are expanded to be univariate, creating extra columns.

*Returns:* Data frame of observation variables

**Method** `known_states()`: Vector of known states

*Usage:*

```
Observation$known_states(mat = TRUE)
```

*Arguments:*

`mat` Logical.

**Method** `fixpar()`: Fixed parameters

*Usage:*

```
Observation$fixpar(all = FALSE)
```

*Arguments:*

`all` Logical. If FALSE, only user-specified fixed parameters are returned, but not parameters that are fixed for some other reason (e.g., size of binomial distribution)

**Method** `update_par()`: Update parameters

Updates the 'par' attribute to the list passed as input, and updates the intercept elements of 'coeff\_fe' using the list passed as input

*Usage:*

```
Observation$update_par(par)
```

*Arguments:*

`par` New list of parameters

**Method** `update_coeff_fe()`: Update coefficients for fixed effect parameters

*Usage:*

```
Observation$update_coeff_fe(coeff_fe)
```

*Arguments:*

`coeff_fe` New vector of coefficients for fixed effect parameters

**Method** `update_coeff_re()`: Update random effect parameters

*Usage:*

```
Observation$update_coeff_re(coeff_re)
```

*Arguments:*

`coeff_re` New vector of coefficients for random effect parameters

**Method** `update_X_fe()`: Update fixed effect design matrix

*Usage:*

```
Observation$update_X_fe(X_fe)
```

*Arguments:*

`X_fe` New fixed effect design matrix

**Method** `update_X_re()`: Update random effect design matrix

*Usage:*

```
Observation$update_X_re(X_re)
```

*Arguments:*

`X_re` New random effect design matrix

**Method** `update_lambda()`: Update smoothness parameters

*Usage:*

```
Observation$update_lambda(lambda)
```

*Arguments:*

`lambda` New smoothness parameter vector

**Method** `update_data()`: Update data

*Usage:*

```
Observation$update_data(data)
```

*Arguments:*

`data` New data frame

**Method** `update_fixpar()`: Update information about fixed parameters

*Usage:*

```
Observation$update_fixpar(fixpar)
```

*Arguments:*

`fixpar` New list of fixed parameters, in the same format expected by `Observation$new()`

**Method** `make_mat()`: Make model matrices

*Usage:*

```
Observation$make_mat(new_data = NULL)
```

*Arguments:*

`new_data` Optional new data set, including covariates for which the design matrices should be created. If this argument is not specified, the design matrices are based on the original data frame.

*Returns:* A list with elements:

- `X_fe` Design matrix for fixed effects
- `X_re` Design matrix for random effects
- `SSmoothness` matrix for random effects
- `ncol_fe` Number of columns of `X_fe` for each parameter
- `ncol_re` Number of columns of `X_re` and `S` for each random effect

Design matrices for grid of covariates

**Method** `make_newdata_grid()`:

*Usage:*

```
Observation$make_newdata_grid(var, covs = NULL, n_grid = 1000)
```

*Arguments:*

`var` Name of variable

`covs` Optional named list for values of covariates (other than 'var') that should be used in the plot (or dataframe with single row). If this is not specified, the mean value is used for numeric variables, and the first level for factor variables.

`n_grid` Grid size (number of points). Default: 1000.

*Returns:* A list with the same elements as the output of `make_mat`, plus a data frame of covariates values.

**Method** `n2w()`: Natural to working parameter transformation

This function applies the link functions of the distribution parameters, to transform parameters from their natural scale to the working scale (i.e., linear predictor scale)

*Usage:*

```
Observation$n2w(par)
```

*Arguments:*

`par` List of parameters on natural scale

*Returns:* Vector of parameters on working scale

**Method** `w2n()`: Working to natural parameter transformation

This function applies the inverse link functions of the distribution parameters, to transform parameters from the working scale (i.e., linear predictor scale) to their natural scale.

*Usage:*

```
Observation$w2n(wpar)
```

*Arguments:*

wpar Vector of parameters on working scale

*Returns:* List of parameters on natural scale

**Method** linpred(): Compute linear predictor

*Usage:*

```
Observation$linpred()
```

**Method** obs\_probs(): Observation likelihoods

*Usage:*

```
Observation$obs_probs(data = NULL)
```

*Arguments:*

data Optional dataframe to include in form of obs\_var() output

*Returns:* Matrix of likelihoods of observations, with one row for each time step, and one column for each state.

**Method** suggest\_initial(): Suggest initial observation parameters

The K-means algorithm is used to define clusters of observations (supposed to approximate the HMM states). Then, for each cluster, the parapprox function of the relevant Dist object is used to obtain parameter values.

*Usage:*

```
Observation$suggest_initial()
```

*Returns:* List of initial parameters for each observation variable

*Examples:*

```
# Load data set from MSwM package
data(energy, package = "MSwM")

# Initial observation parameters
par0 <- list(Price = list(mean = c(3, 6), sd = c(2, 2)))

# Model "energy" with normal distributions
obs <- Observation$new(data = energy,
                      dists = list(Price = "norm"),
                      par = par0,
                      n_states = 2)

# Print observation parameters
obs$par()

# Suggest initial parameters
par0_new <- obs$suggest_initial()
par0_new

# Update model parameters to suggested
obs$update_par(par = par0_new)
obs$par()
```

**Method** `plot_dist()`: Plot histogram of data and pdfs

Plot histogram of observations for the variable specified by the argument name, overlaid with the pdf of the specified distribution for that data stream. Helpful to select initial parameter values for model fitting, or to visualise fitted state-dependent distributions.

*Usage:*

```
Observation$plot_dist(var, weights = NULL, t = 1)
```

*Arguments:*

`var` Name of response variable for which the histogram and pdfs should be plotted.

`weights` Optional vector of length the number of pdfs that are plotted. Useful to visualise a mixture of distributions weighted by the proportion of time spent in the different states.

`t` Index of time step to use for covariates (default: 1).

*Returns:* A ggplot object

**Method** `formulation()`: Print model formulation

*Usage:*

```
Observation$formulation()
```

**Method** `print()`: Print Observation object Check constructor arguments

*Usage:*

```
Observation$print()
```

**Method** `clone()`: The objects of this class are cloneable with this method.

*Usage:*

```
Observation$clone(deep = FALSE)
```

*Arguments:*

`deep` Whether to make a deep clone.

## Examples

```
## -----
## Method `Observation$new`
## -----

# Load data set from MSwM package
data(energy, package = "MSwM")

# Initial observation parameters
par0 <- list(Price = list(mean = c(3, 6), sd = c(2, 2)))

# Model "energy" with normal distributions
obs <- Observation$new(data = energy,
                      dists = list(Price = "norm"),
                      par = par0,
                      n_states = 2)

# Model "energy" with gamma distributions
```

```

obs <- Observation$new(data = energy,
                      dists = list(Price = "gamma2"),
                      par = par0,
                      n_states = 2)

# Model with non-linear effect of EurDol on mean price
f <- list(Price = list(mean = ~ s(EurDol, k = 5, bs = "cs")))
obs <- Observation$new(data = energy,
                      dists = list(Price = "norm"),
                      par = par0,
                      n_states = 2,
                      formula = f)

## -----
## Method `Observation$par`
## -----

# Load data set from MSwM package
data(energy, package = "MSwM")

# Initial observation parameters
par0 <- list(Price = list(mean = c(3, 6), sd = c(2, 2)))

# Model with linear effect of EurDol on mean price
f <- list(Price = list(mean = ~ EurDol))
obs <- Observation$new(data = energy,
                      dists = list(Price = "norm"),
                      par = par0,
                      n_states = 2,
                      formula = f)

# Set slope coefficients
obs$update_coeff_fe(coeff_fe = c(3, 2, 6, -2, log(2), log(2)))

# Observation parameter values for given data rows
obs$par(t = c(1, 10, 20))

## -----
## Method `Observation$suggest_initial`
## -----

# Load data set from MSwM package
data(energy, package = "MSwM")

# Initial observation parameters
par0 <- list(Price = list(mean = c(3, 6), sd = c(2, 2)))

# Model "energy" with normal distributions
obs <- Observation$new(data = energy,
                      dists = list(Price = "norm"),
                      par = par0,
                      n_states = 2)

```

```

# Print observation parameters
obs$par()

# Suggest initial parameters
par0_new <- obs$suggest_initial()
par0_new

# Update model parameters to suggested
obs$update_par(par = par0_new)
obs$par()

```

---

```
prec_to_cov
```

*Get covariance matrix from precision matrix*

---

### Description

The covariance matrix is the inverse of the precision matrix. By default, the function `solve` is used for inversion. If it fails (e.g., singular system), then `MASS::ginv` is used instead, and returns the Moore-Penrose generalised inverse of the precision matrix.

### Usage

```
prec_to_cov(prec_mat)
```

### Arguments

prec_mat	Precision matrix (either of 'matrix' type or sparse matrix on which <code>as.matrix</code> can be used)
----------	---

### Value

Precision matrix

---

```
quad_pos_solve
```

*Solve for positive root of quadratic  $ax^2 + bx + c = 0$  when it exists*

---

### Description

Solve for positive root of quadratic  $ax^2 + bx + c = 0$  when it exists

### Usage

```
quad_pos_solve(a, b, c)
```

**Arguments**

a	coefficient of $x^2$
b	coefficient of $x$
c	scalar coefficient

**Value**

real positive root if it exists

---

strip_comments	<i>Strip comments marked with a hash from a character vector</i>
----------------	--

---

**Description**

Strip comments marked with a hash from a character vector

**Usage**

```
strip_comments(str)
```

**Arguments**

str	the character vector
-----	----------------------

**Value**

character vector with comments removed (and lines with only comments completely removed)

---

update.HMM	<i>Update a model to a new model by changing one formula</i>
------------	--

---

**Description**

Update a model to a new model by changing one formula

**Usage**

```
## S3 method for class 'HMM'
update(object, type, i, j, change, fit = TRUE, silent = FALSE, ...)
```



**Arguments**

object	HMM model object
type	Character string for the part of the model that is updated (either "hid" or "obs")
i	If type = "hid" then i is the row of the formula containing the change. If type = "obs" then i is the observation variable name.
j	If type = "hid" then j is the column of the formula containing the change. If type = "obs" then j is the parameter whose formula is to be changed.
change	The change to make to the formula, see ?update.formula for details.
fit	If FALSE then change is made but model is not re-fit.
silent	If TRUE then no model fitting output is given
...	Additional arguments are ignored (for compatibility with generic S3 method)

**Examples**

```
# Load data set from MSwM package
data(energy, package = "MSwM")

# Create hidden state and observation models
hid <- MarkovChain$new(data = energy, n_states = 2)
par0 <- list(Price = list(mean = c(3, 6), sd = c(2, 3)))
obs <- Observation$new(data = energy, n_states = 2,
                      dists = list(Price = "norm"),
                      par = par0)

# Create HMM (no covariate effects)
hmm <- HMM$new(hid = hid, obs = obs)
hmm$hid()$formula()
hmm$obs()$formulas()

# Update transition probability formulas (one at a time)
hmm <- update(hmm, type = "hid", i = 1, j = 2,
             change = ~ . + Oil, fit = FALSE)
hmm <- update(hmm, type = "hid", i = 2, j = 1,
             change = ~ . + Gas + Coal, fit = FALSE)
hmm$hid()$formula()

# Update observation parameter formulas (one at a time)
hmm <- update(hmm, type = "obs", i = "Price", j = "mean",
             change = ~ . + EurDol, fit = FALSE)
hmm$obs()$formulas()
```

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