

# Package ‘msm’

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**Title** Multi-State Markov and Hidden Markov Models in Continuous Time

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**Description** Functions for fitting continuous-time Markov and hidden Markov multi-state models to longitudinal data. Designed for processes observed at arbitrary times in continuous time (panel data) but some other observation schemes are supported. Both Markov transition rates and the hidden Markov output process can be modelled in terms of covariates, which may be constant or piecewise-constant in time.

**License** GPL (>= 2)

**Imports** survival,mvtnorm,expm

**Suggests** mstate,minqa,doParallel,foreach,numDeriv,testthat,flexsurv

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2phase	<i>Coxian phase-type distribution with two phases</i>
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## Description

Density, distribution, quantile functions and other utilities for the Coxian phase-type distribution with two phases.

## Usage

```
d2phase(x, l1, mu1, mu2, log=FALSE)
p2phase(q, l1, mu1, mu2, lower.tail=TRUE, log.p=FALSE)
q2phase(p, l1, mu1, mu2, lower.tail=TRUE, log.p=FALSE)
r2phase(n, l1, mu1, mu2)
h2phase(x, l1, mu1, mu2, log=FALSE)
```

## Arguments

x, q	vector of quantiles.
p	vector of probabilities.
n	number of observations. If <code>length(n) &gt; 1</code> , the length is taken to be the number required.
l1	Intensity for transition between phase 1 and phase 2.
mu1	Intensity for transition from phase 1 to exit.
mu2	Intensity for transition from phase 2 to exit.
log	logical; if TRUE, return log density or log hazard.
log.p	logical; if TRUE, probabilities p are given as <code>log(p)</code> .
lower.tail	logical; if TRUE (default), probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .

### Details

This is the distribution of the time to reach state 3 in a continuous-time Markov model with three states and transitions permitted from state 1 to state 2 (with intensity  $\lambda_1$ ) state 1 to state 3 (intensity  $\mu_1$ ) and state 2 to state 3 (intensity  $\mu_2$ ). States 1 and 2 are the two "phases" and state 3 is the "exit" state.

The density is

$$f(t|\lambda_1, \mu_1) = e^{-(\lambda_1 + \mu_1)t}(\mu_1 + (\lambda_1 + \mu_1)\lambda_1 t)$$

if  $\lambda_1 + \mu_1 = \mu_2$ , and

$$f(t|\lambda_1, \mu_1, \mu_2) = \frac{(\lambda_1 + \mu_1)e^{-(\lambda_1 + \mu_1)t}(\mu_2 - \mu_1) + \mu_2\lambda_1 e^{-\mu_2 t}}{\lambda_1 + \mu_1 - \mu_2}$$

otherwise. The distribution function is

$$F(t|\lambda_1, \mu_1) = 1 - e^{-(\lambda_1 + \mu_1)t}(1 + \lambda_1 t)$$

if  $\lambda_1 + \mu_1 = \mu_2$ , and

$$F(t|\lambda_1, \mu_1, \mu_2) = 1 - \frac{e^{-(\lambda_1 + \mu_1)t}(\mu_2 - \mu_1) + \lambda_1 e^{-\mu_2 t}}{\lambda_1 + \mu_1 - \mu_2}$$

otherwise. Quantiles are calculated by numerically inverting the distribution function.

The mean is  $(1 + \lambda_1/\mu_2)/(\lambda_1 + \mu_1)$ .

The variance is  $(2 + 2\lambda_1(\lambda_1 + \mu_1 + \mu_2)/\mu_2^2 - (1 + \lambda_1/\mu_2)^2)/(\lambda_1 + \mu_1)^2$ .

If  $\mu_1 = \mu_2$  it reduces to an exponential distribution with rate  $\mu_1$ , and the parameter  $\lambda_1$  is redundant. Or also if  $\lambda_1 = 0$ .

The hazard at  $x = 0$  is  $\mu_1$ , and smoothly increasing if  $\mu_1 < \mu_2$ . If  $\lambda_1 + \mu_1 \geq \mu_2$  it increases to an asymptote of  $\mu_2$ , and if  $\lambda_1 + \mu_1 \leq \mu_2$  it increases to an asymptote of  $\lambda_1 + \mu_1$ . The hazard is decreasing if  $\mu_1 > \mu_2$ , to an asymptote of  $\mu_2$ .

### Value

d2phase gives the density, p2phase gives the distribution function, q2phase gives the quantile function, r2phase generates random deviates, and h2phase gives the hazard.

### Alternative parameterisation

An individual following this distribution can be seen as coming from a mixture of two populations:

1) "short stayers" whose mean sojourn time is  $M_1 = 1/(\lambda_1 + \mu_1)$  and sojourn distribution is exponential with rate  $\lambda_1 + \mu_1$ .

2) "long stayers" whose mean sojourn time  $M_2 = 1/(\lambda_1 + \mu_1) + 1/\mu_2$  and sojourn distribution is the sum of two exponentials with rate  $\lambda_1 + \mu_1$  and  $\mu_2$  respectively. The individual is a "long stayer" with probability  $p = \lambda_1/(\lambda_1 + \mu_1)$ .

Thus a two-phase distribution can be more intuitively parameterised by the short and long stay means  $M_1 < M_2$  and the long stay probability  $p$ . Given these parameters, the transition intensities are  $\lambda_1 = p/M_1$ ,  $\mu_1 = (1 - p)/M_1$ , and  $\mu_2 = 1/(M_2 - M_1)$ . This can be useful for choosing intuitively reasonable initial values for procedures to fit these models to data.

The hazard is increasing at least if  $M_2 < 2M_1$ , and also only if  $(M_2 - 2M_1)/(M_2 - M_1) < p$ .

For increasing hazards with  $\lambda_1 + \mu_1 \leq \mu_2$ , the maximum hazard ratio between any time  $t$  and time 0 is  $1/(1 - p)$ .

For increasing hazards with  $\lambda_1 + \mu_1 \geq \mu_2$ , the maximum hazard ratio is  $M_1/((1 - p)(M_2 - M_1))$ . This is the minimum hazard ratio for decreasing hazards.

### General phase-type distributions

This is a special case of the n-phase Coxian phase-type distribution, which in turn is a special case of the (general) phase-type distribution. The **actuar** R package implements a general n-phase distribution defined by the time to absorption of a general continuous-time Markov chain with a single absorbing state, where the process starts in one of the transient states with a given probability.

### Author(s)

C. H. Jackson <chris.jackson@mrc-bsu.cam.ac.uk>

### References

C. Dutang, V. Goulet and M. Pigeon (2008). actuar: An R Package for Actuarial Science. Journal of Statistical Software, vol. 25, no. 7, 1-37. URL <http://www.jstatsoft.org/v25/i07>

---

aneur

*Aortic aneurysm progression data*

---

### Description

This dataset contains longitudinal measurements of grades of aortic aneurysms, measured by ultrasound examination of the diameter of the aorta.

### Usage

aneur

### Format

A data frame containing 4337 rows, with each row corresponding to an ultrasound scan from one of 838 men over 65 years of age.

ptnum	(numeric)	Patient identification number
age	(numeric)	Recipient age at examination (years)
diam	(numeric)	Aortic diameter
state	(numeric)	State of aneurysm.

The states represent successive degrees of aneurysm severity, as indicated by the aortic diameter.

State 1	Aneurysm-free	< 30 cm
State 2	Mild aneurysm	30-44 cm
State 3	Moderate aneurysm	45-54 cm
State 4	Severe aneurysm	> 55 cm

683 of these men were aneurysm-free at age 65 and were re-screened every two years. The remaining men were aneurysmal at entry and had successive screens with frequency depending on the state of the aneurysm. Severe aneurysms are repaired by surgery.

### Source

The Chichester, U.K. randomised controlled trial of screening for abdominal aortic aneurysms by ultrasonography.

### References

Jackson, C.H., Sharples, L.D., Thompson, S.G. and Duffy, S.W. and Couto, E. Multi-state Markov models for disease progression with classification error. *The Statistician*, 52(2): 193–209 (2003)

Couto, E. and Duffy, S. W. and Ashton, H. A. and Walker, N. M. and Myles, J. P. and Scott, R. A. P. and Thompson, S. G. (2002) *Probabilities of progression of aortic aneurysms: estimates and implications for screening policy* Journal of Medical Screening 9(1):40–42

---

boot.msm

*Bootstrap resampling for multi-state models*

---

### Description

Draw a number of bootstrap resamples, refit a `msm` model to the resamples, and calculate statistics on the refitted models.

### Usage

```
boot.msm(x, stat=pmatrx.msm, B=1000, file=NULL, cores=NULL)
```

### Arguments

x	A fitted <code>msm</code> model, as output by <code>msm</code> .
stat	A function to call on each refitted <code>msm</code> model. By default this is <code>pmatrx.msm</code> , returning the transition probability matrix in one time unit. If <code>NULL</code> then no function is computed.
B	Number of bootstrap resamples.

file	Name of a file in which to save partial results after each replicate. This is saved using <code>save</code> and can be restored using <code>load</code> , producing an object called <code>boot.list</code> containing the partial results. Not supported when using parallel processing.
cores	Number of processor cores to use for parallel processing. Requires the <b>doParallel</b> package to be installed. If not specified, parallel processing is not used. If <code>cores</code> is set to the string "default", the default methods of <code>makeCluster</code> (on Windows) or <code>registerDoParallel</code> (on Unix-like) are used.

## Details

The bootstrap datasets are computed by resampling independent transitions between pairs of states (for non-hidden models without censoring), or independent individual series (for hidden models or models with censoring). Therefore this approach doesn't work if, for example, the data for a HMM consist of a series of observations from just one individual, and is inaccurate for small numbers of independent transitions or individuals.

Confidence intervals or standard errors for the corresponding statistic can be calculated by summarising the returned list of B replicated outputs. This is currently implemented for most of the output functions `qmatrix.msm`, `ematrix.msm`, `qratio.msm`, `pmatrix.msm`, `pmatrix.piecewise.msm`, `totlos.msm` and `prevalence.msm`. For other outputs, users will have to write their own code to summarise the output of `boot.msm`.

Most of **msm**'s output functions present confidence intervals based on asymptotic standard errors calculated from the Hessian. These are expected to be underestimates of the true standard errors (Cramer-Rao lower bound). Some of these functions use a further approximation, the delta method (see `deltamethod`) to obtain standard errors of transformed parameters. Bootstrapping should give a more accurate estimate of the uncertainty.

An alternative method which is less accurate though faster than bootstrapping, but more accurate than the delta method, is to draw a sample from the asymptotic multivariate normal distribution implied by the maximum likelihood estimates (and covariance matrix), and summarise the transformed estimates. See `pmatrix.msm`.

All objects used in the original call to `msm` which produced `x`, such as the `qmatrix`, should be in the working environment, or else `boot.msm` will produce an "object not found" error. This enables `boot.msm` to refit the original model to the replicate datasets. However there is currently a limitation. In the original call to `msm`, the "formula" argument should be specified directly, as, for example,

```
msm(state ~ time, data = ...)
```

and not, for example,

```
form = data$state ~ data$time
```

```
msm(formula=form, data = ...)
```

otherwise `boot.msm` will be unable to draw the replicate datasets.

`boot.msm` will also fail with an incomprehensible error if the original call to `msm` used a user-defined object whose name is the same as a built-in R object, or an object in any other loaded package. For example, if you have called a Q matrix `q`, when `q()` is the built-in function for quitting R.

If `stat` is `NULL`, then B different `msm` model objects will be stored in memory. This is inadvisable, as `msm` objects tend to be large, since they contain the original data used for the `msm` fit, so this will be wasteful of memory.

To specify more than one statistic, write a function consisting of a list of different function calls, for example,

```
stat = function(x) list (pmatrix.msm(x,t=1),pmatrix.msm(x,t=2))
```

### Value

A list with B components, containing the result of calling function `stat` on each of the refitted models. If `stat` is NULL, then each component just contains the refitted model. If one of the B model fits was unsuccessful and resulted in an error, then the corresponding list component will contain the error message.

### Author(s)

C.H.Jackson <chris.jackson@mrc-bsu.cam.ac.uk>

### References

Efron, B. and Tibshirani, R.J. (1993) *An Introduction to the Bootstrap*, Chapman and Hall.

### See Also

[qmatrix.msm](#), [qratio.msm](#), [sojourn.msm](#), [ematrix.msm](#), [pmatrix.msm](#), [pmatrix.piecewise.msm](#), [totlos.msm](#), [prevalence.msm](#).

### Examples

```
## Not run:
## Psoriatic arthritis example
data(psor)
psor.q <- rbind(c(0,0.1,0,0),c(0,0,0.1,0),c(0,0,0,0.1),c(0,0,0,0))
psor.msm <- msm(state ~ months, subject=ptnum, data=psor, qmatrix =
  psor.q, covariates = ~ollwsdrt+hieffusn,
  constraint = list(hieffusn=c(1,1,1),ollwsdrt=c(1,1,2)),
  control = list(REPORT=1,trace=2), method="BFGS")
## Bootstrap the baseline transition intensity matrix. This will take a long time.
q.list <- boot.msm(psor.msm, function(x)x$Qmatrices$baseline)
## Manipulate the resulting list of matrices to calculate bootstrap standard errors.
apply(array(unlist(q.list), dim=c(4,4,5)), c(1,2), sd)
## Similarly calculate a bootstrap 95% confidence interval
apply(array(unlist(q.list), dim=c(4,4,5)), c(1,2),
  function(x)quantile(x, c(0.025, 0.975)))
## Bootstrap standard errors are larger than the asymptotic standard
## errors calculated from the Hessian
psor.msm$QmatricesSE$baseline

## End(Not run)
```



---

bos *Bronchiolitis obliterans syndrome after lung transplants*

---

### Description

A dataset containing histories of bronchiolitis obliterans syndrome (BOS) from lung transplant recipients. BOS is a chronic decline in lung function, often observed after lung transplantation. The condition is classified into four stages of severity: none, mild, moderate and severe.

### Usage

bos

### Format

A data frame containing 638 rows, grouped by patient, including histories of 204 patients. The first observation for each patient is defined to be stage 1, no BOS, at six months after transplant. Subsequent observations denote the entry times into stages 2, 3, 4, representing mild, moderate and severe BOS respectively, and stage 5, representing death.

ptnum	(numeric)	Patient identification number
time	(numeric)	Months after transplant
state	(numeric)	BOS state entered at this time

### Details

The entry time of each patient into each stage of BOS was estimated by clinicians, based on their history of lung function measurements and acute rejection and infection episodes. BOS is only assumed to occur beyond six months after transplant. In the first six months the function of each patient's new lung stabilises. Subsequently BOS is diagnosed by comparing the lung function against the "baseline" value.

The objects bos3 and bos4 contain the same data, but with mild/moderate/severe combined, and moderate/severe combined, to give 3 and 4-state representations respectively.

### Source

Papworth Hospital, U.K.

### References

Heng, D. et al. (1998). Bronchiolitis Obliterans Syndrome: Incidence, Natural History, Prognosis, and Risk Factors. *Journal of Heart and Lung Transplantation* 17(12)1255–1263.

cav

*Heart transplant monitoring data***Description**

A series of approximately yearly angiographic examinations of heart transplant recipients. The state at each time is a grade of cardiac allograft vasculopathy (CAV), a deterioration of the arterial walls.

**Usage**

cav

**Format**

A data frame containing 2846 rows. There are 622 patients, the rows are grouped by patient number and ordered by years after transplant, with each row representing an examination and containing additional covariates.

PTNUM	(numeric)	Patient identification number
age	(numeric)	Recipient age at examination (years)
years	(numeric)	Examination time (years after transplant)
dage	(numeric)	Age of heart donor (years)
sex	(numeric)	sex (0=male, 1=female)
pdiag	(factor)	Primary diagnosis (reason for transplant) IHD=ischaemic heart disease, IDC=idiopathic dilated cardiomyopathy.
cumrej	(numeric)	Cumulative number of acute rejection episodes
state	(numeric)	State at the examination. State 1 represents no CAV, state 2 is mild/moderate CAV and state 3 is severe CAV. State 4 indicates death.
firstobs	(numeric)	0 = record represents an angiogram or date of death. 1 = record represents transplant (patient's first observation)
statemax	(numeric)	Maximum observed state so far for this patient (added in version 1.5.1)

**Source**

Papworth Hospital, U.K.

**References**

Sharples, L.D. and Jackson, C.H. and Parameshwar, J. and Wallwork, J. and Large, S.R. (2003). Diagnostic accuracy of coronary angiopathy and risk factors for post-heart-transplant cardiac allograft vasculopathy. *Transplantation* 76(4):679-82

---

cmodel.object

*Developer documentation: censoring model object*


---

**Description**

A list giving information about censored states, their labels in the data and what true states they represent.

**Value**

ncens	The number of distinct values used for censored observations in the state data supplied to <a href="#">msm</a> .
censor	A vector of length ncens, giving the labels used for censored states in the data.
states	A vector obtained by <code>unlist()</code> ing a list with ncens elements, each giving the set of true states that an observation with this label could be.
index	Index into states for the first state corresponding to each censor, plus an extra <code>length(states)+1</code> .

**See Also**

[msm.object](#).

---

coef.msm

*Extract model coefficients*


---

**Description**

Extract the estimated log transition intensities and the corresponding linear effects of each covariate.

**Usage**

```
## S3 method for class 'msm'
coef(object, ...)
```

**Arguments**

object	A fitted multi-state model object, as returned by <a href="#">msm</a> .
...	(unused) further arguments passed to or from other methods.

**Value**

If there is no misclassification, `coef.msm` returns a list of matrices. The first component, labelled `logbaseline`, is a matrix containing the estimated transition intensities on the log scale with any covariates fixed at their means in the data. Each remaining component is a matrix giving the linear effects of the labelled covariate on the matrix of log intensities.

For misclassification models, `coef.msm` returns a list of lists. The first component, `Qmatrices`, is a list of matrices as described in the previous paragraph. The additional component `Ematrices` is a list of similar format containing the logit-misclassification probabilities and any estimated covariate effects.

**Author(s)**

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**See Also**

[msm](#)

---

crudeinits.msm

*Calculate crude initial values for transition intensities*

---

**Description**

Calculates crude initial values for transition intensities by assuming that the data represent the exact transition times of the Markov process.

**Usage**

```
crudeinits.msm(formula, subject, qmatrix, data=NULL, censor=NULL, censor.states=NULL)
```

**Arguments**

formula	A formula giving the vectors containing the observed states and the corresponding observation times. For example, <code>state ~ time</code> Observed states should be in the set $1, \dots, n$ , where $n$ is the number of states. Note hidden Markov models are not supported by this function.
subject	Vector of subject identification numbers for the data specified by formula. If missing, then all observations are assumed to be on the same subject. These must be sorted so that all observations on the same subject are adjacent.
qmatrix	Matrix of indicators for the allowed transitions. An initial value will be estimated for each value of <code>qmatrix</code> that is greater than zero. Transitions are taken as disallowed for each entry of <code>qmatrix</code> that is 0.
data	An optional data frame in which the variables represented by <code>subject</code> and <code>state</code> can be found.

`censor` A state, or vector of states, which indicates censoring. See [msm](#).

`censor.states` Specifies the underlying states which censored observations can represent. See [msm](#).

### Details

Suppose we want a crude estimate of the transition intensity  $q_{rs}$  from state  $r$  to state  $s$ . If we observe  $n_{rs}$  transitions from state  $r$  to state  $s$ , and a total of  $n_r$  transitions from state  $r$ , then  $q_{rs}/q_{rr}$  can be estimated by  $n_{rs}/n_r$ . Then, given a total of  $T_r$  years spent in state  $r$ , the mean sojourn time  $1/q_{rr}$  can be estimated as  $T_r/n_r$ . Thus,  $n_{rs}/T_r$  is a crude estimate of  $q_{rs}$ .

If the data do represent the exact transition times of the Markov process, then these are the exact maximum likelihood estimates.

Observed transitions which are incompatible with the given `qmatrix` are ignored. Censored states are ignored.

### Value

The estimated transition intensity matrix. This can be used as the `qmatrix` argument to [msm](#).

### Author(s)

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### See Also

[statetable.msm](#)

### Examples

```
data(cav)
twoway4.q <- rbind(c(-0.5, 0.25, 0, 0.25), c(0.166, -0.498, 0.166, 0.166),
c(0, 0.25, -0.5, 0.25), c(0, 0, 0, 0))
statetable.msm(state, PTNUM, data=cav)
crudeinits.msm(state ~ years, PTNUM, data=cav, qmatrix=twoway4.q)
```

---

deltamethod

*The delta method*

---

### Description

Delta method for approximating the standard error of a transformation  $g(X)$  of a random variable  $X = (x_1, x_2, \dots)$ , given estimates of the mean and covariance matrix of  $X$ .

### Usage

```
deltamethod(g, mean, cov, ses=TRUE)
```

**Arguments**

g	A formula representing the transformation. The variables must be labelled $x_1, x_2, \dots$ . For example, $\sim 1 / (x_1 + x_2)$ If the transformation returns a vector, then a list of formulae representing $(g_1, g_2, \dots)$ can be provided, for example <code>list( ~ x1 + x2, ~ x1 / (x1 + x2) )</code>
mean	The estimated mean of $X$
cov	The estimated covariance matrix of $X$
ses	If TRUE, then the standard errors of $g_1(X), g_2(X), \dots$ are returned. Otherwise the covariance matrix of $g(X)$ is returned.

**Details**

The delta method expands a differentiable function of a random variable about its mean, usually with a first-order Taylor approximation, and then takes the variance. For example, an approximation to the covariance matrix of  $g(X)$  is given by

$$Cov(g(X)) = g'(\mu)Cov(X)[g'(\mu)]^T$$

where  $\mu$  is an estimate of the mean of  $X$ . This function uses symbolic differentiation via [deriv](#).

A limitation of this function is that variables created by the user are not visible within the formula `g`. To work around this, it is necessary to build the formula as a string, using functions such as `sprintf`, then to convert the string to a formula using `as.formula`. See the example below.

If you can spare the computational time, bootstrapping is a more accurate method of calculating confidence intervals or standard errors for transformations of parameters. See [boot.msm](#). Simulation from the asymptotic distribution of the MLEs (see e.g. Mandel 2013) is also a convenient alternative.

**Value**

A vector containing the standard errors of  $g_1(X), g_2(X), \dots$  or a matrix containing the covariance of  $g(X)$ .

**Author(s)**

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>

**References**

Oehlert, G. W. (1992) *A note on the delta method*. *American Statistician* 46(1).

Mandel, M. (2013) *Simulation based confidence intervals for functions with complicated derivatives*. *The American Statistician* 67(2):76-81.

## Examples

```
## Simple linear regression, E(y) = alpha + beta x
x <- 1:100
y <- rnorm(100, 4*x, 5)
toy.lm <- lm(y ~ x)
estmean <- coef(toy.lm)
estvar <- summary(toy.lm)$cov.unscaled * summary(toy.lm)$sigma^2

## Estimate of (1 / (alphahat + betahat))
1 / (estmean[1] + estmean[2])
## Approximate standard error
deltamethod (~ 1 / (x1 + x2), estmean, estvar)

## We have a variable z we would like to use within the formula.
z <- 1
## deltamethod (~ z / (x1 + x2), estmean, estvar) will not work.
## Instead, build up the formula as a string, and convert to a formula.
form <- sprintf("~ %f / (x1 + x2)", z)
form
deltamethod(as.formula(form), estmean, estvar)
```

---

draic.msm

*Criteria for comparing two multi-state models with nested state spaces*


---

## Description

A modification of Akaike's information criterion, and a leave-one-out likelihood cross-validation criterion, for comparing the predictive ability of two Markov multi-state models with nested state spaces. This is evaluated based on the restricted or aggregated data which the models have in common.

Note that standard AIC can be computed for one or more fitted `msm` models  $x, y, \dots$  using `AIC(x, y, \dots)`, and this can be used to compare models fitted to the same data. `draic.msm` and `dr1cv.msm` are designed for models fitted to data with differently-aggregated state spaces.

## Usage

```
draic.msm(msm.full, msm.coarse, likelihood.only=FALSE,
          information=c("expected", "observed"), tl=0.95)
dr1cv.msm(msm.full, msm.coarse, tl=0.95, cores=NULL,
          verbose=TRUE, outfile=NULL)
```

## Arguments

`msm.full`            Model on the bigger state space.

msm.coarse	<p>Model on the smaller state space.</p> <p>The two models must both be non-hidden Markov models without censored states.</p> <p>The two models must be fitted to the same datasets, except that the state space of the coarse model must be an aggregated version of the state space of the full model. That is, every state in the full dataset must correspond to a unique state in the coarse dataset. For example, for the full state variable <math>c(1, 1, 2, 2, 3, 4)</math>, the corresponding coarse states could be <math>c(1, 1, 2, 2, 3)</math>, but not <math>c(1, 2, 3, 4, 4, 4)</math>.</p> <p>The structure of allowed transitions in the coarse model must also be a collapsed version of the big model structure, but no check is currently made for this in the code.</p> <p>To use these functions, all objects which were used in the calls to fit <code>msm.full</code> and <code>msm.coarse</code> must be in the working environment, for example, datasets and definitions of transition matrices.</p>
likelihood.only	Don't calculate Hessians and trace term (DRAIC).
information	Use observed or expected information in the DRAIC trace term. Expected is the default, and much faster, though is only available for models fitted to pure panel data (all <code>obstype=1</code> in the call to <code>msm</code> , thus not exact transition times or exact death times)
t1	Width of symmetric tracking interval, by default 0.95 for a 95% interval.
cores	Number of processor cores to use in <code>dr1cv</code> for cross-validation by parallel processing. Requires the <b>doParallel</b> package to be installed. If not specified, parallel processing is not used. If <code>cores</code> is set to the string "default", the default methods of <code>makeCluster</code> (on Windows) or <code>registerDoParallel</code> (on Unix-like) are used.
verbose	Print intermediate results of each iteration of cross-validation to the console while running. May not work with parallel processing.
outfile	Output file to print intermediate results of cross-validation. Useful to track execution speed when using parallel processing, where output to the console may not work.

## Details

The difference in restricted AIC (Liquet and Commenges, 2011), as computed by this function, is defined as

$$D_{RAIC} = l(\gamma_n | \mathbf{x}'') - l(\theta_n | \mathbf{x}'') + \text{trace}(J(\theta_n | \mathbf{x}'') J(\theta_n | \mathbf{x})^{-1} - J(\gamma_n | \mathbf{x}'') J(\gamma_n | \mathbf{x}')^{-1})$$

where  $\gamma$  and  $\theta$  are the maximum likelihood estimates of the smaller and bigger models, fitted to the smaller and bigger data, respectively.

$l(\gamma_n | \mathbf{x}'')$  represents the likelihood of the simpler model evaluated on the restricted data.

$l(\theta_n | \mathbf{x}'')$  represents the likelihood of the complex model evaluated on the restricted data. This is a hidden Markov model, with a misclassification matrix and initial state occupancy probabilities as described by Thom et al (2014).



$J()$  are the corresponding (expected or observed, as specified by the user) information matrices.

$\mathbf{x}$  is the expanded data, to which the bigger model was originally fitted, and  $\mathbf{x}'$  is the data to which the smaller model was originally fitted.  $\mathbf{x}''$  is the restricted data which the two models have in common.  $\mathbf{x}'' = \mathbf{x}'$  in this implementation, so the models are nested.

The difference in likelihood cross-validatory criteria (Liquet and Commenges, 2011) is defined as

$$D_{RLCV} = 1/n \sum_{i=1}^n \log(h_{X''}(x''_i|\gamma_{-i})/g_{X''}(x''_i|\theta_{-i}))$$

where  $\gamma_{-i}$  and  $\theta_{-i}$  are the maximum likelihood estimates from the smaller and bigger models fitted to datasets with subject  $i$  left out,  $g()$  and  $h()$  are the densities of the corresponding models, and  $x''_i$  is the restricted data from subject  $i$ .

Tracking intervals are analogous to confidence intervals, but not strictly the same, since the quantity which D\_RAIC aims to estimate, the difference in expected Kullback-Leibler discrepancy for predicting a replicate dataset, depends on the sample size. See the references.

Positive values for these criteria indicate the coarse model is preferred, while negative values indicate the full model is preferred.

### Value

A list containing  $D_{RAIC}$  (`draic.msm`) or  $D_{RLCV}$  (`dr1cv.msm`), its component terms, and tracking intervals.

### Author(s)

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>, H. H. Z. Thom <[howard.thom@bristol.ac.uk](mailto:howard.thom@bristol.ac.uk)>

### References

Thom, H. and Jackson, C. and Commenges, D. and Sharples, L. (2015) State selection in multistate models with application to quality of life in psoriatic arthritis. *Statistics In Medicine* 34(16) 2381 - 2480.

Liquet, B. and Commenges D. (2011) Choice of estimators based on different observations: Modified AIC and LCV criteria. *Scandinavian Journal of Statistics*; 38:268-287.

### See Also

[logLik.msm](#)

---

ecmodel.object	<i>Developer documentation: model for covariates on misclassification probabilities</i>
----------------	---

---

### Description

A list representing the model for covariates on misclassification probabilities.

### Value

npars	Number of covariate effect parameters. This is defined as the number of covariates on misclassification (with factors expanded as contrasts) multiplied by the number of allowed misclassifications in the model.
ndpars	Number of distinct covariate effect parameters, as npars, but after any equality constraints have been applied.
ncovs	Number of covariates on misclassification, with factors expanded as contrasts.
constr	List of equality constraints on these covariate effects, as supplied in the <code>misconstraint</code> argument to <code>msm</code> .
covlabels	Names / labels of these covariates in the model matrix (see <code>model.matrix.msm</code> ).
inits	Initial values for these covariate effects, as a vector formed from the <code>misccovinits</code> list supplied to <code>msm</code> .
covmeans	Means of these covariates in the data (excluding data not required to fit the model, such as observations with missing data in other elements or subjects' last observations). This includes means of 0/1 factor contrasts as well as continuous covariates (for historic reasons, which may not be sensible).

### See Also

[msm.object](#).

---

efpt.msm	<i>Expected first passage time</i>
----------	------------------------------------

---

### Description

Expected time until first reaching a particular state or set of states in a Markov model.

### Usage

```
efpt.msm(x=NULL, qmatrix=NULL, tostate, start="all", covariates="mean",
         ci=c("none", "normal", "bootstrap"), cl=0.95, B=1000,
         cores=NULL, ...)
```

**Arguments**

<code>x</code>	A fitted multi-state model, as returned by <a href="#">msm</a> .
<code>qmatrix</code>	Instead of <code>x</code> , you can simply supply a transition intensity matrix in <code>qmatrix</code> .
<code>tostate</code>	State, or set of states supplied as a vector, for which to estimate the first passage time into. Can be integer, or character matched to the row names of the <code>Q</code> matrix.
<code>start</code>	Starting state (integer). By default ( <code>start="all"</code> ), this will return a vector of expected passage times from each state in turn.  Alternatively, this can be used to obtain the expected first passage time from a <i>set</i> of states, rather than single states. To achieve this, <code>state</code> is set to a vector of weights, with length equal to the number of states in the model. These weights should be proportional to the probability of starting in each of the states in the desired set, so that weights of zero are supplied for other states. The function will calculate the weighted average of the expected passage times from each of the corresponding states.
<code>covariates</code>	Covariate values defining the intensity matrix for the fitted model <code>x</code> , as supplied to <a href="#">qmatrix.msm</a> .
<code>ci</code>	If "normal", then calculate a confidence interval by simulating <code>B</code> random vectors from the asymptotic multivariate normal distribution implied by the maximum likelihood estimates (and covariance matrix) of the log transition intensities and covariate effects.  If "bootstrap" then calculate a confidence interval by non-parametric bootstrap refitting. This is 1-2 orders of magnitude slower than the "normal" method, but is expected to be more accurate. See <a href="#">boot.msm</a> for more details of bootstrapping in <a href="#">msm</a> .  If "none" (the default) then no confidence interval is calculated.
<code>c1</code>	Width of the symmetric confidence interval, relative to 1.
<code>B</code>	Number of bootstrap replicates.
<code>cores</code>	Number of cores to use for bootstrapping using parallel processing. See <a href="#">boot.msm</a> for more details.
<code>...</code>	Arguments to pass to <a href="#">MatrixExp</a> .

**Details**

The expected first passage times from each of a set of states  $\mathbf{i}$  to the remaining set of states  $\bar{\mathbf{i}}$  in the state space, for a model with transition intensity matrix  $Q$ , are

$$-Q_{\mathbf{i},\mathbf{i}}^{-1} \mathbf{1}$$

where  $\mathbf{1}$  is a vector of ones, and  $Q_{\mathbf{i},\mathbf{i}}$  is the square subset of  $Q$  pertaining to states  $\mathbf{i}$ .

It is equal to the sum of mean sojourn times for all states between the "from" and "to" states in a unidirectional model. If there is non-zero chance of reaching an absorbing state before reaching `tostate`, then it is infinite. It is trivially zero if the "from" state equals `tostate`.

This function currently only handles time-homogeneous Markov models. For time-inhomogeneous models it will assume that  $Q$  equals the average intensity matrix over all times and observed covariates. Simulation might be used to handle time dependence.

Note this is the *expectation* of first passage time, and the confidence intervals are CIs for this mean, not predictive intervals for the first passage time. The full distribution of the first passage time to a set of states can be obtained by setting the rows of the intensity matrix  $Q$  corresponding to that set of states to zero to make a model where those states are absorbing. The corresponding transition probability matrix  $Exp(Qt)$  then gives the probabilities of having hit or passed that state by a time  $t$  (see the example below). This is implemented in [ppass.msm](#).

### Value

A vector of expected first passage times, or "hitting times", from each state to the desired state.

### Author(s)

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>

### References

Norris, J. R. (1997) Markov Chains. Cambridge University Press.

### See Also

[sojourn.msm](#), [totlos.msm](#), [boot.msm](#).

### Examples

```

twoway4.q <- rbind(c(-0.5, 0.25, 0, 0.25), c(0.166, -0.498, 0.166, 0.166),
                 c(0, 0.25, -0.5, 0.25), c(0, 0, 0, 0))
efpt.msm(qmatrix=twoway4.q, tostate=3)
# given in state 1, expected time to reaching state 3 is infinite
# since may die (state 4) before entering state 3

# If we remove the death state from the model, EFPTs become finite
Q <- twoway4.q[1:3,1:3]; diag(Q) <- 0; diag(Q) <- -rowSums(Q)
efpt.msm(qmatrix=Q, tostate=3)

# Suppose we cannot die or regress while in state 2, can only go to state 3
Q <- twoway4.q; Q[2,4] <- Q[2,1] <- 0; diag(Q) <- 0; diag(Q) <- -rowSums(Q)
efpt.msm(qmatrix=Q, tostate=3)
# The expected time from 2 to 3 now equals the mean sojourn time in 2.
-1/Q[2,2]

# Calculate cumulative distribution of the first passage time
# into state 3 for the following three-state model
Q <- twoway4.q[1:3,1:3]; diag(Q) <- 0; diag(Q) <- -rowSums(Q)
# Firstly form a model where the desired hitting state is absorbing
Q[3,] <- 0
MatrixExp(Q, t=10)[,3]
ppass.msm(qmatrix=Q, tot=10)
# Given in state 1 at time 0, P(hit 3 by time 10) = 0.479
MatrixExp(Q, t=50)[,3] # P(hit 3 by time 50) = 0.98
ppass.msm(qmatrix=Q, tot=50)

```

ematrix.msm

*Misclassification probability matrix***Description**

Extract the estimated misclassification probability matrix, and corresponding confidence intervals, from a fitted multi-state model at a given set of covariate values.

**Usage**

```
ematrix.msm(x, covariates="mean", ci=c("delta","normal","bootstrap","none"),
           cl=0.95, B=1000, cores=NULL)
```

**Arguments**

x	A fitted multi-state model, as returned by <a href="#">msm</a> .
covariates	The covariate values for which to estimate the misclassification probability matrix. This can either be: <ul style="list-style-type: none"> <li>the string "mean", denoting the means of the covariates in the data (this is the default),</li> <li>the number 0, indicating that all the covariates should be set to zero,</li> <li>or a list of values, with optional names. For example               <pre>list(60,1)</pre>               where the order of the list follows the order of the covariates originally given in the model formula, or a named list,               <pre>list(age = 60, sex = 1)</pre> </li> </ul>
ci	If "delta" (the default) then confidence intervals are calculated by the delta method, or by simple transformation of the Hessian in the very simplest cases. If "normal", then calculate a confidence interval by simulating B random vectors from the asymptotic multivariate normal distribution implied by the maximum likelihood estimates (and covariance matrix) of the multinomial-logit-transformed misclassification probabilities and covariate effects, then transforming back. If "bootstrap" then calculate a confidence interval by non-parametric bootstrap refitting. This is 1-2 orders of magnitude slower than the "normal" method, but is expected to be more accurate. See <a href="#">boot.msm</a> for more details of bootstrapping in <a href="#">msm</a> .
cl	Width of the symmetric confidence interval to present. Defaults to 0.95.
B	Number of bootstrap replicates, or number of normal simulations from the distribution of the MLEs
cores	Number of cores to use for bootstrapping using parallel processing. See <a href="#">boot.msm</a> for more details.

**Details**

Misclassification probabilities and covariate effects are estimated on the multinomial-logit scale by [msm](#). A covariance matrix is estimated from the Hessian of the maximised log-likelihood. From these, the delta method can be used to obtain standard errors of the probabilities on the natural scale at arbitrary covariate values. Confidence intervals are estimated by assuming normality on the multinomial-logit scale.

**Value**

A list with components:

estimate	Estimated misclassification probability matrix. The rows correspond to true states, and columns observed states.
SE	Corresponding approximate standard errors.
L	Lower confidence limits.
U	Upper confidence limits.

Or if `ci="none"`, then `ematrix.msm` just returns the estimated misclassification probability matrix.

The default print method for objects returned by `ematrix.msm` presents estimates and confidence limits. To present estimates and standard errors, do something like

```
ematrix.msm(x)[c("estimates", "SE")]
```

**Author(s)**

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>

**See Also**

[qmatrix.msm](#)

---

emodel.object

*Developer documentation: misclassification model structure object*

---

**Description**

A list giving information about the misclassifications assumed in a multi-state model fitted with the `ematrix` argument of `msm`. Returned in a fitted `msm` model object. This information is converted internally to a `hmodel` object (see [hmodel.object](#)) for use in likelihood computations.

**Value**

<code>nstates</code>	Number of states (same as <code>qmodel\$nstates</code> ).
<code>npars</code>	Number of allowed misclassifications, equal to <code>sum(imatrix)</code> .
<code>imatrix</code>	Indicator matrix for allowed misclassifications. This has $(r, s)$ entry 1 if misclassification of true state $r$ as observed state $s$ is possible. diagonal entries are arbitrarily set to 0.
<code>ematrix</code>	Matrix of initial values for the misclassification probabilities, supplied as the <code>ematrix</code> argument of <code>msm</code> .
<code>inits</code>	Vector of these initial values, reading across rows of <code>qmatrix</code> and excluding the diagonal and disallowed transitions.
<code>constr</code>	Indicators for equality constraints on baseline misclassification probabilities, taken from the <code>econstraint</code> argument to <code>msm</code> , and mapped if necessary to the set (1,2,3,...)
<code>ndpars</code>	Number of distinct misclassification probabilities, after applying equality constraints.
<code>nipars</code>	Number of initial state occupancy probabilities being estimated. This is zero if <code>est.iniprbs=FALSE</code> , otherwise equal to the number of states.
<code>iniprbs</code>	Initial state occupancy probabilities, as supplied to <code>msm</code> (initial values before estimation, if <code>est.iniprbs=TRUE</code> .)
<code>est.iniprbs</code>	Are initial state occupancy probabilities estimated (TRUE or FALSE), as supplied in the <code>est.iniprbs</code> argument of <code>msm</code> .

**See Also**

[msm.object](#), [qmodel.object](#), [hmodel.object](#).

---

fev

*FEV1 measurements from lung transplant recipients*


---

**Description**

A series of measurements of the forced expiratory volume in one second (FEV1) from lung transplant recipients, from six months onwards after their transplant.

**Usage**

```
fev
```

**Format**

A data frame containing 5896 rows. There are 204 patients, the rows are grouped by patient number and ordered by days after transplant. Each row represents an examination and containing an additional covariate.

ptnum	(numeric)	Patient identification number.
days	(numeric)	Examination time (days after transplant).
fev	(numeric)	Percentage of baseline FEV1. A code of 999 indicates the patient's date of death.
acute	(numeric)	0/1 indicator for whether the patient suffered an acute infection or rejection within 14 days of the visit.

### Details

A baseline "normal" FEV1 for each individual is calculated using measurements from the first six months after transplant. After six months, as presented in this dataset, FEV1 is expressed as a percentage of the baseline value.

FEV1 is monitored to diagnose bronchiolitis obliterans syndrome (BOS), a long-term lung function decline, thought to be a form of chronic rejection. Acute rejections and infections also affect the lung function in the short term.

### Source

Papworth Hospital, U.K.

### References

Jackson, C.H. and Sharples, L.D. Hidden Markov models for the onset and progression of bronchiolitis obliterans syndrome in lung transplant recipients *Statistics in Medicine*, 21(1): 113–128 (2002).

---

hazard.msm	<i>Calculate tables of hazard ratios for covariates on transition intensities</i>
------------	---

---

### Description

Hazard ratios are computed by exponentiating the estimated covariate effects on the log-transition intensities. This function is called by [summary.msm](#).

### Usage

```
hazard.msm(x, hazard.scale = 1, cl = 0.95)
```

### Arguments

x	Output from <a href="#">msm</a> representing a fitted multi-state model.
hazard.scale	Vector with same elements as number of covariates on transition rates. Corresponds to the increase in each covariate used to calculate its hazard ratio. Defaults to all 1.
cl	Width of the symmetric confidence interval to present. Defaults to 0.95.



**Value**

A list of tables containing hazard ratio estimates, one table for each covariate. Each table has three columns, containing the hazard ratio, and an approximate upper and lower confidence limit respectively (assuming normality on the log scale), for each Markov chain transition intensity.

**Author(s)**

C. H. Jackson <chris.jackson@mrc-bsu.cam.ac.uk>

**See Also**

[msm](#), [summary.msm](#), [odds.msm](#)

---

hmm-dists

*Hidden Markov model constructors*


---

**Description**

These functions are used to specify the distribution of the response conditionally on the underlying state in a hidden Markov model. A list of these function calls, with one component for each state, should be used for the `hmodel` argument to `msm`. The initial values for the parameters of the distribution should be given as arguments. Note the initial values should be supplied as literal values - supplying them as variables is currently not supported.

**Usage**

```

hmmCat(prob, basecat)
hmmIdent(x)
hmmUnif(lower, upper)
hmmNorm(mean, sd)
hmmLNorm(meanlog, sdlog)
hmmExp(rate)
hmmGamma(shape, rate)
hmmWeibull(shape, scale)
hmmPois(rate)
hmmBinom(size, prob)
hmmTNorm(mean, sd, lower, upper)
hmmMETNorm(mean, sd, lower, upper, sderr, meanerr=0)
hmmMEUnif(lower, upper, sderr, meanerr=0)
hmmNBinom(dis, prob)
hmmBetaBinom(size, meanp, sdp)
hmmBeta(shape1, shape2)
hmmT(mean, scale, df)

```

**Arguments**

prob	(hmmCat) Vector of probabilities of observing category 1, 2, . . . , length(prob) respectively. Or the probability governing a binomial or negative binomial distribution.
basecat	(hmmCat) Category which is considered to be the "baseline", so that during estimation, the probabilities are parameterised as probabilities relative to this baseline category. By default, the category with the greatest probability is used as the baseline.
x	(hmmIdent) Code in the data which denotes the exactly-observed state.
mean	(hmmNorm, hmmLNorm, hmmTNorm) Mean defining a Normal, or truncated Normal distribution.
sd	(hmmNorm, hmmLNorm, hmmTNorm) Standard deviation defining a Normal, or truncated Normal distribution.
meanlog	(hmmNorm, hmmLNorm, hmmTNorm) Mean on the log scale, for a log Normal distribution.
sdlog	(hmmNorm, hmmLNorm, hmmTNorm) Standard deviation on the log scale, for a log Normal distribution.
rate	(hmmPois, hmmExp, hmmGamma) Rate of a Poisson, Exponential or Gamma distribution (see <a href="#">dpois</a> , <a href="#">dexp</a> , <a href="#">dgamma</a> ).
shape	(hmmPois, hmmExp, hmmGamma) Shape parameter of a Gamma or Weibull distribution (see <a href="#">dgamma</a> , <a href="#">dweibull</a> ).
shape1, shape2	First and second parameters of a beta distribution (see <a href="#">dbeta</a> ).
scale	(hmmGamma) Scale parameter of a Gamma distribution (see <a href="#">dgamma</a> ), or unstandardised Student t distribution.
df	Degrees of freedom of the Student t distribution.
size	Order of a Binomial distribution (see <a href="#">dbinom</a> ).
disp	Dispersion parameter of a negative binomial distribution, also called size or order. (see <a href="#">dnbinom</a> ).
meanp	Mean outcome probability in a beta-binomial distribution
sdp	Standard deviation describing the overdispersion of the outcome probability in a beta-binomial distribution
lower	(hmmUnif, hmmTNorm, hmmMEUnif) Lower limit for an Uniform or truncated Normal distribution.
upper	(hmmUnif, hmmTNorm, hmmMEUnif) Upper limit for an Uniform or truncated Normal distribution.
sderr	(hmmMETNorm, hmmUnif) Standard deviation of the Normal measurement error distribution.
meanerr	(hmmMETNorm, hmmUnif) Additional shift in the measurement error, fixed to 0 by default. This may be modelled in terms of covariates.

## Details

`hmmCat` represents a categorical response distribution on the set  $1, 2, \dots, \text{length}(\text{prob})$ . The Markov model with misclassification is an example of this type of model. The categories in this case are (some subset of) the underlying states.

The `hmmIdent` distribution is used for underlying states which are observed exactly without error. For hidden Markov models with multiple outcomes, (see `hmmMV`), the outcome in the data which takes the special `hmmIdent` value must be the first of the multiple outcomes.

`hmmUnif`, `hmmNorm`, `hmmLNorm`, `hmmExp`, `hmmGamma`, `hmmWeibull`, `hmmPois`, `hmmBinom`, `hmmTNorm`, `hmmNBinom` and `hmmBeta` represent Uniform, Normal, log-Normal, exponential, Gamma, Weibull, Poisson, Binomial, truncated Normal, negative binomial and beta distributions, respectively, with parameterisations the same as the default parameterisations in the corresponding base R distribution functions.

`hmmT` is the Student t distribution with general mean  $\mu$ , scale  $\sigma$  and degrees of freedom  $df$ . The variance is  $\sigma^2 df / (df + 2)$ . Note the t distribution in base R `dt` is a standardised one with mean 0 and scale 1. These allow any positive (integer or non-integer)  $df$ . By default, all three parameters, including  $df$ , are estimated when fitting a hidden Markov model, but in practice,  $df$  might need to be fixed for identifiability - this can be done using the `fixedpars` argument to `msm`.

The `hmmMETNorm` and `hmmMEUnif` distributions are truncated Normal and Uniform distributions, but with additional Normal measurement error on the response. These are generalisations of the distributions proposed by Satten and Longini (1996) for modelling the progression of CD4 cell counts in monitoring HIV disease. See `medists` for density, distribution, quantile and random generation functions for these distributions. See also `tnorm` for density, distribution, quantile and random generation functions for the truncated Normal distribution.

See the PDF manual ‘`msm-manual.pdf`’ in the ‘`doc`’ subdirectory for algebraic definitions of all these distributions. New hidden Markov model response distributions can be added to `msm` by following the instructions in Section 2.17.1.

Parameters which can be modelled in terms of covariates, on the scale of a link function, are as follows.

PARAMETER NAME	LINK FUNCTION
<code>mean</code>	identity
<code>meanlog</code>	identity
<code>rate</code>	log
<code>scale</code>	log
<code>meanerr</code>	identity
<code>meanp</code>	logit
<code>prob</code>	logit or multinomial logit

Parameters `basecat`, `lower`, `upper`, `size`, `meanerr` are fixed at their initial values. All other parameters are estimated while fitting the hidden Markov model, unless the appropriate `fixedpars` argument is supplied to `msm`.

For categorical response distributions (`hmmCat`) the outcome probabilities initialized to zero are fixed at zero, and the probability corresponding to `basecat` is fixed to one minus the sum of the remaining probabilities. These remaining probabilities are estimated, and can be modelled in terms of covariates via multinomial logistic regression (relative to `basecat`).

**Value**

Each function returns an object of class `hmodel`, which is a list containing information about the model. The only component which may be useful to end users is `r`, a function of one argument `n` which returns a random sample of size `n` from the given distribution.

**Author(s)**

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>

**References**

Satten, G.A. and Longini, I.M. Markov chains with measurement error: estimating the 'true' course of a marker of the progression of human immunodeficiency virus disease (with discussion) *Applied Statistics* 45(3): 275-309 (1996).

Jackson, C.H. and Sharples, L.D. Hidden Markov models for the onset and progression of bronchiolitis obliterans syndrome in lung transplant recipients *Statistics in Medicine*, 21(1): 113-128 (2002).

Jackson, C.H., Sharples, L.D., Thompson, S.G. and Duffy, S.W. and Couto, E. Multi-state Markov models for disease progression with classification error. *The Statistician*, 52(2): 193-209 (2003).

**See Also**

[msm](#)

---

hmmMV

*Multivariate hidden Markov models*

---

**Description**

Constructor for a a multivariate hidden Markov model (HMM) where each of the `n` variables observed at the same time has a (potentially different) standard univariate distribution conditionally on the underlying state. The `n` outcomes are independent conditionally on the hidden state.

If a particular state in a HMM has such an outcome distribution, then a call to `hmmMV` is supplied as the corresponding element of the `hmodel` argument to `msm`. See Example 2 below.

A multivariate HMM where multiple outcomes at the same time are generated from the *same* distribution is specified in the same way as the corresponding univariate model, so that `hmmMV` is not required. The outcome data are simply supplied as a matrix instead of a vector. See Example 1 below.

The outcome data for such models are supplied as a matrix, with number of columns equal to the maximum number of arguments supplied to the `hmmMV` calls for each state. If some but not all of the variables are missing (NA) at a particular time, then the observed data at that time still contribute to the likelihood. The missing data are assumed to be missing at random. The Viterbi algorithm may be used to predict the missing values given the fitted model and the observed data.

Typically the outcome model for each state will be from the same family or set of families, but with different parameters. Theoretically, different numbers of distributions may be supplied for different

states. If a particular state has fewer outcomes than the maximum, then the data for that state are taken from the first columns of the response data matrix. However this is not likely to be a useful model, since the number of observations will probably give information about the underlying state, violating the missing at random assumption.

Models with outcomes that are dependent conditionally on the hidden state (e.g. correlated multivariate normal observations) are not currently supported.

## Usage

```
hmmMV(...)
```

## Arguments

... The number of arguments supplied should equal the maximum number of observations made at one time. Each argument represents the univariate distribution of that outcome conditionally on the hidden state, and should be the result of calling a univariate hidden Markov model constructor (see [hmm-dists](#)).

## Value

A list of objects, each of class `hmmdist` as returned by the univariate HMM constructors documented in [hmm-dists](#). The whole list has class `hmmMVdist`, which inherits from `hmmdist`.

## Author(s)

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>

## References

Jackson, C. H., Su, L., Gladman, D. D. and Farewell, V. T. (2015) On modelling minimal disease activity. *Arthritis Care and Research* (early view).

## See Also

[hmm-dists](#), [msm](#)

## Examples

```
## Simulate data from a Markov model
nsubj <- 30; nobspt <- 5
sim.df <- data.frame(subject = rep(1:nsubj, each=nobspt),
                     time = seq(0, 20, length=nobspt))

set.seed(1)
two.q <- rbind(c(-0.1, 0.1), c(0, 0))
dat <- simmulti.msm(sim.df[,1:2], qmatrix=two.q, drop.absorb=FALSE)

### EXAMPLE 1
## Generate two observations at each time from the same outcome
## distribution:
## Bin(40, 0.1) for state 1, Bin(40, 0.5) for state 2
dat$obs1[dat$state==1] <- rbinom(sum(dat$state==1), 40, 0.1)
```

```

dat$obs2[dat$state==1] <- rbinom(sum(dat$state==1), 40, 0.1)
dat$obs1[dat$state==2] <- rbinom(sum(dat$state==2), 40, 0.5)
dat$obs2[dat$state==2] <- rbinom(sum(dat$state==2), 40, 0.5)
dat$obs <- cbind(obs1 = dat$obs1, obs2 = dat$obs2)

## Fitted model should approximately recover true parameters
msm(obs ~ time, subject=subject, data=dat, qmatrix=two.q,
     hmodel = list(hmmBinom(size=40, prob=0.2),
                   hmmBinom(size=40, prob=0.2)))

### EXAMPLE 2
## Generate two observations at each time from different
## outcome distributions:
## Bin(40, 0.1) and Bin(40, 0.2) for state 1,
dat$obs1 <- dat$obs2 <- NA
dat$obs1[dat$state==1] <- rbinom(sum(dat$state==1), 40, 0.1)
dat$obs2[dat$state==1] <- rbinom(sum(dat$state==1), 40, 0.2)

## Bin(40, 0.5) and Bin(40, 0.6) for state 2
dat$obs1[dat$state==2] <- rbinom(sum(dat$state==2), 40, 0.6)
dat$obs2[dat$state==2] <- rbinom(sum(dat$state==2), 40, 0.5)
dat$obs <- cbind(obs1 = dat$obs1, obs2 = dat$obs2)

## Fitted model should approximately recover true parameters
msm(obs ~ time, subject=subject, data=dat, qmatrix=two.q,
     hmodel = list(hmmMV(hmmBinom(size=40, prob=0.3),
                         hmmBinom(size=40, prob=0.3)),
                   hmmMV(hmmBinom(size=40, prob=0.3),
                         hmmBinom(size=40, prob=0.3))),
     control=list(maxit=10000))

```

---

hmodel.object

*Developer documentation: hidden Markov model structure object*


---

## Description

A list giving information about the models for the outcome data conditionally on the states of a hidden Markov model. Used in internal computations, and returned in a fitted [msm](#) model object.

## Value

hidden	TRUE for hidden Markov models, FALSE otherwise.
nstates	Number of states, the same as <code>qmodel\$nstates</code> .
fitted	TRUE if the parameter values in <code>pars</code> are the maximum likelihood estimates, FALSE if they are the initial values.
models	The outcome distribution for each hidden state. A vector of length <code>nstates</code> whose $r$ th entry is the index of the state $r$ outcome distributions in the vector of supported distributions. The vector of supported distributions is given in full by <code>msm:::msm.HMODELS</code> : the first few are 1 for categorical outcome, 2 for identity, 3 for uniform and 4 for normal.

labels	String identifying each distribution in models.
npars	Vector of length <code>nstates</code> giving the number of parameters in each outcome distribution, excluding covariate effects.
nipars	Number of initial state occupancy probabilities being estimated. This is zero if <code>est.initprobs=FALSE</code> , otherwise equal to the number of states.
totpars	Total number of parameters, equal to <code>sum(npars)</code> .
pars	A vector of length <code>totpars</code> , made from concatenating a list of length <code>nstates</code> whose $r$ th component is vector of the parameters for the state $r$ outcome distribution.
plabs	List with the names of the parameters in <code>pars</code> .
parstate	A vector of length <code>totpars</code> , whose $i$ th element is the state corresponding to the $i$ th parameter.
firstpar	A vector of length <code>nstates</code> giving the index in <code>pars</code> of the first parameter for each state.
locpars	Index in <code>pars</code> of parameters which can have covariates on them.
initprobs	Initial state occupancy probabilities, as supplied to <code>msm</code> (initial values before estimation, if <code>est.initprobs=TRUE</code> .)
est.initprobs	Are initial state occupancy probabilities estimated (TRUE or FALSE), as supplied in the <code>est.initprobs</code> argument of <code>msm</code> .
ncovs	Number of covariate effects per parameter in <code>pars</code> , with, e.g. factor contrasts expanded.
coveffect	Vector of covariate effects, of length <code>sum(ncovs)</code> .
covlabels	Labels of these effects.
coveffstate	Vector indicating state corresponding to each element of <code>coveffect</code> .
ncoveffs	Number of covariate effects on HMM outcomes, equal to <code>sum(ncovs)</code> .
nicovs	Vector of length <code>nstates-1</code> giving the number of covariate effects on each initial state occupancy probability (log relative to the baseline probability).
icoveffect	Vector of length <code>sum(nicovs)</code> giving covariate effects on initial state occupancy probabilities.
nicoveffs	Number of covariate effects on initial state occupancy probabilities, equal to <code>sum(nicovs)</code> .
constr	Constraints on (baseline) hidden Markov model outcome parameters, as supplied in the <code>hconstraint</code> argument of <code>msm</code> , excluding covariate effects, converted to a vector and mapped to the set 1,2,3,... if necessary.
covconstr	Vector of constraints on covariate effects in hidden Markov outcome models, as supplied in the <code>hconstraint</code> argument of <code>msm</code> , excluding baseline parameters, converted to a vector and mapped to the set 1,2,3,... if necessary.
ranges	Matrix of range restrictions for HMM parameters, including those given to the <code>hranges</code> argument to <code>msm</code> .
foundse	TRUE if standard errors are available for the estimates.
initpmat	Matrix of initial state occupancy probabilities with one row for each subject (estimated if <code>est.initprobs=TRUE</code> ).
ci	Confidence intervals for baseline HMM outcome parameters.
covci	Confidence intervals for covariate effects in HMM outcome models.

**See Also**

[msm.object](#), [qmodel.object](#), [emodel.object](#).

---

logLik.msm

*Extract model log-likelihood*

---

**Description**

Extract the log-likelihood and the number of parameters of a model fitted with [msm](#).

**Usage**

```
## S3 method for class 'msm'  
logLik(object, by.subject=FALSE, ...)
```

**Arguments**

object	A fitted multi-state model object, as returned by <a href="#">msm</a> .
by.subject	Return vector of subject-specific log-likelihoods, which should sum to the total log-likelihood.
...	(unused) further arguments passed to or from other methods.

**Value**

The log-likelihood of the model represented by 'object' evaluated at the maximum likelihood estimates.

Akaike's information criterion can also be computed using [AIC\(object\)](#).

**Author(s)**

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>

**See Also**

[msm](#), [lrtest.msm](#).



---

lrtest.msm	<i>Likelihood ratio test</i>
------------	------------------------------

---

**Description**

Likelihood ratio test between two or more fitted multi-state models

**Usage**

```
lrtest.msm(...)
```

**Arguments**

... Two or more fitted multi-state models, as returned by [msm](#), ordered by increasing numbers of parameters.

**Value**

A matrix with three columns, giving the likelihood ratio statistic, difference in degrees of freedom and the chi-squared p-value for a comparison of the first model supplied with each subsequent model.

**Warning**

The comparison between models will only be valid if they are fitted to the same dataset. This may be a problem if there are missing values and R's default of 'na.action = na.omit' is used.

The likelihood ratio statistic only has the indicated chi-squared distribution if the models are nested. An alternative for comparing non-nested models is Akaike's information criterion. This can be computed for one or more fitted msm models  $x, y, \dots$  using [AIC\(x, y, \dots\)](#).

**See Also**

[logLik.msm,msm](#)

---

MatrixExp	<i>Matrix exponential</i>
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---

**Description**

Calculates the exponential of a square matrix.

**Usage**

```
MatrixExp(mat, t = 1, method=NULL, ...)
```

**Arguments**

mat	A square matrix
t	An optional scaling factor for mat.
method	Under the default of NULL, this simply wraps the <code>expm</code> function from the <b>expm</b> package. This is recommended. Options to <code>expm</code> can be supplied to <code>MatrixExp</code> , including <code>method</code> .  Otherwise, for backwards compatibility, the following options, which use code in the <b>msm</b> package, are available: "pade" for a Pade approximation method, "series" for the power series approximation, or "analytic" for the analytic formulae for simpler Markov model intensity matrices (see below). These options are only used if <code>mat</code> has repeated eigenvalues, thus the usual eigen-decomposition method cannot be used.
...	Arguments to pass to <code>expm</code> .

**Details**

See the `expm` documentation for details of the algorithms it uses.

Generally the exponential  $E$  of a square matrix  $M$  can often be calculated as

$$E = U \exp(D) U^{-1}$$

where  $D$  is a diagonal matrix with the eigenvalues of  $M$  on the diagonal,  $\exp(D)$  is a diagonal matrix with the exponentiated eigenvalues of  $M$  on the diagonal, and  $U$  is a matrix whose columns are the eigenvectors of  $M$ .

This method of calculation is used if "pade" or "series" is supplied but  $M$  has distinct eigenvalues. If  $M$  has repeated eigenvalues, then its eigenvector matrix may be non-invertible. In this case, the matrix exponential is calculated using the Pade approximation defined by Moler and van Loan (2003), or the less robust power series approximation,

$$\exp(M) = I + M + M^2/2 + M^3/3! + M^4/4! + \dots$$

For a continuous-time homogeneous Markov process with transition intensity matrix  $Q$ , the probability of occupying state  $s$  at time  $u + t$  conditional on occupying state  $r$  at time  $u$  is given by the  $(r, s)$  entry of the matrix  $\exp(tQ)$ .

If `mat` is a valid transition intensity matrix for a continuous-time Markov model (i.e. diagonal entries non-positive, off-diagonal entries non-negative, rows sum to zero), then for certain simpler model structures, there are analytic formulae for the individual entries of the exponential of `mat`. These structures are listed in the PDF manual and the formulae are coded in the **msm** source file `src/analyticp.c`. These formulae are only used if `method="analytic"`. This is more efficient, but it is not the default in `MatrixExp` because the code is not robust to extreme values. However it is the default when calculating likelihoods for models fitted by `msm`.

The implementation of the Pade approximation used by `method="pade"` was taken from JAGS by Martyn Plummer (<http://mcmc-jags.sourceforge.net>).

**Value**

The exponentiated matrix  $\exp(mat)$ . Or, if  $t$  is a vector of length 2 or more, an array of exponentiated matrices.

**References**

Cox, D. R. and Miller, H. D. *The theory of stochastic processes*, Chapman and Hall, London (1965)  
 Moler, C and van Loan, C (2003). Nineteen dubious ways to compute the exponential of a matrix, twenty-five years later. *SIAM Review* **45**, 3–49.

---

 medists

*Measurement error distributions*


---

**Description**

Truncated Normal and Uniform distributions, where the response is also subject to a Normally distributed measurement error.

**Usage**

```
dmenorm(x, mean=0, sd=1, lower=-Inf, upper=Inf, sderr=0, meanerr=0,
        log = FALSE)
pmenorm(q, mean=0, sd=1, lower=-Inf, upper=Inf, sderr=0, meanerr=0,
        lower.tail = TRUE, log.p = FALSE)
qmenorm(p, mean=0, sd=1, lower=-Inf, upper=Inf, sderr=0, meanerr=0,
        lower.tail = TRUE, log.p = FALSE)
rmenorm(n, mean=0, sd=1, lower=-Inf, upper=Inf, sderr=0, meanerr=0)
dmeunif(x, lower=0, upper=1, sderr=0, meanerr=0, log = FALSE)
pmeunif(q, lower=0, upper=1, sderr=0, meanerr=0, lower.tail = TRUE,
        log.p = FALSE)
qmeunif(p, lower=0, upper=1, sderr=0, meanerr=0, lower.tail = TRUE,
        log.p = FALSE)
rmeunif(n, lower=0, upper=1, sderr=0, meanerr=0)
```

**Arguments**

<code>x, q</code>	vector of quantiles.
<code>p</code>	vector of probabilities.
<code>n</code>	number of observations. If <code>length(n) &gt; 1</code> , the length is taken to be the number required.
<code>mean</code>	vector of means.
<code>sd</code>	vector of standard deviations.
<code>lower</code>	lower truncation point.
<code>upper</code>	upper truncation point.

<code>sderr</code>	Standard deviation of measurement error distribution.
<code>meanerr</code>	Optional shift for the measurement error distribution.
<code>log, log.p</code>	logical; if TRUE, probabilities $p$ are given as $\log(p)$ , or log density is returned.
<code>lower.tail</code>	logical; if TRUE (default), probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .

## Details

The normal distribution with measurement error has density

$$\frac{\Phi(u, \mu_2, \sigma_3) - \Phi(l, \mu_2, \sigma_3)}{\Phi(u, \mu_0, \sigma_0) - \Phi(l, \mu_0, \sigma_0)} \phi(x, \mu_0 + \mu_\epsilon, \sigma_2)$$

where

$$\begin{aligned}\sigma_2^2 &= \sigma_0^2 + \sigma_\epsilon^2, \\ \sigma_3 &= \sigma_0 \sigma_\epsilon / \sigma_2, \\ \mu_2 &= (x - \mu_\epsilon) \sigma_0^2 + \mu_0 \sigma_\epsilon^2,\end{aligned}$$

$\mu_0$  is the mean of the original Normal distribution before truncation,

$\sigma_0$  is the corresponding standard deviation,

$u$  is the upper truncation point,

$l$  is the lower truncation point,

$\sigma_\epsilon$  is the standard deviation of the additional measurement error,

$\mu_\epsilon$  is the mean of the measurement error (usually 0).

$\phi(x)$  is the density of the corresponding normal distribution, and

$\Phi(x)$  is the distribution function of the corresponding normal distribution.

The uniform distribution with measurement error has density

$$(\Phi(x, \mu_\epsilon + l, \sigma_\epsilon) - \Phi(x, \mu_\epsilon + u, \sigma_\epsilon)) / (u - l)$$

These are calculated from the original truncated Normal or Uniform density functions  $f(\cdot | \mu, \sigma, l, u)$  as

$$\int f(y | \mu, \sigma, l, u) \phi(x, y + \mu_\epsilon, \sigma_\epsilon) dy$$

If `sderr` and `meanerr` are not specified they assume the default values of 0, representing no measurement error variance, and no constant shift in the measurement error, respectively.

Therefore, for example with no other arguments, `dmenorm(x)`, is simply equivalent to `dtnorm(x)`, which in turn is equivalent to `dnorm(x)`.

These distributions were used by Satten and Longini (1996) for CD4 cell counts conditionally on hidden Markov states of HIV infection, and later by Jackson and Sharples (2002) for FEV1 measurements conditionally on states of chronic lung transplant rejection.

These distribution functions are just provided for convenience, and are not optimised for numerical accuracy or speed. To fit a hidden Markov model with these response distributions, use a [hmmMETNorm](#) or [hmmMEUnif](#) constructor. See the [hmm-dists](#) help page for further details.

**Value**

dmenorm, dmeunif give the density, pmenorm, pmeunif give the distribution function, qmenorm, qmeunif give the quantile function, and rmenorm, rmeunif generate random deviates, for the Normal and Uniform versions respectively.

**Author(s)**

C. H. Jackson <chris.jackson@mrc-bsu.cam.ac.uk>

**References**

Satten, G.A. and Longini, I.M. Markov chains with measurement error: estimating the 'true' course of a marker of the progression of human immunodeficiency virus disease (with discussion) *Applied Statistics* 45(3): 275-309 (1996)

Jackson, C.H. and Sharples, L.D. Hidden Markov models for the onset and progression of bronchiolitis obliterans syndrome in lung transplant recipients *Statistics in Medicine*, 21(1): 113–128 (2002).

**See Also**

[dnorm](#), [dunif](#), [dtnorm](#)

**Examples**

```
## what does the distribution look like?
x <- seq(50, 90, by=1)
plot(x, dnorm(x, 70, 10), type="l", ylim=c(0,0.06)) ## standard Normal
lines(x, dtnorm(x, 70, 10, 60, 80), type="l")      ## truncated Normal
## truncated Normal with small measurement error
lines(x, dmenorm(x, 70, 10, 60, 80, sderr=3), type="l")
```

---

model.frame.msm

*Extract original data from msm objects.*

---

**Description**

Extract the data from a multi-state model fitted with msm.

**Usage**

```
## S3 method for class 'msm'
model.frame(formula, agg=FALSE, ...)
## S3 method for class 'msm'
model.matrix(object, model="intens", state=1, ...)
```

**Arguments**

formula	A fitted multi-state model object, as returned by <a href="#">msm</a> .
agg	Return the model frame in the efficient aggregated form used to calculate the likelihood internally for non-hidden Markov models. This has one row for each unique combination of from-state, to-state, time lag, covariate value and observation type. The variable named "(nocc)" counts how many observations of that combination there are in the original data.
object	A fitted multi-state model object, as returned by <a href="#">msm</a> .
model	"intens" to return the design matrix for covariates on intensities, "misc" for misclassification probabilities, "hmm" for a general hidden Markov model, and "inits" for initial state probabilities in hidden Markov models.
state	State corresponding to the required covariate design matrix in a hidden Markov model.
...	Further arguments (not used).

**Value**

`model.frame` returns a data frame with all the original variables used for the model fit, with any missing data removed (see `na.action` in [msm](#)). The state, time, subject, obstype and obstrue variables are named "(state)", "(time)", "(subject)", "(obstype)" and "(obstrue)" respectively (note the brackets). A variable called "(obs)" is the observation number from the original data before any missing data were dropped. The variable "(pcomb)" is used for computing the likelihood for hidden Markov models, and identifies which distinct time difference, obstype and covariate values (thus which distinct interval transition probability matrix) each observation corresponds to.

The model frame object has some other useful attributes, including "usernames" giving the user's original names for these variables (used for model refitting, e.g. in bootstrapping or cross validation) and "covnames" identifying which ones are covariates.

`model.matrix` returns a design matrix for a part of the model that includes covariates. The required part is indicated by the "model" argument.

For time-inhomogeneous models fitted with "pci", these datasets will have imputed observations at each time change point, indicated where the variable "(pci.imp)" in the model frame is 1. The model matrix for intensities will have factor contrasts for the timeperiod covariate.

**Author(s)**

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>

**See Also**

[msm](#), [model.frame](#), [model.matrix](#).

**Description**

Fit a continuous-time Markov or hidden Markov multi-state model by maximum likelihood. Observations of the process can be made at arbitrary times, or the exact times of transition between states can be known. Covariates can be fitted to the Markov chain transition intensities or to the hidden Markov observation process.

**Usage**

```
msm ( formula, subject=NULL, data = list(), qmatrix, gen.inits = FALSE,
      ematrix=NULL, hmodel=NULL, obstype=NULL, obstrue=NULL,
      covariates = NULL, covinits = NULL, constraint = NULL,
      misccovariates = NULL, misccovinits = NULL, miscconstraint = NULL,
      hcovariates = NULL, hcovinits = NULL, hconstraint = NULL, hranges=NULL,
      qconstraint=NULL, econstraint=NULL, initprobs = NULL,
      est.initprobs=FALSE, initcovariates = NULL, initcovinits = NULL,
      deathexact = NULL, death=NULL, exacttimes = FALSE, censor=NULL,
      censor.states=NULL, pci=NULL, phase.states=NULL, phase.inits=NULL,
      cl = 0.95, fixedpars = NULL, center=TRUE,
      opt.method="optim", hessian=NULL, use.deriv=TRUE,
      use.expm=TRUE, analyticp=TRUE, na.action=na.omit, ... )
```

**Arguments**

formula	A formula giving the vectors containing the observed states and the corresponding observation times. For example, state ~ time Observed states should be numeric variables in the set $1, \dots, n$ , where $n$ is the number of states. Factors are allowed only if their levels are called "1", ..., "n". The times can indicate different types of observation scheme, so be careful to choose the correct obstype. For hidden Markov models, state refers to the outcome variable, which need not be a discrete state. It may also be a matrix, giving multiple observations at each time (see <a href="#">hmmMV</a> ).
subject	Vector of subject identification numbers for the data specified by formula. If missing, then all observations are assumed to be on the same subject. These must be sorted so that all observations on the same subject are adjacent.
data	Optional data frame in which to interpret the variables supplied in formula, subject, covariates, misccovariates, hcovariates, obstype and obstrue.
qmatrix	Matrix which indicates the allowed transitions in the continuous-time Markov chain, and optionally also the initial values of those transitions. If an instantaneous transition is not allowed from state $r$ to state $s$ , then <code>qmatrix</code> should have $(r, s)$ entry 0, otherwise it should be non-zero.

If supplying initial values yourself, then the non-zero entries should be those values. If using `gen.inits=TRUE` then the non-zero entries can be anything you like (conventionally 1). Any diagonal entry of `qmatrix` is ignored, as it is constrained to be equal to minus the sum of the rest of the row.

For example,

```
rbind( c( 0, 0.1, 0.01 ), c( 0.1, 0, 0.2 ), c( 0, 0, 0 ) )
```

represents a 'health - disease - death' model, with initial transition intensities 0.1 from health to disease, 0.01 from health to death, 0.1 from disease to health, and 0.2 from disease to death.

If the states represent ordered levels of severity of a disease, then this matrix should usually only allow transitions between adjacent states. For example, if someone was observed in state 1 ("mild") at their first observation, followed by state 3 ("severe") at their second observation, they are assumed to have passed through state 2 ("moderate") in between, and the 1,3 entry of `qmatrix` should be zero.

The initial intensities given here are with any covariates set to their means in the data (or set to zero, if `center = FALSE`). If any intensities are constrained to be equal using `qconstraint`, then the initial value is taken from the first of these (reading across rows).

`gen.inits` If TRUE, then initial values for the transition intensities are generated automatically using the method in `crudeinits.msm`. The non-zero entries of the supplied `qmatrix` are assumed to indicate the allowed transitions of the model. This is not available for hidden Markov models, including models with misclassified states.

`ematrix` If misclassification between states is to be modelled, this should be a matrix of initial values for the misclassification probabilities. The rows represent underlying states, and the columns represent observed states. If an observation of state  $s$  is not possible when the subject occupies underlying state  $r$ , then `ematrix` should have  $(r, s)$  entry 0. Otherwise `ematrix` should have  $(r, s)$  entry corresponding to the probability of observing  $s$  conditionally on occupying true state  $r$ . The diagonal of `ematrix` is ignored, as rows are constrained to sum to 1. For example,

```
rbind( c( 0, 0.1, 0 ), c( 0.1, 0, 0.1 ), c( 0, 0.1, 0 ) )
```

represents a model in which misclassifications are only permitted between adjacent states.

If any probabilities are constrained to be equal using `econstraint`, then the initial value is taken from the first of these (reading across rows).

For an alternative way of specifying misclassification models, see `hmodel`.

`hmodel` Specification of the hidden Markov model (HMM). This should be a list of return values from HMM constructor functions. Each element of the list corresponds to the outcome model conditionally on the corresponding underlying state. Univariate constructors are described in the `hmm-dists` help page. These



may also be grouped together to specify a multivariate HMM with a set of conditionally independent univariate outcomes at each time, as described in [hmmMV](#).

For example, consider a three-state hidden Markov model. Suppose the observations in underlying state 1 are generated from a Normal distribution with mean 100 and standard deviation 16, while observations in underlying state 2 are Normal with mean 54 and standard deviation 18. Observations in state 3, representing death, are exactly observed, and coded as 999 in the data. This model is specified as

```
hmodel = list(hmmNorm(mean=100, sd=16), hmmNorm(mean=54, sd=18), hmmIdent(999))
```

The mean and standard deviation parameters are estimated starting from these initial values. If multiple parameters are constrained to be equal using `hconstraint`, then the initial value is taken from the value given on the first occasion that parameter appears in `hmodel`.

See the [hmm-dists](#) help page for details of the constructor functions for each univariate distribution.

A misclassification model, that is, a hidden Markov model where the outcomes are misclassified observations of the underlying states, can either be specified using a list of [hmmCat](#) or [hmmIdent](#) objects, or by using an `ematrix`.

For example,

```
ematrix = rbind( c( 0, 0.1, 0, 0 ), c( 0.1, 0, 0.1, 0 ), c( 0, 0.1, 0, 0 ), c( 0, 0, 0, 0 )
)
```

is equivalent to

```
hmodel = list( hmmCat(prob=c(0.9, 0.1, 0, 0)), hmmCat(prob=c(0.1, 0.8, 0.1, 0)), hmmCat(prob=
```

`obstype`

A vector specifying the observation scheme for each row of the data. This can be included in the data frame `data` along with the state, time, subject IDs and covariates. Its elements should be either 1, 2 or 3, meaning as follows:

- 1** An observation of the process at an arbitrary time (a "snapshot" of the process, or "panel-observed" data). The states are unknown between observation times.
- 2** An exact transition time, with the state at the previous observation retained until the current observation. An observation may represent a transition to a different state or a repeated observation of the same state (e.g. at the end of follow-up). Note that if all transition times are known, more flexible models could be fitted with packages other than **msm** - see the note under `exacttimes`.  
Note also that if the previous state was censored using  `censor`, for example known only to be state 1 or state 2, then `obstype 2` means that either state 1 is retained or state 2 is retained until the current observation - this does not allow for a change of state in the middle of the observation interval.
- 3** An exact transition time, but the state at the instant before entering this state is unknown. A common example is death times in studies of chronic diseases.

If `obstype` is not specified, this defaults to all 1. If `obstype` is a single number, all observations are assumed to be of this type. The `obstype` value for the first observation from each subject is not used.

This is a generalisation of the `deathexact` and `exacttimes` arguments to allow different schemes per observation. `obstype` overrides both `deathexact` and `exacttimes`.

`exacttimes=TRUE` specifies that all observations are of `obstype` 2.

`deathexact = death.states` specifies that all observations of `death.states` are of type 3. `deathexact = TRUE` specifies that all observations in the final absorbing state are of type 3.

`obstrue` In misclassification models specified with `ematrix`, `obstrue` is a vector of logicals (TRUE or FALSE) or numerics (1 or 0) specifying which observations (TRUE, 1) are observations of the underlying state without error, and which (FALSE, 0) are realisations of a hidden Markov model.

In HMMs specified with `hmodel`, where the hidden state is known at some times, if `obstrue` is supplied it is assumed to contain the actual true state data. Elements of `obstrue` at times when the hidden state is unknown are set to NA. This allows the information from HMM outcomes generated conditionally on the known state to be included in the model, thus improving the estimation of the HMM outcome distributions.

In HMMs where there are also censored states, `obstrue` should be set to 1 for observed states which are *censored* but not *misclassified*.

`covariates` A formula or a list of formulae representing the covariates on the transition intensities via a log-linear model. If a single formula is supplied, like `covariates = ~ age + sex + treatment`

then these covariates are assumed to apply to all intensities. If a named list is supplied, then this defines a potentially different model for each named intensity. For example,

```
covariates = list("1-2" = ~ age, "2-3" = ~ age + treatment)
```

specifies an age effect on the state 1 - state 2 transition, additive age and treatment effects on the state 2 - state 3 transition, but no covariates on any other transitions that are allowed by the `qmatrix`.

If covariates are time dependent, they are assumed to be constant in between the times they are observed, and the transition probability between a pair of times ( $t_1, t_2$ ) is assumed to depend on the covariate value at  $t_1$ .

`covinits` Initial values for log-linear effects of covariates on the transition intensities. This should be a named list with each element corresponding to a covariate. A single element contains the initial values for that covariate on each transition intensity, reading across the rows in order. For a pair of effects constrained to be equal, the initial value for the first of the two effects is used.

For example, for a model with the above `qmatrix` and age and sex covariates, the following initialises all covariate effects to zero apart from the age effect on the 2-1 transition, and the sex effect on the 1-3 transition. `covinits = list(sex=c(0, 0, 0.1, 0), age=c(0, 0.1, 0, 0))`

For factor covariates, name each level by concatenating the name of the covariate with the level name, quoting if necessary. For example, for a covariate `agegroup` with three levels 0-15, 15-60, 60-, use something like

```
covinits = list("agegroup15-60"=c(0,0.1,0,0), "agegroup60-"=c(0.1,0.1,0,0))
```

If not specified or wrongly specified, initial values are assumed to be zero.

constraint

A list of one numeric vector for each named covariate. The vector indicates which covariate effects on intensities are constrained to be equal. Take, for example, a model with five transition intensities and two covariates. Specifying

```
constraint = list (age = c(1,1,1,2,2), treatment = c(1,2,3,4,5))
```

constrains the effect of age to be equal for the first three intensities, and equal for the fourth and fifth. The effect of treatment is assumed to be different for each intensity. Any vector of increasing numbers can be used as indicators. The intensity parameters are assumed to be ordered by reading across the rows of the transition matrix, starting at the first row, ignoring the diagonals.

Negative elements of the vector can be used to indicate that particular covariate effects are constrained to be equal to minus some other effects. For example:

```
constraint = list (age = c(-1,1,1,2,-2), treatment = c(1,2,3,4,5))
```

constrains the second and third age effects to be equal, the first effect to be minus the second, and the fifth age effect to be minus the fourth. For example, it may be realistic that the effect of a covariate on the "reverse" transition rate from state 2 to state 1 is minus the effect on the "forward" transition rate, state 1 to state 2. Note that it is not possible to specify exactly which of the covariate effects are constrained to be positive and which negative. The maximum likelihood estimation chooses the combination of signs which has the higher likelihood.

For categorical covariates, defined as factors, specify constraints as follows:

```
list(..., covnameVALUE1 = c(...), covnameVALUE2 = c(...), ...)
```

where covname is the name of the factor, and VALUE1, VALUE2, ... are the labels of the factor levels (usually excluding the baseline, if using the default contrasts). Make sure the contrasts option is set appropriately, for example, the default

```
options(contrasts=c(contr.treatment, contr.poly))
```

sets the first (baseline) level of unordered factors to zero, then the baseline level is ignored in this specification.

To assume no covariate effect on a certain transition, use the fixedpars argument to fix it at its initial value (which is zero by default) during the optimisation.

misccovariates

A formula representing the covariates on the misclassification probabilities, analogously to covariates, via multinomial logistic regression. Only used if the model is specified using ematrix, rather than hmodel.

This must be a single formula - lists are not supported, unlike covariates. If a different model on each probability is required, include all covariates in this formula, and use fixedpars to fix some of their effects (for particular probabilities) at their default initial values of zero.

misccovinits

Initial values for the covariates on the misclassification probabilities, defined in the same way as covinits. Only used if the model is specified using ematrix.

- misconstraint** A list of one vector for each named covariate on misclassification probabilities. The vector indicates which covariate effects on misclassification probabilities are constrained to be equal, analogously to `constraint`. Only used if the model is specified using `ematrix`.
- hcovariates** List of formulae the same length as `hmodel`, defining any covariates governing the hidden Markov outcome models. The covariates operate on a suitably link-transformed linear scale, for example, log scale for a Poisson outcome model. If there are no covariates for a certain hidden state, then insert a `NULL` in the corresponding place in the list. For example, `hcovariates = list(~acute + age, ~acute, NULL)`.
- hcovinits** Initial values for the hidden Markov model covariate effects. A list of the same length as `hcovariates`. Each element is a vector with initial values for the effect of each covariate on that state. For example, the above `hcovariates` can be initialised with `hcovariates = list(c(-8, 0), -8, NULL)`. Initial values must be given for all or no covariates, if none are given these are all set to zero. The initial value given in the `hmodel` constructor function for the corresponding baseline parameter is interpreted as the value of that parameter with any covariates fixed to their means in the data. If multiple effects are constrained to be equal using `hconstraint`, then the initial value is taken from the first of the multiple initial values supplied.
- hconstraint** A named list. Each element is a vector of constraints on the named hidden Markov model parameter. The vector has length equal to the number of times that class of parameter appears in the whole model.  
 For example consider the three-state hidden Markov model described above, with normally-distributed outcomes for states 1 and 2. To constrain the outcome variance to be equal for states 1 and 2, and to also constrain the effect of acute on the outcome mean to be equal for states 1 and 2, specify  
`hconstraint = list(sd = c(1,1), acute=c(1,1))`  
 Note this excludes initial state occupancy probabilities and covariate effects on those probabilities, which cannot be constrained.
- hranges** Range constraints for hidden Markov model parameters. Supplied as a named list, with each element corresponding to the named hidden Markov model parameter. This element is itself a list with two elements, vectors named "lower" and "upper". These vectors each have length equal to the number of times that class of parameter appears in the whole model, and give the corresponding minimum and maximum allowable values for that parameter. Maximum likelihood estimation is performed with these parameters constrained in these ranges (through a log or logit-type transformation). Lower bounds of `-Inf` and upper bounds of `Inf` can be given if the parameter is unbounded above or below. For example, in the three-state model above, to constrain the mean for state 1 to be between 0 and 6, and the mean of state 2 to be between 7 and 12, supply  
`hranges=list(mean=list(lower=c(0,7), upper=c(6,12)))`  
 These default to the natural ranges, e.g. the positive real line for variance parameters, and `[0,1]` for probabilities. Therefore `hranges` need not be specified for such parameters unless an even stricter constraint is desired. If only one limit is supplied for a parameter, only the first occurrence of that parameter is constrained.

	Initial values should be strictly within any ranges, and not on the range boundary, otherwise optimisation will fail with a "non-finite value" error.
qconstraint	<p>A vector of indicators specifying which baseline transition intensities are equal. For example,</p> <pre>qconstraint = c(1,2,3,3)</pre> <p>constrains the third and fourth intensities to be equal, in a model with four allowed instantaneous transitions. When there are covariates on the intensities and center=TRUE (the default), qconstraint is applied to the intensities with covariates taking the values of the means in the data. When center=FALSE, qconstraint is applied to the intensities with covariates set to zero.</p>
econstraint	A similar vector of indicators specifying which baseline misclassification probabilities are constrained to be equal. Only used if the model is specified using ematrix, rather than hmodel.
initprobs	<p>Only used in hidden Markov models. Underlying state occupancy probabilities at each subject's first observation. Can either be a vector of <i>nstates</i> elements with common probabilities to all subjects, or a <i>nsubjects</i> by <i>nstates</i> matrix of subject-specific probabilities. This refers to observations after missing data and subjects with only one observation have been excluded.</p> <p>If these are estimated (see est.initprobs), then this represents an initial value, and defaults to equal probability for each state. Otherwise this defaults to <code>c(1, rep(0, nstates-1))</code>, that is, in state 1 with a probability of 1. Scaled to sum to 1 if necessary. The state 1 occupancy probability should be non-zero.</p>
est.initprobs	<p>Only used in hidden Markov models. If TRUE, then the underlying state occupancy probabilities at the first observation will be estimated, starting from a vector of initial values supplied in the initprobs argument. Structural zeroes are allowed: if any of these initial values are zero they will be fixed at zero during optimisation, even if est.initprobs=TRUE, and no covariate effects on them are estimated. The exception is state 1, which should have non-zero occupancy probability.</p> <p>Note that the free parameters during this estimation exclude the state 1 occupancy probability, which is fixed at one minus the sum of the other probabilities.</p>
initcovariates	Formula representing covariates on the initial state occupancy probabilities, via multinomial logistic regression. The linear effects of these covariates, observed at the individual's first observation time, operate on the log ratio of the state <i>r</i> occupancy probability to the state 1 occupancy probability, for each <i>r</i> = 2 to the number of states. Thus the state 1 occupancy probability should be non-zero. If est.initprobs is TRUE, these effects are estimated starting from their initial values. If est.initprobs is FALSE, these effects are fixed at their initial values.
initcovinits	Initial values for the covariate effects initcovariates. A named list with each element corresponding to a covariate, as in covinits. Each element is a vector with (1 - number of states) elements, containing the initial values for the linear effect of that covariate on the log odds of that state relative to state 1, from state 2 to the final state. If initcovinits is not specified, all covariate effects are initialised to zero.
deathexact	Vector of indices of absorbing states whose time of entry is known exactly, but the individual is assumed to be in an unknown transient state ("alive") at the

previous instant. This is the usual situation for times of death in chronic disease monitoring data. For example, if you specify `deathexact = c(4, 5)` then states 4 and 5 are assumed to be exactly-observed death states.

See the `obstype` argument. States of this kind correspond to `obstype=3`. `deathexact = TRUE` indicates that the final absorbing state is of this kind, and `deathexact = FALSE` or `deathexact = NULL` (the default) indicates that there is no state of this kind.

The `deathexact` argument is overridden by `obstype` or `exacttimes`.

Note that you do not always supply a `deathexact` argument, even if there are states that correspond to deaths, because they do not necessarily have `obstype=3`. If the state is known between the time of death and the previous observation, then you should specify `obstype=2` for the death times, or `exacttimes=TRUE` if the state is known at all times, and the `deathexact` argument is ignored.

`death` Old name for the `deathexact` argument. Overridden by `deathexact` if both are supplied. Deprecated.

`censor` A state, or vector of states, which indicates censoring. Censoring means that the observed state is known only to be one of a particular set of states. For example, `censor=999` indicates that all observations of 999 in the vector of observed states are censored states. By default, this means that the true state could have been any of the transient (non-absorbing) states. To specify corresponding true states explicitly, use a `censor.states` argument.

Note that in contrast to the usual terminology of survival analysis, here it is the *state* which is considered to be censored, rather than the *event time*. If at the end of a study, an individual has not died, but their true state is *known*, then `censor` is unnecessary, since the standard multi-state model likelihood is applicable. Also a "censored" state here can be at any time, not just at the end.

Note in particular that general time-inhomogeneous Markov models with piecewise constant transition intensities can be constructed using the `censor` facility. If the true state is unknown on occasions when a piecewise constant covariate is known to change, then censored states can be inserted in the data on those occasions. The covariate may represent time itself, in which case the `pci` option to `msm` can be used to perform this trick automatically, or some other time-dependent variable.

`censor.states` Specifies the underlying states which censored observations can represent. If `censor` is a single number (the default) this can be a vector, or a list with one element. If `censor` is a vector with more than one element, this should be a list, with each element a vector corresponding to the equivalent element of `censor`. For example

```
censor = c(99, 999), censor.states = list(c(2, 3), c(3, 4))
```

means that observations coded 99 represent either state 2 or state 3, while observations coded 999 are really either state 3 or state 4.

`pci` Model for piecewise-constant intensities. Vector of cut points defining the times, since the start of the process, at which intensities change for all subjects. For example

```
pci = c(5, 10)
```

specifies that the intensity changes at time points 5 and 10. This will automatically construct a model with a categorical (factor) covariate called `timeperiod`,

with levels " $[-\text{Inf}, 5)$ ", " $[5, 10)$ " and " $[10, \text{Inf})$ ", where the first level is the baseline. This covariate defines the time period in which the observation was made. Initial values and constraints on covariate effects are specified the same way as for a model with a covariate of this name, for example,

```
covinits = list("timeperiod[5,10)"=c(0.1,0.1), "timeperiod[10,Inf)"=c(0.1,0.1))
```

Thus if `pci` is supplied, you cannot have a previously-existing variable called `timeperiod` as a covariate in any part of a `msm` model.

To assume piecewise constant intensities for some transitions but not others with `pci`, use the `fixedpars` argument to fix the appropriate covariate effects at their default initial values of zero.

Internally, this works by inserting censored observations in the data at times when the intensity changes but the state is not observed.

If the supplied times are outside the range of the time variable in the data, `pci` is ignored and a time-homogeneous model is fitted.

After fitting a time-inhomogeneous model, `qmatrix.msm` can be used to obtain the fitted intensity matrices for each time period, for example,

```
qmatrix.msm(example.msm, covariates=list(timeperiod="[5, Inf)"))
```

This facility does not support interactions between time and other covariates. Such models need to be specified "by hand", using a state variable with censored observations inserted. Note that the data component of the `msm` object returned from a call to `msm` with `pci` supplied contains the states with inserted censored observations and time period indicators. These can be used to construct such models.

Note that you do not need to use `pci` in order to model the effect of a time-dependent covariate in the data. `msm` will automatically assume that covariates are piecewise-constant and change at the times when they are observed. `pci` is for when you want all intensities to change at the same pre-specified times for all subjects.

#### phase.states

Indices of states which have a two-phase sojourn distribution. This defines a semi-Markov model, in which the hazard of an onward transition depends on the time spent in the state.

This uses the technique described by Titman and Sharples (2009). A hidden Markov model is automatically constructed on an expanded state space, where the phases correspond to the hidden states. The "tau" proportionality constraint described in this paper is currently not supported.

Covariates, constraints, `deathexact` and `sensor` are expressed with respect to the expanded state space. If not supplied by hand, `initprobs` is defined automatically so that subjects are assumed to begin in the first of the two phases.

Hidden Markov models can additionally be given phased states. The user supplies an outcome distribution for each original state using `hmodel`, which is expanded internally so that it is assumed to be the same within each of the phased states. `initprobs` is interpreted on the expanded state space. Misclassification models defined using `ematrix` are not supported, and these must be defined using `hmmCat` or `hmmIdent` constructors, as described in the `hmodel` section of this help page. Or the HMM on the expanded state space can be defined by hand.

Output functions are presented as it were a hidden Markov model on the expanded state space, for example, transition probabilities between states, covari-

ate effects on transition rates, or prevalence counts, are not aggregated over the hidden phases.

Numerical estimation will be unstable when there is weak evidence for a two-phase sojourn distribution, that is, if the model is close to Markov.

See [d2phase](#) for the definition of the two-phase distribution and the interpretation of its parameters.

This is an experimental feature, and some functions are not implemented. Please report any experiences of using this feature to the author!

phase.inits	<p>Initial values for phase-type models. A list with one component for each "two-phased" state. Each component is itself a list of two elements. The first of these elements is a scalar defining the transition intensity from phase 1 to phase 2. The second element is a matrix, with one row for each potential destination state from the two-phased state, and two columns. The first column is the transition rate from phase 1 to the destination state, and the second column is the transition rate from phase 2 to the destination state. If there is only one destination state, then this may be supplied as a vector.</p> <p>In phase type models, the initial values for transition rates out of non-phased states are taken from the <code>qmatrix</code> supplied to <code>msm</code>, and entries of this matrix corresponding to transitions out of phased states are ignored.</p>
exacttimes	<p>By default, the transitions of the Markov process are assumed to take place at unknown occasions in between the observation times. If <code>exacttimes</code> is set to <code>TRUE</code>, then the observation times are assumed to represent the exact times of transition of the process. The subject is assumed to be in the same state between these times. An observation may represent a transition to a different state or a repeated observation of the same state (e.g. at the end of follow-up). This is equivalent to every row of the data having <code>obstype = 2</code>. See the <code>obstype</code> argument. If both <code>obstype</code> and <code>exacttimes</code> are specified then <code>exacttimes</code> is ignored.</p> <p>Note that the complete history of the multi-state process is known with this type of data. The models which <code>msm</code> fits have the strong assumption of constant (or piecewise-constant) transition rates. Knowing the exact transition times allows more realistic models to be fitted with other packages. For example parametric models with sojourn distributions more flexible than the exponential can be fitted with the <code>flexsurv</code> package, or semi-parametric models can be implemented with <code>survival</code> in conjunction with <code>mstate</code>.</p>
c1	<p>Width of symmetric confidence intervals for maximum likelihood estimates, by default 0.95.</p>
fixedpars	<p>Vector of indices of parameters whose values will be fixed at their initial values during the optimisation. These are given in the order: transition intensities (reading across rows of the transition matrix), covariates on intensities (ordered by intensities within covariates), hidden Markov model parameters, including misclassification probabilities (ordered by parameters within states), hidden Markov model covariate parameters (ordered by covariates within parameters within states), initial state occupancy probabilities (excluding the first probability, which is fixed at one minus the sum of the others).</p> <p>If there are equality constraints on certain parameters, then <code>fixedpars</code> indexes the set of unique parameters, excluding those which are constrained to be equal</p>



	<p>to previous parameters.</p> <p>To fix all parameters, specify <code>fixedpars = TRUE</code>.</p> <p>This can be useful for profiling likelihoods, and building complex models stage by stage.</p>
<code>center</code>	<p>If <code>TRUE</code> (the default, unless <code>fixedpars=TRUE</code>) then covariates are centered at their means during the maximum likelihood estimation. This usually improves stability of the numerical optimisation.</p>
<code>opt.method</code>	<p>If "optim", "nlm" or "bobyqa", then the corresponding R function will be used for maximum likelihood estimation. <code>optim</code> is the default. "bobyqa" requires the package <b>minqa</b> to be installed. See the help of these functions for further details. Advanced users can also add their own optimisation methods, see the source for <code>optim.R</code> in <code>msm</code> for some examples.</p> <p>If "fisher", then a specialised Fisher scoring method is used (Kalbfleisch and Lawless, 1985) which can be faster than the generic methods, though less robust. This is only available for Markov models with panel data (<code>obstype=1</code>), that is, not for models with censored states, hidden Markov models, exact observation or exact death times (<code>obstype=2, 3</code>).</p>
<code>hessian</code>	<p>If <code>TRUE</code> then standard errors and confidence intervals are obtained from a numerical estimate of the Hessian (the observed information matrix). This is the default when maximum likelihood estimation is performed. If all parameters are fixed at their initial values and no optimisation is performed, then this defaults to <code>FALSE</code>. If requested, the actual Hessian is returned in <code>x\$paramdata\$opt\$hessian</code>, where <code>x</code> is the fitted model object.</p> <p>If <code>hessian</code> is set to <code>FALSE</code>, then standard errors and confidence intervals are obtained from the Fisher (expected) information matrix, if this is available. This may be preferable if the numerical estimation of the Hessian is computationally intensive, or if the resulting estimate is non-invertible or not positive definite.</p>
<code>use.deriv</code>	<p>If <code>TRUE</code> then analytic first derivatives are used in the optimisation of the likelihood, where available and an appropriate quasi-Newton optimisation method, such as BFGS, is being used. Analytic derivatives are not available for all models.</p>
<code>use.expm</code>	<p>If <code>TRUE</code> then any matrix exponentiation needed to calculate the likelihood is done using the <b>expm</b> package. Otherwise the original routines used in <b>msm</b> 1.2.4 and earlier are used. Set to <code>FALSE</code> for backward compatibility, and let the package maintainer know if this gives any substantive differences.</p>
<code>analyticp</code>	<p>By default, the likelihood for certain simpler 3, 4 and 5 state models is calculated using an analytic expression for the transition probability (P) matrix. For all other models, matrix exponentiation is used to obtain P. To revert to the original method of using the matrix exponential for all models, specify <code>analyticp=FALSE</code>. See the PDF manual for a list of the models for which analytic P matrices are implemented.</p>
<code>na.action</code>	<p>What to do with missing data: either <code>na.omit</code> to drop it and carry on, or <code>na.fail</code> to stop with an error. Missing data includes all NAs in the states, times, subject or <code>obstrue</code>, all NAs at the first observation for a subject for covariates in <code>initcovariates</code>, all NAs in other covariates (excluding the last observation for a subject), all NAs in <code>obstype</code> (excluding the first observation for a subject), and any subjects with only one observation (thus no observed transitions).</p>

... Optional arguments to the general-purpose R optimisation routine, `optim` by default. For example `method="Nelder-Mead"` to change the optimisation algorithm from the "BFGS" method that `msm` calls by default.

It is often worthwhile to normalize the optimisation using `control=list(fnscale = a)`, where `a` is the a number of the order of magnitude of the -2 log likelihood.

If 'false' convergence is reported and the standard errors cannot be calculated due to a non-positive-definite Hessian, then consider tightening the tolerance criteria for convergence. If the optimisation takes a long time, intermediate steps can be printed using the `trace` argument of the control list. See `optim` for details.

For the Fisher scoring method, a control list can be supplied in the same way, but the only supported options are `reltol`, `trace` and `damp`. The first two are used in the same way as for `optim`. If the algorithm fails with a singular information matrix, adjust `damp` from the default of zero (to, e.g. 1). This adds a constant identity matrix multiplied by `damp` to the information matrix during optimisation.

## Details

For full details about the methodology behind the `msm` package, refer to the PDF manual '`msm-manual.pdf`' in the 'doc' subdirectory of the package. This includes a tutorial in the typical use of `msm`. The paper by Jackson (2011) in *Journal of Statistical Software* presents the material in this manual in a more concise form.

`msm` was designed for fitting *continuous-time* Markov models, processes where transitions can occur at any time. These models are defined by *intensities*, which govern both the time spent in the current state and the probabilities of the next state. In *discrete-time models*, transitions are known in advance to only occur at multiples of some time unit, and the model is purely governed by the probability distributions of the state at the next time point, conditionally on the state at the current time. These can also be fitted in `msm`, assuming that there is a continuous-time process underlying the data. Then the fitted transition probability matrix over one time period, as returned by `pmatrix.msm(..., t=1)` is equivalent to the matrix that governs the discrete-time model. However, these can be fitted more efficiently using multinomial logistic regression, for example, using `multinom` from the R package `nnet` (Venables and Ripley, 2002).

For simple continuous-time multi-state Markov models, the likelihood is calculated in terms of the transition intensity matrix  $Q$ . When the data consist of observations of the Markov process at arbitrary times, the exact transition times are not known. Then the likelihood is calculated using the transition probability matrix  $P(t) = \exp(tQ)$ , where  $\exp$  is the matrix exponential. If state  $i$  is observed at time  $t$  and state  $j$  is observed at time  $u$ , then the contribution to the likelihood from this pair of observations is the  $i, j$  element of  $P(u - t)$ . See, for example, Kalbfleisch and Lawless (1985), Kay (1986), or Gentleman *et al.* (1994).

For hidden Markov models, the likelihood for an individual with  $k$  observations is calculated directly by summing over the unknown state at each time, producing a product of  $k$  matrices. The calculation is a generalisation of the method described by Satten and Longini (1996), and also by Jackson and Sharples (2002), and Jackson *et al.* (2003).

There must be enough information in the data on each state to estimate each transition rate, otherwise the likelihood will be flat and the maximum will not be found. It may be appropriate to reduce

the number of states in the model, the number of allowed transitions, or the number of covariate effects, to ensure convergence. Hidden Markov models, and situations where the value of the process is only known at a series of snapshots, are particularly susceptible to non-identifiability, especially when combined with a complex transition matrix. Choosing an appropriate set of initial values for the optimisation can also be important. For flat likelihoods, 'informative' initial values will often be required. See the PDF manual for other tips.

## Value

To obtain summary information from models fitted by the `msm` function, it is recommended to use extractor functions such as `qmatrix.msm`, `pmatrix.msm`, `sojourn.msm`, `msm.form.qoutput`. These provide estimates and confidence intervals for quantities such as transition probabilities for given covariate values.

For advanced use, it may be necessary to directly use information stored in the object returned by `msm`. This is documented in the help page `msm.object`.

Printing a `msm` object by typing the object's name at the command line implicitly invokes `print.msm`. This formats and prints the important information in the model fit, and also returns that information in an R object. This includes estimates and confidence intervals for the transition intensities and (log) hazard ratios for the corresponding covariates. When there is a hidden Markov model, the chief information in the `hmodel` component is also formatted and printed. This includes estimates and confidence intervals for each parameter.

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## References

- Jackson, C.H. (2011). Multi-State Models for Panel Data: The `msm` Package for R., *Journal of Statistical Software*, 38(8), 1-29. URL <http://www.jstatsoft.org/v38/i08/>.
- Kalbfleisch, J., Lawless, J.F., The analysis of panel data under a Markov assumption *Journal of the Americal Statistical Association* (1985) 80(392): 863–871.
- Kay, R. A Markov model for analysing cancer markers and disease states in survival studies. *Biometrics* (1986) 42: 855–865.
- Gentleman, R.C., Lawless, J.F., Lindsey, J.C. and Yan, P. Multi-state Markov models for analysing incomplete disease history data with illustrations for HIV disease. *Statistics in Medicine* (1994) 13(3): 805–821.
- Satten, G.A. and Longini, I.M. Markov chains with measurement error: estimating the 'true' course of a marker of the progression of human immunodeficiency virus disease (with discussion) *Applied Statistics* 45(3): 275-309 (1996)
- Jackson, C.H. and Sharples, L.D. Hidden Markov models for the onset and progression of bronchiolitis obliterans syndrome in lung transplant recipients *Statistics in Medicine*, 21(1): 113–128 (2002).
- Jackson, C.H., Sharples, L.D., Thompson, S.G. and Duffy, S.W. and Couto, E. Multi-state Markov models for disease progression with classification error. *The Statistician*, 52(2): 193–209 (2003)
- Titman, A.C. and Sharples, L.D. Semi-Markov models with phase-type sojourn distributions. *Biometrics* 66, 742-752 (2009).

Venables, W.N. and Ripley, B.D. (2002) *Modern Applied Statistics with S*, second edition. Springer.

### See Also

[simmulti.msm](#), [plot.msm](#), [summary.msm](#), [qmatrix.msm](#), [pmatrix.msm](#), [sojourn.msm](#).

### Examples

```
### Heart transplant data
### For further details and background to this example, see
### Jackson (2011) or the PDF manual in the doc directory.
print(cav[1:10,])
tway4.q <- rbind(c(-0.5, 0.25, 0, 0.25), c(0.166, -0.498, 0.166, 0.166),
c(0, 0.25, -0.5, 0.25), c(0, 0, 0, 0))
statetable.msm(state, PTNUM, data=cav)
crudeinits.msm(state ~ years, PTNUM, data=cav, qmatrix=tway4.q)
cav.msm <- msm( state ~ years, subject=PTNUM, data = cav,
               qmatrix = tway4.q, deathexact = 4,
               control = list ( trace = 2, REPORT = 1 ) )

cav.msm
qmatrix.msm(cav.msm)
pmatrix.msm(cav.msm, t=10)
sojourn.msm(cav.msm)
```

---

msm.form.qoutput

*Extract msm model parameter estimates in compact format*

---

### Description

Extract estimates and confidence intervals for transition intensities (or misclassification probabilities), and their covariate effects, in a tidy matrix format with one row per transition. This is used by the print method ([print.msm](#)) for `msm` objects. Covariate effects are returned as hazard or odds ratios, not on the log scale.

### Usage

```
msm.form.qoutput(x, covariates="mean", cl=0.95, digits=4, ...)
msm.form.eoutput(x, covariates="mean", cl=0.95, digits=4, ...)
```

### Arguments

<code>x</code>	A fitted multi-state model object, as returned by <a href="#">msm</a> .
<code>covariates</code>	Covariate values defining the "baseline" parameters (see <a href="#">qmatrix.msm</a> ).
<code>cl</code>	Width of the symmetric confidence interval to present. Defaults to 0.95.
<code>digits</code>	Minimum number of significant digits for the formatted character matrix returned as an attribute. This is passed to <a href="#">format</a> . Defaults to 4.
<code>...</code>	Other arguments to be passed to <a href="#">format</a> .

**Value**

A numeric matrix with one row per transition, and one column for each estimate or confidence limit. The "formatted" attribute contains the same results formatted for pretty printing. `msm.form.qoutput` returns the transition intensities and their covariates, and `msm.form.eoutput` returns the misclassification probabilities and their covariates.

**Author(s)**

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**See Also**

[print.msm](#)

---

msm.object

*Fitted msm model objects*

---

**Description**

The `msm` function returns a list with the following components. These are intended for developers and confident users. To extract results from fitted model objects, functions such as [qmatrix.msm](#) or [print.msm](#) should be used instead.

**Value**

<code>call</code>	The original call to <code>msm</code> , as returned by <a href="#">match.call</a> .
<code>Qmatrices</code>	A list of matrices. The first component, labelled <code>logbaseline</code> , is a matrix containing the estimated transition intensities on the log scale with any covariates fixed at their means in the data (or at zero, if <code>center=FALSE</code> ). The component labelled <code>baseline</code> is the equivalent on the untransformed scale. Each remaining component is a matrix giving the linear effects of the labelled covariate on the matrix of log intensities. To extract an estimated intensity matrix on the natural scale, at an arbitrary combination of covariate values, use the function <a href="#">qmatrix.msm</a> .
<code>QmatricesSE</code>	The standard error matrices corresponding to <code>Qmatrices</code> .
<code>QmatricesL, QmatricesU</code>	Corresponding lower and upper symmetric confidence limits, of width 0.95 unless specified otherwise by the <code>cl</code> argument.
<code>Ematrices</code>	A list of matrices. The first component, labelled <code>logitbaseline</code> , is the estimated misclassification probability matrix (expressed as log odds relative to the probability of the true state) with any covariates fixed at their means in the data (or at zero, if <code>center=FALSE</code> ). The component labelled <code>baseline</code> is the equivalent on the untransformed scale. Each remaining component is a matrix giving the linear effects of the labelled covariate on the matrix of logit misclassification probabilities. To extract an estimated misclassification probability matrix on the natural scale, at an arbitrary combination of covariate values, use the function <a href="#">ematrix.msm</a> .

<code>EmatricesSE</code>	The standard error matrices corresponding to <code>Ematrices</code> .
<code>EmatricesL,EmatricesU</code>	Corresponding lower and upper symmetric confidence limits, of width 0.95 unless specified otherwise by the <code>cl</code> argument.
<code>minus2loglik</code>	Minus twice the maximised log-likelihood.
<code>deriv</code>	Derivatives of the minus twice log-likelihood at its maximum.
<code>estimates</code>	Vector of untransformed maximum likelihood estimates returned from <code>optim</code> . Transition intensities are on the log scale and misclassification probabilities are given as log odds relative to the probability of the true state.
<code>estimates.t</code>	Vector of transformed maximum likelihood estimates with intensities and probabilities on their natural scales.
<code>fixedpars</code>	Indices of <code>estimates</code> which were fixed during the maximum likelihood estimation.
<code>center</code>	Indicator for whether the estimation was performed with covariates centered on their means in the data.
<code>covmat</code>	Covariance matrix corresponding to <code>estimates</code> .
<code>ci</code>	Matrix of confidence intervals corresponding to <code>estimates.t</code>
<code>opt</code>	Return value from the optimisation routine (such as <code>optim</code> or <code>nlm</code> ), giving information about the results of the optimisation.
<code>foundse</code>	Logical value indicating whether the Hessian was positive-definite at the supposed maximum of the likelihood. If not, the covariance matrix of the parameters is unavailable. In these cases the optimisation has probably not converged to a maximum.
<code>data</code>	A list giving the data used for the model fit, for use in post-processing. To extract it, use the methods <code>model.frame.msm</code> or <code>model.matrix.msm</code> . The format of this element changed in version 1.4 of <code>msm</code> , so that it now contains a <code>model.frame</code> object <code>mf</code> with all the variables used in the model. The previous format (an ad-hoc list of vectors and matrices) can be obtained with the function <code>recreate.olddata(msmobject)</code> , where <code>msmobject</code> is the object returned by <code>msm</code> .
<code>qmodel</code>	A list of objects representing the transition matrix structure and options for likelihood calculation. See <code>qmodel.object</code> for documentation of the components.
<code>emodel</code>	A list of objects representing the misclassification model structure, for models specified using the <code>ematrix</code> argument to <code>msm</code> . See <code>emodel.object</code> .
<code>qcmode1</code>	A list of objects representing the model for covariates on transition intensities. See <code>qcmode1.object</code> .
<code>ecmode1</code>	A list of objects representing the model for covariates on transition intensities. See <code>ecmode1.object</code> .
<code>hmodel</code>	A list of objects representing the hidden Markov model structure. See <code>hmodel.object</code> .
<code>cmode1</code>	A list giving information about censored states. See <code>cmode1.object</code> .
<code>pci</code>	Cut points for time-varying intensities, as supplied to <code>msm</code> , but excluding any that are outside the times observed in the data.

paramdata	A list giving information about the parameters of the multi-state model. See <a href="#">paramdata.object</a> .
cl	Confidence interval width, as supplied to <a href="#">msm</a> .
covariates	Formula for covariates on intensities, as supplied to <a href="#">msm</a> .
misccovariates	Formula for covariates on misclassification probabilities, as supplied to <a href="#">msm</a> .
hcovariates	Formula for covariates on hidden Markov model outcomes, as supplied to <a href="#">msm</a> .
initcovariates	Formula for covariates on initial state occupancy probabilities in hidden Markov models, as supplied to <a href="#">msm</a> .
sojourn	A list as returned by <a href="#">sojourn.msm</a> , with components: mean = estimated mean sojourn times in the transient states, with covariates fixed at their means (if center=TRUE) or at zero (if center=FALSE). se = corresponding standard errors.

---

msm.summary

*Summarise a fitted multi-state model*


---

## Description

Summary method for fitted [msm](#) models. This is simply a wrapper around [prevalence.msm](#) which produces a table of observed and expected state prevalences for each time, and for models with covariates, [hazard.msm](#) to print hazard ratios with 95% confidence intervals for covariate effects.

## Usage

```
## S3 method for class 'msm'
summary(object, hazard.scale=1, ...)
```

## Arguments

object	A fitted multi-state model object, as returned by <a href="#">msm</a> .
hazard.scale	Vector with same elements as number of covariates on transition rates. Corresponds to the increase in each covariate used to calculate its hazard ratio. Defaults to all 1.
...	Further arguments passed to <a href="#">prevalence.msm</a> .

## Value

A list of class `summary.msm`, with components:

prevalences	Output from <a href="#">prevalence.msm</a> .
hazard	Output from <a href="#">hazard.msm</a> .
hazard.scale	Value of the <code>hazard.scale</code> argument.

**Author(s)**

C. H. Jackson <chris.jackson@mrc-bsu.cam.ac.uk>

**See Also**

[msm](#), [prevalence.msm](#), [hazard.msm](#)

---

msm2Surv

*Convert data for 'msm' to data for 'survival', 'mstate' or 'flexsurv' analysis*

---

**Description**

Converts longitudinal data for a [msm](#) model fit, where observations represent the exact transition times of the process, to counting process data. This enables, for example, flexible parametric multi-state models to be fitted with [flexsurvreg](#) from the **flexsurv** package, or semiparametric models to be implemented with [coxph](#) and the **mstate** package.

**Usage**

```
msm2Surv(data, subject, time, state, covs, Q)
```

**Arguments**

data	Data frame in the format expected by a <a href="#">msm</a> model fit with <code>exacttimes=TRUE</code> or all <code>obstype=2</code> . Each row represents an observation of a state, and the time variable contains the exact and complete transition times of the underlying process. This is explained in more detail in the help page for <a href="#">msm</a> , section <code>obstype=2</code> .
subject	Name of the subject ID in the data (character format, i.e. quoted).
time	Name of the time variable in the data (character).
state	Name of the state variable in the data (character).
covs	Vector of covariate names to carry through (character). If not supplied, this is taken to be all remaining variables in the data.
Q	Transition intensity matrix. This should have number of rows and number of columns both equal to the number of states. If an instantaneous transition is not allowed from state $r$ to state $s$ , then Q should have $(r, s)$ entry 0, otherwise it should be non-zero. The diagonal entries are ignored.

**Details**

For example, if the data supplied to [msm](#) look like this:

subj	days	status	age	treat
1	0	1	66	1
1	27	2	66	1
1	75	3	66	1



1	97	4	66	1
1	1106	4	69	1
2	0	1	49	0
2	90	2	49	0
2	1037	2	51	0

then the output of `msm2Surv` will be a data frame looking like this:

id	from	to	Tstart	Tstop	time	status	age	treat	trans
1	1	2	0	27	27	1	66	1	1
1	1	4	0	27	27	0	66	1	2
1	2	3	27	75	48	1	66	1	3
1	2	4	27	75	48	0	66	1	4
1	3	4	75	97	22	1	69	1	5
2	1	2	0	90	90	1	49	0	1
2	1	4	0	90	90	0	49	0	2
2	2	3	90	1037	947	0	49	0	3
2	2	4	90	1037	947	0	49	0	4

At 27 days, subject 1 is observed to move from state 1 to state 2 (first row, status 1), which means that their potential transition from state 1 to state 4 is censored (second row, status 0).

See the `mstate` package and the references below for more details of this data format and using it for semi-parametric multi-state modelling.

### Value

A data frame of class "msdata", with rows representing observed or censored transitions. There will be one row for each observed transition in the original data, and additional rows for every potential transition that could have occurred out of each observed state.

The data frame will have columns called:

id	Subject ID
from	Starting state of the transition
to	Finishing state of the transition
Tstart	The starting time of the transition
Tstop	The finishing time of the transition
time	The time difference = Tstop - Tstart
status	Event or censoring indicator, with 1 indicating an observed transition, and 0 indicating censoring
trans	Transition number

and any remaining columns will represent covariates. Any covariates whose names clash with the standard variables in the returned data ("id", "from", "to", "Tstart", "Tstop", "time", "status" or "trans") have ".2" appended to their names.

The transition matrix in **mstate** format is stored in the `trans` attribute of the returned object. See the example code below.

### Author(s)

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### References

Putter H, Fiocco M, Geskus RB (2007). Tutorial in biostatistics: Competing risks and multi-state models. *Statistics in Medicine* 26: 2389-2430.

Liesbeth C. de Wreede, Marta Fiocco, Hein Putter (2011). **mstate**: An R Package for the Analysis of Competing Risks and Multi-State Models. *Journal of Statistical Software*, 38(7), 1-30. <http://www.jstatsoft.org/v38/i07>

Jackson, C. H. (2014). flexsurv: Flexible parametric survival and multi-state models. R package version 0.5.

### See Also

[msprep](#), in **mstate**, which produces data in a similar format, given data in "wide" format with one row per subject.

### Examples

```
msmdat <- data.frame(
  subj = c(1, 1, 1, 1, 1, 2, 2, 2),
  days = c(0, 27, 75, 97, 1106, 0, 90, 1037),
  status = c(1, 2, 3, 4, 4, 1, 2, 2),
  age = c(66, 66, 66, 66, 69, 49, 49, 51),
  treat = c(1, 1, 1, 1, 1, 0, 0, 0)
)
# transitions only allowed to next state up or state 4
Q <- rbind(c(1, 1, 0, 1),
           c(0, 1, 1, 1),
           c(0, 0, 1, 1),
           c(0, 0, 0, 0))
dat <- msm2Surv(data=msmdat, subject="subj", time="days", state="status",
               Q=Q)
dat
attr(dat, "trans")
```

---

odds.msm

*Calculate tables of odds ratios for covariates on misclassification probabilities*

---

### Description

Odds ratios are computed by exponentiating the estimated covariate effects on the logit-misclassification probabilities.

**Usage**

```
odds.msm(x, odds.scale = 1, cl = 0.95)
```

**Arguments**

**x** Output from [msm](#) representing a fitted multi-state model.

**odds.scale** Vector with same elements as number of covariates on misclassification probabilities. Corresponds to the increase in each covariate used to calculate its odds ratio. Defaults to all 1.

**cl** Width of the symmetric confidence interval to present. Defaults to 0.95.

**Value**

A list of tables containing odds ratio estimates, one table for each covariate. Each table has three columns, containing the odds ratio, and an approximate upper 95% and lower 95% confidence limit respectively (assuming normality on the log scale), for each misclassification probability.

**Author(s)**

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**See Also**

[msm](#), [hazard.msm](#)

---

paramdata.object

*Developer documentation: internal msm parameters object*

---

**Description**

An object giving information about the parameters of the multi-state model. Used internally during maximum likelihood estimation and arranging results. Returned as the `paramdata` component of a fitted [msm](#) model object.

**Value**

**inits** Vector of initial values for distinct parameters which are being estimated. These have been transformed to the real line (e.g. by log), and exclude parameters being fixed at their initial values, parameters defined to be always fixed (e.g. binomial denominators) and parameters constrained to equal previous ones.

**plabs** Names of parameters in `allinits`.

**allinits** Vector of parameter values before estimation, including those which are fixed or constrained to equal other parameters, and transformed to the real line.

**hmpars** Indices of `allinits` which represent baseline parameters of hidden Markov outcome models (thus excluding covariate effects in HMMs and initial state occupancy probabilities).

<code>fixed</code>	TRUE if all parameters are fixed, FALSE otherwise.
<code>fixedpars</code>	Indices of parameters in <code>allinits</code> which are fixed, either by definition or as requested by the user in the <code>fixedpars</code> argument to <code>msm</code> . Excludes parameters fixed by constraining to equal other parameters.
<code>notfixed</code>	Indices of parameters which are not fixed by the definition of <code>fixedpars</code> .
<code>optpars</code>	Indices of parameters in <code>allinits</code> being estimated, thus those included in <code>inits</code> .
<code>auxpars</code>	Indices of "auxiliary" parameters which are always fixed, for example, binomial denominators ( <code>hmmBinom</code> ) and the <code>which</code> parameter in <code>hmmIdent</code> .
<code>constr</code>	Vector of integers, of length <code>npars</code> , indicating which sets of parameters are constrained to be equal to each other. If two of these integers are equal the corresponding parameters are equal. A negative element indicates that parameter is defined to be minus some other parameter (this is used for covariate effects on transition intensities).
<code>npars</code>	Total number of parameters, equal to <code>length(allinits)</code> .
<code>nfix</code>	Number of fixed parameters, equal to <code>length(fixedpars)</code> .
<code>nopt</code>	Number of parameters being estimated, equal to <code>length(inits)</code> and <code>length(optpars)</code> .
<code>ndup</code>	Number of parameters defined as duplicates of previous parameters by equality constraints (currently unused).
<code>ranges</code>	Matrix of defined ranges for each parameter on the natural scale (e.g. 0 to infinity for rate parameters).
<code>opt</code>	Object returned by the optimisation routine (such as <code>optim</code> ).
<code>foundse</code>	TRUE if standard errors are available after optimisation. If FALSE the optimisation probably hasn't converged.
<code>lik</code>	Minus twice the log likelihood at the parameter estimates.
<code>deriv</code>	Derivatives of the minus twice log likelihood at the parameter estimates, if available.
<code>information</code>	Corresponding expected information matrix at the parameter estimates, if available.
<code>params</code>	Vector of parameter values after maximum likelihood estimation, corresponding to <code>allinits</code> , still on the real-line transformed scale.
<code>covmat</code>	Covariance matrix corresponding to <code>params</code> .
<code>ci</code>	Matrix of confidence intervals corresponding to <code>params</code> , with nominal coverage (default 0.95) defined by the <code>cl</code> argument of <code>msm</code> .
<code>estimates.t</code>	Vector of parameter estimates, as <code>params</code> but with parameters on their natural scales.

**See Also**

[msm.object](#)

---

pearson.msm	<i>Pearson-type goodness-of-fit test</i>
-------------	--

---

## Description

Pearson-type goodness-of-fit test for multi-state models fitted to panel-observed data.

## Usage

```
pearson.msm(x, transitions=NULL, timegroups=3, intervalgroups=3,
            covgroups=3, groups=NULL, boot=FALSE, B=500,
            next.obstime=NULL, N=100, indep.cens=TRUE,
            maxtimes=NULL, pval=TRUE)
```

## Arguments

- |                |   |
|----------------|---|
| x              | A fitted multi-state model, as returned by <a href="#">msm</a> .  |
| transitions    | <p>This should be an integer vector indicating which interval transitions should be grouped together in the contingency table. Its length should be the number of allowed interval transitions, excluding transitions from absorbing states to absorbing states.</p> <p>The allowed interval transitions are the set of pairs of states <math>(a, b)</math> for which it is possible to observe <math>a</math> at one time and <math>b</math> at any later time. For example, in a "well-disease-death" model with allowed <i>instantaneous</i> 1-2, 2-3 transitions, there are 5 allowed <i>interval</i> transitions. In numerical order, these are 1-1, 1-2, 1-3, 2-2 and 2-3, excluding absorbing-absorbing transitions.</p> <p>Then, to group transitions 1-1,1-2 together, and transitions 2-2,2-3 together, specify</p> <pre>transitions = c(1, 1, 2, 3, 3).</pre> <p>Only transitions from the same state may be grouped. By default, each interval transition forms a separate group.</p> |
| timegroups     | Number of groups based on quantiles of the time since the start of the process.   |
| intervalgroups | Number of groups based on quantiles of the time interval between observations, within time groups   |
| covgroups      | <p>Number of groups based on quantiles of <math>\sum_r q_{irr}</math>, where <math>q_{irr}</math> are the diagonal entries of the transition intensity matrix for the <math>i</math>th transition. These are a function of the covariate effects and the covariate values at the <math>i</math>th transition: <math>q_{irr}</math> is minus the sum of the off-diagonal entries <math>q_{rs}^{(0)} \exp(\beta_{rs}^T z_i)</math> on the <math>r</math>th row.</p> <p>Thus covgroups summarises the impact of covariates at each observation, by calculating the overall rate of progression through states at that observation.</p> <p>For time-inhomogeneous models specified using the <code>pci</code> argument to <a href="#">msm</a>, if the only covariate is the time period, covgroups is set to 1, since timegroups ensures that transitions are grouped by time.</p>  |

groups	<p>A vector of arbitrary groups in which to categorise each transition. This can be an integer vector or a factor. This can be used to diagnose specific areas of poor fit. For example, the contingency table might be grouped by arbitrary combinations of covariates to detect types of individual for whom the model fits poorly.</p> <p>The length of groups should be <math>x\\$data\\$n</math>, the number of observations used in the model fit, which is the number of observations in the original dataset with any missing values excluded. The value of groups at observation <math>i</math> is used to categorise the transition which <i>ends</i> at observation <math>i</math>. Values of groups at the first observation for each subject are ignored.</p>
boot	<p>Estimate an "exact" p-value using a parametric bootstrap.</p> <p>All objects used in the original call to <code>msm</code> which produced <code>x</code>, such as the <code>qmatrix</code>, should be in the working environment, or else an "object not found" error will be given. This enables the original model to be refitted to the replicate datasets. Note that groups cannot be used with bootstrapping, as the simulated observations will not be in the same categories as the original observations.</p>
B	Number of bootstrap replicates.
next.obstime	<p>This is a vector of length <math>x\\$data\\$n</math> (the number of observations used in the model fit) giving the time to the next <i>scheduled</i> observation following each time point. This is only used when times to death are known exactly.</p> <p>For individuals who died (entered an absorbing state) before the next scheduled observation, and the time of death is known exactly, <code>next.obstime</code> would be <i>greater</i> than the observed death time.</p> <p>If the individual did not die, and a scheduled observation did follow that time point, <code>next.obstime</code> should just be the same as the time to that observation.</p> <p><code>next.obstime</code> is used to determine a grouping of the time interval between observations, which should be based on scheduled observations. If exact times to death were used in the grouping, then shorter intervals would contain excess deaths, and the goodness-of-fit statistic would be biased.</p> <p>If <code>next.obstime</code> is unknown, it is multiply-imputed using a product-limit estimate based on the intervals to observations other than deaths. The resulting tables of transitions are averaged over these imputations. This may be slow.</p>
N	Number of imputations for the estimation of the distribution of the next scheduled observation time, when there are exact death times.
indep.cens	If TRUE, then times to censoring are included in the estimation of the distribution to the next scheduled observation time. If FALSE, times to censoring are assumed to be systematically different from other observation times.
maxtimes	A vector of length $x\$data\$n$ , or a common scalar, giving an upper bound for the next scheduled observation time. Used in the multiple imputation when times to death are known exactly. If a value greater than <code>maxtimes</code> is simulated, then the next scheduled observation is taken as censored. This should be supplied, if known. If not supplied, this is taken to be the maximum interval occurring in the data, plus one time unit. For observations which are not exact death times, this should be the time since the previous observation.
pval	Calculate a p-value using the improved approximation of Titman (2009). This is optional since it is not needed during bootstrapping, and it is computationally

non-trivial. Only available currently for non-hidden Markov models for panel data without exact death times. Also not available for models with censoring, including time-homogeneous models fitted with the `pci` option to `msm`.

## Details

This method (Aguirre-Hernandez and Farewell, 2002) is intended for data which represent observations of the process at arbitrary times ("snapshots", or "panel-observed" data). For data which represent the exact transition times of the process, `prevalence.msm` can be used to assess fit, though without a formal test.

When times of death are known exactly, states are misclassified, or an individual's final observation is a censored state, the modification by Titman and Sharples (2008) is used. The only form of censoring supported is a state at the end of an individual's series which represents an unknown transient state (i.e. the individual is only known to be alive at this time). Other types of censoring are omitted from the data before performing the test.

See the references for further details of the methods. The method used for censored states is a modification of the method in the appendix to Titman and Sharples (2008), described at <http://www.mrc-bsu.cam.ac.uk/wp-content/uploads/robustcensoring.pdf> (Titman, 2007).

Groupings of the time since initiation, the time interval and the impact of covariates are based on equally-spaced quantiles. The number of groups should be chosen that there are not many cells with small expected numbers of transitions, since the deviance statistic will be unstable for sparse contingency tables. Ideally, the expected numbers of transitions in each cell of the table should be no less than about 5. Conversely, the power of the test is reduced if there are too few groups. Therefore, some sensitivity analysis of the test results to the grouping is advisable.

Saved model objects fitted with previous versions of R (versions less than 1.2) will need to be refitted under the current R for use with `pearson.msm`.

## Value

A list whose first two elements are contingency tables of observed transitions  $O$  and expected transitions  $E$ , respectively, for each combination of groups. The third element is a table of the deviances  $(O - E)^2/E$  multiplied by the sign of  $O - E$ . If the expected number of transitions is zero then the deviance is zero. Entries in the third matrix will be bigger in magnitude for groups for which the model fits poorly.

"test" the fourth element of the list, is a data frame with one row containing the Pearson-type goodness-of-fit test statistic `stat`. The test statistic is the sum of the deviances. For panel-observed data without exact death times, misclassification or censored observations, `p` is the p-value for the test statistic calculated using the improved approximation of Titman (2009).  
For these models, for comparison with older versions of the package, `test` also presents `p.lower` and `p.upper`, which are theoretical lower and upper limits for the p-value of the test statistic, based on  $\chi^2$  distributions with `df.lower` and `df.upper` degrees of freedom, respectively. `df.upper` is the number of independent cells in the contingency table, and `df.lower` is `df.upper` minus the number of estimated parameters in the model.

"intervalq"	(not printed by default) contains the definition of the grouping of the intervals between observations. These groups are defined by quantiles within the groups corresponding to the time since the start of the process.
"sim"	If there are exact death times, this contains simulations of the contingency tables and test statistics for each imputation of the next scheduled sampling time. These are averaged over to produce the presented tables and test statistic. This element is not printed by default.  With exact death times, the null variance of the test statistic (formed by taking mean of simulated test statistics) is less than twice the mean (Titman, 2008), and the null distribution is not $\chi^2$ . In this case, p.upper is an upper limit for the true asymptotic p-value, but p.lower is not a lower limit, and is not presented.
"boot"	If the bootstrap has been used, the element will contain the bootstrap replicates of the test statistics (not printed by default).
"lambda"	If the Titman (2009) p-value has been calculated, this contains the weights defining the null distribution of the test statistic as a weighted sum of $\chi_1^2$ random variables (not printed by default).

**Author(s)**

Andrew Titman <a.titman@lancaster.ac.uk>, Chris Jackson <chris.jackson@mrc-bsu.cam.ac.uk>

**References**

- Aguirre-Hernandez, R. and Farewell, V. (2002) A Pearson-type goodness-of-fit test for stationary and time-continuous Markov regression models. *Statistics in Medicine* 21:1899-1911.
- Titman, A. and Sharples, L. (2008) A general goodness-of-fit test for Markov and hidden Markov models. *Statistics in Medicine* 27(12):2177-2195
- Titman, A. (2009) Computation of the asymptotic null distribution of goodness-of-fit tests for multi-state models. *Lifetime Data Analysis* 15(4):519-533.
- Titman, A. (2008) Model diagnostics in multi-state models of biological systems. PhD thesis, University of Cambridge.

**See Also**

[msm](#), [prevalence.msm](#), [scorereresid.msm](#),

**Examples**

```
psor.q <- rbind(c(0,0.1,0,0),c(0,0,0.1,0),c(0,0,0,0.1),c(0,0,0,0))
psor.msm <- msm(state ~ months, subject=ptnum, data=psor,
               qmatrix = psor.q, covariates = ~ollwsdrt+hieffusn,
               constraint = list(hieffusn=c(1,1,1),ollwsdrt=c(1,1,2)))
pearson.msm(psor.msm, timegroups=2, intervalgroups=2, covgroups=2)
# More 1-2, 1-3 and 1-4 observations than expected in shorter time
# intervals - the model fits poorly.
# A random effects model might accommodate such fast progressors.
```



pexp

*Exponential distribution with piecewise-constant rate***Description**

Density, distribution function, quantile function and random generation for a generalisation of the exponential distribution, in which the rate changes at a series of times.

**Usage**

```
dpexp(x, rate=1, t=0, log = FALSE)
ppexp(q, rate=1, t=0, lower.tail = TRUE, log.p = FALSE)
qpexp(p, rate=1, t=0, lower.tail = TRUE, log.p = FALSE)
rpexp(n, rate=1, t=0)
```

**Arguments**

x, q	vector of quantiles.
p	vector of probabilities.
n	number of observations. If <code>length(n) &gt; 1</code> , the length is taken to be the number required.
rate	vector of rates.
t	vector of the same length as <code>rate</code> , giving the times at which the rate changes. The first element of <code>t</code> should be 0, and <code>t</code> should be in increasing order.
log, log.p	logical; if TRUE, probabilities <code>p</code> are given as <code>log(p)</code> , or log density is returned.
lower.tail	logical; if TRUE (default), probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .

**Details**

Consider the exponential distribution with rates  $r_1, \dots, r_n$  changing at times  $t_1, \dots, t_n$ , with  $t_1 = 0$ . Suppose  $t_k$  is the maximum  $t_i$  such that  $t_i < x$ . The density of this distribution at  $x > 0$  is  $f(x)$  for  $k = 1$ , and

$$\prod_{i=1}^k (1 - F(t_i - t_{i-1}, r_i)) f(x - t_k, r_k)$$

for  $k > 1$ .

where  $F()$  and  $f()$  are the distribution and density functions of the standard exponential distribution.

If `rate` is of length 1, this is just the standard exponential distribution. Therefore, for example, `dpexp(x)`, with no other arguments, is simply equivalent to `dexp(x)`.

Only `rpexp` is used in the `msm` package, to simulate from Markov processes with piecewise-constant intensities depending on time-dependent covariates. These functions are merely provided for completion, and are not optimized for numerical stability or speed.

**Value**

dpexp gives the density, ppexp gives the distribution function, qpexp gives the quantile function, and rpexp generates random deviates.

**Author(s)**

C. H. Jackson <chris.jackson@mrc-bsu.cam.ac.uk>

**See Also**

[dexp](#), [sim.msm](#).

**Examples**

```
x <- seq(0.1, 50, by=0.1)
rate <- c(0.1, 0.2, 0.05, 0.3)
t <- c(0, 10, 20, 30)
## standard exponential distribution
plot(x, dexp(x, 0.1), type="l")
## distribution with piecewise constant rate
lines(x, dpexp(x, rate, t), type="l", lty=2)
## standard exponential distribution
plot(x, pexp(x, 0.1), type="l")
## distribution with piecewise constant rate
lines(x, ppexp(x, rate, t), type="l", lty=2)
```

---

phasemeans.msm

*Parameters of phase-type models in mixture form*

---

**Description**

Parameters of fitted two-phase models, in mixture model parameterisation.

**Usage**

```
phasemeans.msm(x, covariates="mean", ci=c("none", "normal", "bootstrap"),
               cl=0.95, B=1000, cores=NULL)
```

**Arguments**

x	A fitted multi-state model, as returned by <a href="#">msm</a> .
covariates	Covariate values, see <a href="#">qmatrix.msm</a> .
ci	If "none" (the default) no confidence intervals are calculated. Otherwise "normal", or "boot" as described by <a href="#">qmatrix.msm</a> .
cl	Width of the symmetric confidence interval, relative to 1.
B	Number of bootstrap replicates, or number of normal simulations from the distribution of the MLEs.
cores	Number of cores to use for bootstrapping using parallel processing. See <a href="#">boot.msm</a> for more details.

**Value**

Matrix with one row for each state that has a two-phase distribution, and three columns: the short-stay mean, long-stay mean and long-stay probability. These are functions of the transition intensities of the expanded hidden Markov model, defined in [d2phase](#).

**Author(s)**

C. H. Jackson <chris.jackson@mrc-bsu.cam.ac.uk>

**See Also**

[d2phase](#).

---

plot.msm

*Plots of multi-state models*

---

**Description**

This produces a plot of the expected probability of survival against time, from each transient state. Survival is defined as not entering an absorbing state.

**Usage**

```
## S3 method for class 'msm'
plot(x, from, to, range, covariates, legend.pos, xlab="Time",
      ylab="Fitted survival probability", lwd=1, ...)
```

**Arguments**

x	Output from <a href="#">msm</a> , representing a fitted multi-state model object.
from	States from which to consider survival. Defaults to the complete set of transient states.
to	Absorbing state to consider. Defaults to the highest-labelled absorbing state.
range	Vector of two elements, giving the range of times to plot for.
covariates	Covariate values for which to evaluate the expected probabilities. This can either be: <ul style="list-style-type: none"> <li>the string "mean", denoting the means of the covariates in the data (this is the default),</li> <li>the number 0, indicating that all the covariates should be set to zero,</li> <li>or a list of values, with optional names. For example <pre>list(60, 1)</pre>           where the order of the list follows the order of the covariates originally given in the model formula, or a named list, <pre>list(age = 60, sex = 1)</pre> </li> </ul>

legend.pos	Vector of the $x$ and $y$ position, respectively, of the legend.
xlab	x axis label.
ylab	y axis label.
lwd	Line width. See <a href="#">par</a> .
...	Other arguments to be passed to the generic <a href="#">plot</a> and <a href="#">lines</a> functions.

**Author(s)**

C. H. Jackson <chris.jackson@mrc-bsu.cam.ac.uk>

**See Also**

[msm](#)

---

plot.prevalence.msm *Plot of observed and expected prevalences*

---

**Description**

Provides a rough indication of goodness of fit of a multi-state model, by estimating the observed numbers of individuals occupying a state at a series of times, and plotting these against forecasts from the fitted model, for each state. Observed prevalences are indicated as solid lines, expected prevalences as dashed lines.

**Usage**

```
## S3 method for class 'prevalence.msm'
plot(x, mintime=NULL, maxtime=NULL, timezero=NULL,
      initstates=NULL, interp=c("start","midpoint"),
      censtime=Inf, subset=NULL,
      covariates="population", misccovariates="mean",
      piecewise.times=NULL, piecewise.covariates=NULL,
      xlab="Times",ylab="Prevalence (%)", lwd.obs=1,
      lwd.exp=1, lty.obs=1, lty.exp=2, col.obs="blue",
      col.exp="red", legend.pos=NULL,
      ...)
```

**Arguments**

x	A fitted multi-state model produced by <a href="#">msm</a> .
mintime	Minimum time at which to compute the observed and expected prevalences of states.
maxtime	Maximum time at which to compute the observed and expected prevalences of states.
timezero	Initial time of the Markov process. Expected values are forecasted from here. Defaults to the minimum of the observation times given in the data.

initstates	Optional vector of the same length as the number of states. Gives the numbers of individuals occupying each state at the initial time, to be used for forecasting expected prevalences. The default is those observed in the data. These should add up to the actual number of people in the study at the start.
interp	Interpolation method for observed states, see <a href="#">prevalence.msm</a> .
censtime	Subject-specific maximum follow-up times, see <a href="#">prevalence.msm</a> .
subset	Vector of the subject identifiers to calculated observed prevalences for.
covariates	Covariate values for which to forecast expected state occupancy. See <a href="#">prevalence.msm</a> — if this function runs too slowly, as it may if there are continuous covariates, replace covariates="population" with covariates="mean".
misccovariates	(Misclassification models only) Values of covariates on the misclassification probability matrix. See <a href="#">prevalence.msm</a> .
piecewise.times	Times at which piecewise-constant intensities change. See <a href="#">prevalence.msm</a> .
piecewise.covariates	Covariates on which the piecewise-constant intensities depend. See <a href="#">prevalence.msm</a> .
xlab	x axis label.
ylab	y axis label.
lwd.obs	Line width for observed prevalences. See <a href="#">par</a> .
lwd.exp	Line width for expected prevalences. See <a href="#">par</a> .
lty.obs	Line type for observed prevalences. See <a href="#">par</a> .
lty.exp	Line type for expected prevalences. See <a href="#">par</a> .
col.obs	Line colour for observed prevalences. See <a href="#">par</a> .
col.exp	Line colour for expected prevalences. See <a href="#">par</a> .
legend.pos	Vector of the $x$ and $y$ position, respectively, of the legend.
...	Further arguments to be passed to the generic <a href="#">plot</a> function.

### Details

See [prevalence.msm](#) for details of the assumptions underlying this method.

Observed prevalences are plotted with a solid line, and expected prevalences with a dotted line.

### References

Gentleman, R.C., Lawless, J.F., Lindsey, J.C. and Yan, P. Multi-state Markov models for analysing incomplete disease history data with illustrations for HIV disease. *Statistics in Medicine* (1994) 13(3): 805–821.

### See Also

[prevalence.msm](#)

---

plot.survfit.msm      *Plot empirical and fitted survival curves*

---

### Description

Plot a Kaplan-Meier estimate of the survival probability and compare it with the fitted survival probability from a `msm` model.

### Usage

```
## S3 method for class 'survfit.msm'
plot(x, from=1, to=NULL, range=NULL, covariates="mean",
      interp=c("start", "midpoint"), ci=c("none", "normal", "bootstrap"), B=100,
      legend.pos=NULL, xlab="Time", ylab="Survival probability",
      lty=1, lwd=1, col="red", lty.ci=2, lwd.ci=1, col.ci="red",
      mark.time=TRUE, col.surv="blue", lty.surv=2, lwd.surv=1,
      survdata=FALSE,
      ...)
```

### Arguments

<code>x</code>	Output from <code>msm</code> , representing a fitted multi-state model object.
<code>from</code>	Non-absorbing state from which to consider survival. Defaults to state 1. The fitted probabilities will then be calculated as the transition probabilities from this state to <code>to</code> . The empirical survival curve plots survival from the first observation of <code>from</code> (where this exists) to the first entry time into <code>to</code> .
<code>to</code>	Absorbing state to consider. Defaults to the highest-labelled absorbing state.
<code>range</code>	Vector of two elements, giving the range of times to plot for.
<code>covariates</code>	Covariate values for which to evaluate the expected probabilities. This can either be: <ul style="list-style-type: none"> <li>the string "mean", denoting the means of the covariates in the data (this is the default),</li> <li>the number 0, indicating that all the covariates should be set to zero,</li> <li>or a list of values, with optional names. For example <pre>list(60, 1)</pre>           where the order of the list follows the order of the covariates originally given in the model formula, or a named list, <pre>list(age = 60, sex = 1)</pre>           but note the empirical curve is plotted for the full population. To consider subsets for the empirical curve, set <code>survdata=TRUE</code> to extract the survival data and build a survival plot by hand using <code>plot.survfit</code>.</li> </ul>

ci	If "none" (the default) no confidence intervals are plotted. If "normal" or "bootstrap", confidence intervals are plotted based on the respective method in <a href="#">pmatrix.msm</a> . This is very computationally-intensive, since intervals must be computed at a series of times.
B	Number of bootstrap or normal replicates for the confidence interval. The default is 100 rather than the usual 1000, since these plots are for rough diagnostic purposes.
interp	If <code>interp="start"</code> (the default) then the entry time into the absorbing state is assumed to be the time it is first observed in the data. If <code>interp="midpoint"</code> then the entry time into the absorbing state is assumed to be halfway between the time it is first observed and the previous observation time. This is generally more reasonable for "progressive" models with observations at arbitrary times.
legend.pos	Vector of the $x$ and $y$ position, respectively, of the legend.
xlab	x axis label.
ylab	y axis label.
lty	Line type for the fitted curve. See <a href="#">par</a> .
lwd	Line width for the fitted curve. See <a href="#">par</a> .
col	Colour for the fitted curve. See <a href="#">par</a> .
lty.ci	Line type for the fitted curve confidence limits. See <a href="#">par</a> .
lwd.ci	Line width for the fitted curve confidence limits. See <a href="#">par</a> .
col.ci	Colour for the fitted curve confidence limits. See <a href="#">par</a> .
mark.time	Mark the empirical survival curve at each censoring point, see <a href="#">lines.survfit</a> .
col.surv	Colour for the empirical survival curve, passed to <a href="#">lines.survfit</a> . See <a href="#">par</a> .
lty.surv	Line type for the empirical survival curve, passed to <a href="#">lines.survfit</a> . See <a href="#">par</a> .
lwd.surv	Line width for the empirical survival curve, passed to <a href="#">lines.survfit</a> . See <a href="#">par</a> .
survdata	Set to TRUE to return the survival data frame constructed when plotting the empirical curve. This can be used for constructing survival plots by hand using <a href="#">plot.survfit</a> .
...	Other arguments to be passed to the <a href="#">plot</a> function which draws the fitted curve, or the <a href="#">lines.survfit</a> function which draws the empirical curve.

## Details

If the data represent observations of the process at arbitrary times, then the first occurrence of the absorbing state in the data will usually be greater than the actual first transition time to that state. Therefore the Kaplan-Meier estimate of the survival probability will be an overestimate.

The method of Turnbull (1976) could be used to give a non-parametric estimate of the time to an interval-censored event, and compared to the equivalent estimate from a multi-state model. This is implemented in the CRAN package **interval** (Fay and Shaw 2010).

This currently only handles time-homogeneous models.

## References

Turnbull, B. W. (1976) The empirical distribution function with arbitrarily grouped, censored and truncated data. *J. R. Statist. Soc. B* 38, 290-295.

Fay, MP and Shaw, PA (2010). Exact and Asymptotic Weighted Logrank Tests for Interval Censored Data: The interval R package. *Journal of Statistical Software*. <http://www.jstatsoft.org/v36/i02/>. 36 (2):1-34.

## See Also

[survfit](#), [plot.survfit](#), [plot.prevalence.msm](#)

---

plotprog.msm

*Kaplan Meier estimates of incidence*

---

## Description

Compute and plot Kaplan-Meier estimates of the probability that each successive state has not occurred yet.

## Usage

```
plotprog.msm(formula, subject, data, legend.pos=NULL, xlab="Time",
             ylab="1 - incidence probability", lwd=1, xlim=NULL,
             mark.time=TRUE, ...)
```

## Arguments

formula	A formula giving the vectors containing the observed states and the corresponding observation times. For example, state ~ time Observed states should be in the set $1, \dots, n$ , where $n$ is the number of states.
subject	Vector of subject identification numbers for the data specified by formula. If missing, then all observations are assumed to be on the same subject. These must be sorted so that all observations on the same subject are adjacent.
data	An optional data frame in which the variables represented by state, time and subject can be found.
legend.pos	Vector of the $x$ and $y$ position, respectively, of the legend.
xlab	x axis label.
ylab	y axis label.
lwd	Line width. See <a href="#">par</a> .
xlim	x axis limits, e.g. <code>c(0,10)</code> for an axis ranging from 0 to 10. Default is the range of observation times.
mark.time	Mark the empirical survival curve at each censoring point, see <a href="#">lines.survfit</a> .
...	Other arguments to be passed to the <a href="#">plot</a> and <a href="#">lines.survfit</a> functions.



**Details**

If the data represent observations of the process at arbitrary times, then the first occurrence of the state in the data will usually be greater than the actual first transition time to that state. Therefore the probabilities plotted by [plotprog.msm](#) will be overestimates.

**See Also**

[survfit](#), [plot.survfit](#)

---

pmatrix.msm

*Transition probability matrix*

---

**Description**

Extract the estimated transition probability matrix from a fitted multi-state model for a given time interval, at a given set of covariate values.

**Usage**

```
pmatrix.msm(x=NULL, t=1, t1=0, covariates="mean",
            ci=c("none", "normal", "bootstrap"), cl=0.95, B=1000,
            cores=NULL, qmatrix=NULL,
            ...)
```

**Arguments**

x	A fitted multi-state model, as returned by <a href="#">msm</a> .
t	The time interval to estimate the transition probabilities for, by default one unit.
t1	The starting time of the interval. Used for models x with piecewise-constant intensities fitted using the <code>pci</code> option to <a href="#">msm</a> . The probabilities will be computed on the interval [t1, t1+t].
covariates	The covariate values at which to estimate the transition probabilities. This can either be: <ul style="list-style-type: none"> <li>the string "mean", denoting the means of the covariates in the data (this is the default),</li> <li>the number 0, indicating that all the covariates should be set to zero,</li> <li>or a list of values, with optional names. For example <pre>list(60, 1)</pre> where the order of the list follows the order of the covariates originally given in the model formula, or a named list, <pre>list(age = 60, sex = 1)</pre> If some covariates are specified but not others, the missing ones default to zero.</li> </ul>

	For time-inhomogeneous models fitted using the <code>pci</code> option to <code>msm</code> , "covariates" here include only those specified using the <code>covariates</code> argument to <code>msm</code> , and exclude the artificial covariates representing the time period.
	For time-inhomogeneous models fitted "by hand" by using a time-dependent covariate in the <code>covariates</code> argument to <code>msm</code> , the function <code>pmatrix.piecewise.msm</code> should be used to calculate transition probabilities.
<code>ci</code>	<p>If "normal", then calculate a confidence interval for the transition probabilities by simulating <code>B</code> random vectors from the asymptotic multivariate normal distribution implied by the maximum likelihood estimates (and covariance matrix) of the log transition intensities and covariate effects, then calculating the resulting transition probability matrix for each replicate. See, e.g. Mandel (2013) for a discussion of this approach.</p> <p>If "bootstrap" then calculate a confidence interval by non-parametric bootstrap refitting. This is 1-2 orders of magnitude slower than the "normal" method, but is expected to be more accurate. See <code>boot.msm</code> for more details of bootstrapping in <code>msm</code>.</p> <p>If "none" (the default) then no confidence interval is calculated.</p>
<code>c1</code>	Width of the symmetric confidence interval, relative to 1.
<code>B</code>	Number of bootstrap replicates, or number of normal simulations from the distribution of the MLEs
<code>cores</code>	Number of cores to use for bootstrapping using parallel processing. See <code>boot.msm</code> for more details.
<code>qmatrix</code>	A transition intensity matrix. Either this or a fitted model <code>x</code> must be supplied. No confidence intervals are available if <code>qmatrix</code> is supplied.
<code>...</code>	Optional arguments to be passed to <code>MatrixExp</code> to control the method of computing the matrix exponential.

## Details

For a continuous-time homogeneous Markov process with transition intensity matrix  $Q$ , the probability of occupying state  $s$  at time  $u + t$  conditionally on occupying state  $r$  at time  $u$  is given by the  $(r, s)$  entry of the matrix  $P(t) = \exp(tQ)$ , where  $\exp()$  is the matrix exponential.

For non-homogeneous processes, where covariates and hence the transition intensity matrix  $Q$  are piecewise-constant in time, the transition probability matrix is calculated as a product of matrices over a series of intervals, as explained in `pmatrix.piecewise.msm`.

The `pmatrix.piecewise.msm` function is only necessary for models fitted using a time-dependent covariate in the `covariates` argument to `msm`. For time-inhomogeneous models fitted using "pci", `pmatrix.msm` can be used, with arguments `t` and `t1`, to calculate transition probabilities over any time period.

## Value

The matrix of estimated transition probabilities  $P(t)$  in the given time. Rows correspond to "from-state" and columns to "to-state".

Or if `ci="normal"` or `ci="bootstrap"`, `pmatrix.msm` returns a list with components `estimates` and `ci`, where `estimates` is the matrix of estimated transition probabilities, and `ci` is a list of two matrices containing the upper and lower confidence limits.

**Author(s)**

C. H. Jackson <chris.jackson@mrc-bsu.cam.ac.uk>.

**References**

Mandel, M. (2013). "Simulation based confidence intervals for functions with complicated derivatives." *The American Statistician* 67(2):76-81

**See Also**

[qmatrix.msm](#), [pmatrix.piecewise.msm](#), [boot.msm](#)

---

`pmatrix.piecewise.msm` *Transition probability matrix for processes with piecewise-constant intensities*

---

**Description**

Extract the estimated transition probability matrix from a fitted non-time-homogeneous multi-state model for a given time interval. This is a generalisation of [pmatrix.msm](#) to models with time-dependent covariates. Note that [pmatrix.msm](#) is sufficient to calculate transition probabilities for time-inhomogeneous models fitted using the `pci` argument to [msm](#).

**Usage**

```
pmatrix.piecewise.msm(x=NULL, t1, t2, times, covariates,
ci=c("none", "normal", "bootstrap"), cl=0.95, B=1000, cores=NULL,
qlist=NULL,...)
```

**Arguments**

<code>x</code>	A fitted multi-state model, as returned by <a href="#">msm</a> . This should be a non-homogeneous model, whose transition intensity matrix depends on a time-dependent covariate.
<code>t1</code>	The start of the time interval to estimate the transition probabilities for.
<code>t2</code>	The end of the time interval to estimate the transition probabilities for.
<code>times</code>	Cut points at which the transition intensity matrix changes.
<code>covariates</code>	A list with number of components one greater than the length of <code>times</code> . Each component of the list is specified in the same way as the <code>covariates</code> argument to <a href="#">pmatrix.msm</a> . The components correspond to the covariate values in the intervals ( <code>t1</code> , <code>times[1]</code> ], ( <code>times[1]</code> , <code>times[2]</code> ], ..., ( <code>times[length(times)]</code> , <code>t2</code> ] (assuming that all elements of <code>times</code> are in the interval ( <code>t1</code> , <code>t2</code> )).

ci	<p>If "normal", then calculate a confidence interval for the transition probabilities by simulating B random vectors from the asymptotic multivariate normal distribution implied by the maximum likelihood estimates (and covariance matrix) of the log transition intensities and covariate effects, then calculating the resulting transition probability matrix for each replicate.</p> <p>If "bootstrap" then calculate a confidence interval by non-parametric bootstrap refitting. This is 1-2 orders of magnitude slower than the "normal" method, but is expected to be more accurate. See <a href="#">boot.msm</a> for more details of bootstrapping in <b>msm</b>.</p> <p>If "none" (the default) then no confidence interval is calculated.</p>
c1	Width of the symmetric confidence interval, relative to 1.
B	Number of bootstrap replicates, or number of normal simulations from the distribution of the MLEs
cores	Number of cores to use for bootstrapping using parallel processing. See <a href="#">boot.msm</a> for more details.
qlist	A list of transition intensity matrices, of length one greater than the length of times. Either this or a fitted model x must be supplied. No confidence intervals are available if (just) qlist is supplied.
...	Optional arguments to be passed to <a href="#">MatrixExp</a> to control the method of computing the matrix exponential.

### Details

Suppose a multi-state model has been fitted, in which the transition intensity matrix  $Q(x(t))$  is modelled in terms of time-dependent covariates  $x(t)$ . The transition probability matrix  $P(t_1, t_n)$  for the time interval  $(t_1, t_n)$  cannot be calculated from the estimated intensity matrix as  $\exp((t_n - t_1)Q)$ , because  $Q$  varies within the interval  $t_1, t_n$ . However, if the covariates are piecewise-constant, or can be approximated as piecewise-constant, then we can calculate  $P(t_1, t_n)$  by multiplying together individual matrices  $P(t_i, t_{i+1}) = \exp((t_{i+1} - t_i)Q)$ , calculated over intervals where  $Q$  is constant:

$$P(t_1, t_n) = P(t_1, t_2)P(t_2, t_3) \dots P(t_{n-1}, t_n)$$

### Value

The matrix of estimated transition probabilities  $P(t)$  for the time interval  $[t_1, t_n]$ . That is, the probabilities of occupying state  $s$  at time  $t_n$  conditionally on occupying state  $r$  at time  $t_1$ . Rows correspond to "from-state" and columns to "to-state".

### Author(s)

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>

### See Also

[pmatrix.msm](#)

## Examples

```
## Not run:
## In a clinical study, suppose patients are given a placebo in the
## first 5 weeks, then they begin treatment 1 at 5 weeks, and
## a combination of treatments 1 and 2 from 10 weeks.
## Suppose a multi-state model x has been fitted for the patients'
## progress, with treat1 and treat2 as time dependent covariates.

## Cut points for when treatment covariate changes
times <- c(0, 5, 10)

## Indicators for which treatments are active in the four intervals
## defined by the three cut points
covariates <- list( list (treat1=0, treat2=0), list (treat1=0, treat2=0), list(treat1=1, treat2=0),
list(treat1=1, treat2=1) )

## Calculate transition probabilities from the start of the study to 15 weeks
pmatrix.pieceswise.msm(x, 0, 15, times, covariates)

## End(Not run)
```

---

pnext.msm

*Probability of each state being next*


---

## Description

Compute a matrix of the probability of each state  $s$  being the next state of the process after each state  $r$ . Together with the mean sojourn times in each state ([sojourn.msm](#)), these fully define a continuous-time Markov model.

## Usage

```
pnext.msm(x, covariates = "mean",
          ci=c("normal","bootstrap","delta","none"), cl = 0.95,
          B=1000, cores=NULL)
```

## Arguments

x	A fitted multi-state model, as returned by <a href="#">msm</a> .
covariates	The covariate values at which to estimate the intensities. This can either be: <ul style="list-style-type: none"> <li>the string "mean", denoting the means of the covariates in the data (this is the default),</li> <li>the number 0, indicating that all the covariates should be set to zero,</li> <li>or a list of values, with optional names. For example</li> </ul>

	<pre>list (60, 1)</pre> <p>where the order of the list follows the order of the covariates originally given in the model formula, or a named list,</p> <pre>list (age = 60, sex = 1)</pre>
ci	<p>If "normal" (the default) then calculate a confidence interval by simulating B random vectors from the asymptotic multivariate normal distribution implied by the maximum likelihood estimates (and covariance matrix) of the log transition intensities and covariate effects, then transforming.</p> <p>If "bootstrap" then calculate a confidence interval by non-parametric bootstrap refitting. This is 1-2 orders of magnitude slower than the "normal" method, but is expected to be more accurate. See <a href="#">boot.msm</a> for more details of bootstrapping in <b>msm</b>.</p> <p>If "delta" then confidence intervals are calculated based on the delta method SEs of the log rates, but this is not recommended since it may not respect the constraint that probabilities are less than one.</p>
c1	Width of the symmetric confidence interval to present. Defaults to 0.95.
B	Number of bootstrap replicates, or number of normal simulations from the distribution of the MLEs.
cores	Number of cores to use for bootstrapping using parallel processing. See <a href="#">boot.msm</a> for more details.

## Details

For a continuous-time Markov process in state  $r$ , the probability that the next state is  $s$  is  $-q_{rs}/q_{rr}$ , where  $q_{rs}$  is the transition intensity ([qmatrix.msm](#)).

A continuous-time Markov model is fully specified by these probabilities together with the mean sojourn times  $-1/q_{rr}$  in each state  $r$ . This gives a more intuitively meaningful description of a model than the intensity matrix.

Remember that **msm** deals with continuous-time, not discrete-time models, so these are *not* the same as the probability of observing state  $s$  at a fixed time in the future. Those probabilities are given by [pmatrix.msm](#).

## Value

The matrix of probabilities that the next move of a process in state  $r$  (rows) is to state  $s$  (columns).

## Author(s)

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>

## See Also

[qmatrix.msm](#), [pmatrix.msm](#), [qratio.msm](#)

ppass.msm

*Passage probabilities***Description**

Probabilities of having visited each state by a particular time in a continuous time Markov model.

**Usage**

```
ppass.msm(x=NULL, qmatrix=NULL, tot, start="all", covariates="mean",
          piecewise.times=NULL, piecewise.covariates=NULL,
          ci=c("none", "normal", "bootstrap"), cl=0.95, B=1000,
          cores=NULL, ...)
```

**Arguments**

x	A fitted multi-state model, as returned by <a href="#">msm</a> .
qmatrix	Instead of x, you can simply supply a transition intensity matrix in <code>qmatrix</code> .
tot	Finite time to forecast the passage probabilities for.
start	Starting state (integer). By default ( <code>start="all"</code> ), this will return a matrix one row for each starting state.  Alternatively, this can be used to obtain passage probabilities from a <i>set</i> of states, rather than single states. To achieve this, <code>state</code> is set to a vector of weights, with length equal to the number of states in the model. These weights should be proportional to the probability of starting in each of the states in the desired set, so that weights of zero are supplied for other states. The function will calculate the weighted average of the passage probabilities from each of the corresponding states.
covariates	Covariate values defining the intensity matrix for the fitted model x, as supplied to <a href="#">qmatrix.msm</a> .
piecewise.times	Currently ignored: not implemented for time-inhomogeneous models.
piecewise.covariates	Currently ignored: not implemented for time-inhomogeneous models.
ci	If "normal", then calculate a confidence interval by simulating B random vectors from the asymptotic multivariate normal distribution implied by the maximum likelihood estimates (and covariance matrix) of the log transition intensities and covariate effects.  If "bootstrap" then calculate a confidence interval by non-parametric bootstrap refitting. This is 1-2 orders of magnitude slower than the "normal" method, but is expected to be more accurate. See <a href="#">boot.msm</a> for more details of bootstrapping in <b>msm</b> .  If "none" (the default) then no confidence interval is calculated.
cl	Width of the symmetric confidence interval, relative to 1.

B	Number of bootstrap replicates.
cores	Number of cores to use for bootstrapping using parallel processing. See <a href="#">boot.msm</a> for more details.
...	Arguments to pass to <a href="#">MatrixExp</a> .

### Details

The passage probabilities to state  $s$  are computed by setting the  $s$ th row of the transition intensity matrix  $Q$  to zero, giving an intensity matrix  $Q^*$  for a simplified model structure where state  $s$  is absorbing. The probabilities of passage are then equivalent to row  $s$  of the transition probability matrix  $Exp(tQ^*)$  under this simplified model for  $t = tot$ .

Note this is different from the probability of occupying each state at exactly time  $t$ , given by [pmatrix.msm](#). The passage probability allows for the possibility of having visited the state before  $t$ , but then occupying a different state at  $t$ .

The mean of the passage distribution is the expected first passage time, [efpt.msm](#).

This function currently only handles time-homogeneous Markov models. For time-inhomogeneous models the covariates are held constant at the value supplied, by default the column means of the design matrix over all observations.

### Value

A matrix whose  $r, s$  entry is the probability of having visited state  $s$  at least once before time  $t$ , given the state at time 0 is  $r$ . The diagonal entries should all be 1.

### Author(s)

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>

### References

Norris, J. R. (1997) Markov Chains. Cambridge University Press.

### See Also

[efpt.msm](#), [totlos.msm](#), [boot.msm](#).

### Examples

```
Q <- rbind(c(-0.5, 0.25, 0, 0.25), c(0.166, -0.498, 0.166, 0.166),
          c(0, 0.25, -0.5, 0.25), c(0, 0, 0, 0))

## ppass[1,2](t) converges to 0.5 with t, since given in state 1, the
## probability of going to the absorbing state 4 before visiting state
## 2 is 0.5, and the chance of still being in state 1 at t decreases.

ppass.msm(qmatrix=Q, tot=2)
ppass.msm(qmatrix=Q, tot=20)
ppass.msm(qmatrix=Q, tot=100)
```



```

Q <- Q[1:3,1:3]; diag(Q) <- 0; diag(Q) <- -rowSums(Q)

## Probability of about 1/2 of visiting state 3 by time 10.5, the
## median first passage time

ppass.msm(qmatrix=Q, tot=10.5)

## Mean first passage time from state 2 to state 3 is 10.02: similar
## to the median

efpt.msm(qmatrix=Q, tostate=3)

```

prevalence.msm

*Tables of observed and expected prevalences***Description**

This provides a rough indication of the goodness of fit of a multi-state model, by estimating the observed numbers of individuals occupying each state at a series of times, and comparing these with forecasts from the fitted model.

**Usage**

```

prevalence.msm(x, times=NULL, timezero=NULL, initstates=NULL, covariates="population",
  misccovariates="mean", piecewise.times=NULL, piecewise.covariates=NULL,
  ci=c("none", "normal", "bootstrap"), cl=0.95, B=1000, cores=NULL,
  interp=c("start", "midpoint"), censtime=Inf, subset=NULL, plot=FALSE, ...)

```

**Arguments**

x	A fitted multi-state model produced by <a href="#">msm</a> .
times	Series of times at which to compute the observed and expected prevalences of states.
timezero	Initial time of the Markov process. Expected values are forecasted from here. Defaults to the minimum of the observation times given in the data.
initstates	Optional vector of the same length as the number of states. Gives the numbers of individuals occupying each state at the initial time, to be used for forecasting expected prevalences. The default is those observed in the data. These should add up to the actual number of people in the study at the start.
covariates	Covariate values for which to forecast expected state occupancy. With the default <code>covariates="population"</code> , expected prevalences are produced by summing model predictions over the covariates observed in the original data, for a fair comparison with the observed prevalences. This may be slow, particularly with continuous covariates.  Predictions for fixed covariates can be obtained by supplying covariate values in the standard way, as in <a href="#">qmatrix.msm</a> . Therefore if <code>covariates="population"</code>

	is too slow, using the mean observed values through <code>covariates="mean"</code> may give a reasonable approximation.
	This argument is ignored if <code>piecewise.times</code> is specified. If there are a mixture of time-constant and time-dependent covariates, then the values for all covariates should be supplied in <code>piecewise.covariates</code> .
<code>misccovariates</code>	(Misclassification models only) Values of covariates on the misclassification probability matrix for converting expected true to expected misclassified states. Ignored if <code>covariates="population"</code> , otherwise defaults to the mean values of the covariates in the data set.
<code>piecewise.times</code>	Times at which piecewise-constant intensities change. See <a href="#">pmatrix.piecewise.msm</a> for how to specify this. Ignored if <code>covariates="population"</code> . This is only required for time-inhomogeneous models specified using explicit time-dependent covariates, and should not be used for models specified using "pci".
<code>piecewise.covariates</code>	Covariates on which the piecewise-constant intensities depend. See <a href="#">pmatrix.piecewise.msm</a> for how to specify this. Ignored if <code>covariates="population"</code> .
<code>ci</code>	<p>If "normal", then calculate a confidence interval for the expected prevalences by simulating B random vectors from the asymptotic multivariate normal distribution implied by the maximum likelihood estimates (and covariance matrix) of the log transition intensities and covariate effects, then calculating the expected prevalences for each replicate.</p> <p>If "bootstrap" then calculate a confidence interval by non-parametric bootstrap refitting. This is 1-2 orders of magnitude slower than the "normal" method, but is expected to be more accurate. See <a href="#">boot.msm</a> for more details of bootstrapping in <b>msm</b>.</p> <p>If "none" (the default) then no confidence interval is calculated.</p>
<code>c1</code>	Width of the symmetric confidence interval, relative to 1
<code>B</code>	Number of bootstrap replicates
<code>cores</code>	Number of cores to use for bootstrapping using parallel processing. See <a href="#">boot.msm</a> for more details.
<code>interp</code>	<p>Suppose an individual was observed in states <math>S_{r-1}</math> and <math>S_r</math> at two consecutive times <math>t_{r-1}</math> and <math>t_r</math>, and we want to estimate 'observed' prevalences at a time <math>t</math> between <math>t_{r-1}</math> and <math>t_r</math>.</p> <p>If <code>interp="start"</code>, then individuals are assumed to be in state <math>S_{r-1}</math> at time <math>t</math>, the same state as they were at <math>t_{r-1}</math>.</p> <p>If <code>interp="midpoint"</code> then if <math>t \leq (t_{r-1} + t_r)/2</math>, the midpoint of <math>t_{r-1}</math> and <math>t_r</math>, the state at <math>t</math> is assumed to be <math>S_{r-1}</math>, otherwise <math>S_r</math>. This is generally more reasonable for "progressive" models.</p>
<code>censtime</code>	<p>If the time is greater than <code>censtime</code> and the patient has reached an absorbing state, then that subject will be removed from the risk set. For example, if patients have died but would only have been observed up to this time, then this avoids overestimating the proportion of people who are dead at later times.</p> <p>This can be supplied as a single value, or as a vector with one element per subject (after any subset has been taken), in the same order as the original</p>

data. This vector also only includes subjects with complete data, thus it excludes for example subjects with only one observation (thus no observed transitions), and subjects for whom every observation has missing values. (Note, to help construct this, the complete data used for the model fit can be accessed with `model.frame(x)`, where `x` is the fitted model object)

This is ignored if it is less than the subject's maximum observation time.

`subset` Subset of subjects to calculate observed prevalences for.

`plot` Generate a plot of observed against expected prevalences. See [plot.prevalence.msm](#)

`...` Further arguments to pass to [plot.prevalence.msm](#).

## Details

The fitted transition probability matrix is used to forecast expected prevalences from the state occupancy at the initial time. To produce the expected number in state  $j$  at time  $t$  after the start, the number of individuals under observation at time  $t$  (including those who have died, but not those lost to follow-up) is multiplied by the product of the proportion of individuals in each state at the initial time and the transition probability matrix in the time interval  $t$ . The proportion of individuals in each state at the "initial" time is estimated, if necessary, in the same way as the observed prevalences.

For misclassification models (fitted using an `ematrix`), this aims to assess the fit of the full model for the *observed* states. That is, the combined Markov progression model for the true states and the misclassification model. Thus, expected prevalences of *true* states are estimated from the assumed proportion occupying each state at the initial time using the fitted transition probability matrix. The vector of expected prevalences of true states is then multiplied by the fitted misclassification probability matrix to obtain the expected prevalences of *observed* states.

For general hidden Markov models, the observed state is taken to be the predicted underlying state from the Viterbi algorithm ([viterbi.msm](#)). The goodness of fit of these states to the underlying Markov model is tested.

In any model, if there are censored states, then these are replaced by imputed values of highest probability from the Viterbi algorithm in order to calculate the observed state prevalences.

For an example of this approach, see Gentleman *et al.* (1994).

## Value

A list of matrices, with components:

`Observed` Table of observed numbers of individuals in each state at each time

`Observed percentages` Corresponding percentage of the individuals at risk at each time.

`Expected` Table of corresponding expected numbers.

`Expected percentages` Corresponding percentage of the individuals at risk at each time.

Or if `ci.boot = TRUE`, the component `Expected` is a list with components `estimates` and `ci`. `estimates` is a matrix of the expected prevalences, and `ci` is a list of two matrices, containing the confidence limits. The component `Expected percentages` has a similar format.

**Author(s)**

C. H. Jackson <chris.jackson@mrc-bsu.cam.ac.uk>

**References**

Gentleman, R.C., Lawless, J.F., Lindsey, J.C. and Yan, P. Multi-state Markov models for analysing incomplete disease history data with illustrations for HIV disease. *Statistics in Medicine* (1994) 13(3): 805–821.

Titman, A.C., Sharples, L. D. Model diagnostics for multi-state models. *Statistical Methods in Medical Research* (2010) 19(6):621-651.

**See Also**

[msm](#), [summary.msm](#)

---

print.msm

*Print a fitted msm model object*

---

**Description**

Print a fitted msm model object

**Usage**

```
## S3 method for class 'msm'
print(x,covariates=NULL, digits=4, ...)
printnew.msm(x, covariates=NULL, digits=4, ...)
```

**Arguments**

x	Output from <a href="#">msm</a> , representing a fitted multi-state model object.
covariates	Covariates for which to print “baseline” transition intensities or misclassification probabilities. See <a href="#">qmatrix.msm</a> for more details.
digits	Minimum number of significant digits, passed to <a href="#">format</a> . Defaults to 4.
...	Other arguments to be passed to <a href="#">format</a> .

**Details**

This is the new method of formatting msm objects for printing. The old method was based on printing lists of matrices. That produced a lot of wasted space for parameters which were zero, and it was difficult to match corresponding numbers between matrices. The new method presents all the transition intensities and covariate effects as a single compact table, and likewise for misclassification matrices.

Also in the old method, covariate effects were presented as log hazard ratios or log odds ratios. The log scale is more convenient mathematically, but unnatural to interpret. The new method presents hazard ratios for covariates on transition intensities and odds ratios for misclassification probabilities.

printnew.msm is an alias for print.msm.

**Value**

The object returned by `print.msm` is a numeric matrix with one column for each estimate or confidence limit for intensities and their covariates, in the same arrangement as printed, but with the underlying numbers in full precision. The results formatted for printing are stored in the "formatted" attribute of the object, as a character matrix. These can alternatively be produced by `msm.form.qoutput`, which has no printing side-effect. `msm.form.eoutput` produces the same arrangement for misclassification probabilities instead of intensities.

**Author(s)**

C. H. Jackson <chris.jackson@mrc-bsu.cam.ac.uk>

**See Also**

[msm](#), [printold.msm](#), [msm.form.qoutput](#).

---

printold.msm	<i>Print a fitted msm model object</i>
--------------	--

---

**Description**

Print a fitted msm model object (in old format, from `msm` 1.3.1 and earlier)

**Usage**

```
printold.msm(x,...)
```

**Arguments**

x	Output from <a href="#">msm</a> , representing a fitted multi-state model object.
...	Other arguments to be passed to <a href="#">format</a> .

**Details**

See [print.msm](#) for a better and cleaner output format, and an explanation of the change.

**Author(s)**

C. H. Jackson <chris.jackson@mrc-bsu.cam.ac.uk>

**See Also**

[print.msm](#)

---

psor *Psoriatic arthritis data*

---

### Description

A series of observations of grades of psoriatic arthritis, as indicated by numbers of damaged joints.

### Usage

psor

### Format

A data frame containing 806 observations, representing visits to a psoriatic arthritis (PsA) clinic from 305 patients. The rows are grouped by patient number and ordered by examination time. Each row represents an examination and contains additional covariates.

ptnum	(numeric)	Patient identification number
months	(numeric)	Examination time in months
state	(numeric)	Clinical state of PsA. Patients in states 1, 2, 3 and 4 have 0, 1 to 4, 5 to 9 and 10 or more damaged joints, respectively.
hieffusn	(numeric)	Presence of five or more effusions
ollwsdrt	(character)	Erythrocyte sedimentation rate of less than 15 mm/h

### References

Gladman, D. D. and Farewell, V.T. (1999) Progression in psoriatic arthritis: role of time-varying clinical indicators. *J. Rheumatol.* 26(11):2409-13

### Examples

```
## Four-state progression-only model with high effusion and low
## sedimentation rate as covariates on the progression rates. High
## effusion is assumed to have the same effect on the 1-2, 2-3, and 3-4
## progression rates, while low sedimentation rate has the same effect
## on the 1-2 and 2-3 intensities, but a different effect on the 3-4.

data(psor)
psor.q <- rbind(c(0,0.1,0,0),c(0,0,0.1,0),c(0,0,0,0.1),c(0,0,0,0))
psor.msm <- msm(state ~ months, subject=ptnum, data=psor,
               qmatrix = psor.q, covariates = ~ollwsdrt+hieffusn,
               constraint = list(hieffusn=c(1,1,1),ollwsdrt=c(1,1,2)),
               fixedpars=FALSE, control = list(REPORT=1,trace=2), method="BFGS")
qmatrix.msm(psor.msm)
sojourn.msm(psor.msm)
hazard.msm(psor.msm)
```

---

qcmmodel.object	<i>Developer documentation: model for covariates on transition intensities</i>
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---

## Description

A list representing the model for covariates on transition intensities

## Value

npars	<p>Number of covariate effect parameters. This is defined as the number of covariates on intensities (with factors expanded as contrasts) multiplied by the number of allowed transitions in the model.</p> <p>Note if <code>msm</code> was called with <code>covariates</code> set to a list of different covariates for different intensities, then this will include covariate effects that are implicitly defined as zero by this list. The information in <code>paramdata</code> objects can be used to identify which ones are fixed at zero.</p> <p>This also includes any <code>timeperiod</code> covariates in a time-inhomogeneous model defined by the <code>pci</code> option to <code>msm</code>.</p>
ndpars	Number of distinct covariate effect parameters, as <code>npars</code> , but after any equality constraints have been applied.
ncovs	Number of covariates on intensities, with factors expanded as contrasts.
constr	List of equality constraints on these covariate effects, as supplied in the <code>constraint</code> argument to <code>msm</code> .
covlabels	Names / labels of these covariates in the model matrix (see <code>model.matrix.msm</code> ).
inits	Initial values for these covariate effects, as a vector formed from the <code>covinits</code> list supplied to <code>msm</code> .
covmeans	Means of these covariates in the data (excluding data not required to fit the model, such as observations with missing data in other elements or subjects' last observations). This includes means of 0/1 factor contrasts as well as continuous covariates (for historic reasons, which may not be sensible).

## See Also

[msm.object](#).

qgeneric

*Generic function to find quantiles of a distribution***Description**

Generic function to find the quantiles of a distribution, given the equivalent probability distribution function.

**Usage**

```
qgeneric(pdlist, p, special=NULL, ...)
```

**Arguments**

pdlist	Probability distribution function, for example, <a href="#">pnorm</a> for the normal distribution, which must be defined in the current workspace. This should accept and return vectorised parameters and values. It should also return the correct values for the entire real line, for example a positive distribution should have <code>pdlist(x)==0</code> for $x < 0$ .
p	Vector of probabilities to find the quantiles for.
special	Vector of character strings naming arguments of the distribution function that should not be vectorised over. Used, for example, for the <code>rate</code> and <code>t</code> arguments in <a href="#">qpexp</a> .
...	The remaining arguments define parameters of the distribution <code>pdlist</code> . These <b>MUST</b> be named explicitly.  This may also contain the standard arguments <code>log.p</code> (logical; default FALSE, if TRUE, probabilities <code>p</code> are given as $\log(p)$ ), and <code>lower.tail</code> (logical; if TRUE (default), probabilities are $P[X \leq x]$ otherwise, $P[X > x]$ ).  If the distribution is bounded above or below, then this should contain arguments <code>lbound</code> and <code>ubound</code> respectively, and these will be returned if <code>p</code> is 0 or 1 respectively. Defaults to <code>-Inf</code> and <code>Inf</code> respectively.

**Details**

This function is intended to enable users to define "q" functions for new distributions, in cases where the distribution function `pdlist` is available analytically, but the quantile function is not.

It works by finding the root of the equation  $h(q) = pdlist(q) - p = 0$ . Starting from the interval  $(-1, 1)$ , the interval width is expanded by 50% until  $h()$  is of opposite sign at either end. The root is then found using [uniroot](#).

This assumes a suitably smooth, continuous distribution.

An identical function is provided in the **flexsurv** package.

**Value**

Vector of quantiles of the distribution at `p`.



**Author(s)**

Christopher Jackson <chris.jackson@mrc-bsu.cam.ac.uk>

**Examples**

```
qnorm(c(0.025, 0.975), 0, 1)
qgeneric(pnorm, c(0.025, 0.975), mean=0, sd=1) # must name the arguments
```

---

qmatrix.msm	<i>Transition intensity matrix</i>
-------------	------------------------------------

---

**Description**

Extract the estimated transition intensity matrix, and the corresponding standard errors, from a fitted multi-state model at a given set of covariate values.

**Usage**

```
qmatrix.msm(x, covariates="mean", sojourn=FALSE,
            ci=c("delta", "normal", "bootstrap", "none"), cl=0.95,
            B=1000, cores=NULL)
```

**Arguments**

x	A fitted multi-state model, as returned by <a href="#">msm</a> .
covariates	<p>The covariate values at which to estimate the intensity matrix. This can either be:</p> <p>the string "mean", denoting the means of the covariates in the data (this is the default),</p> <p>the number 0, indicating that all the covariates should be set to zero,</p> <p>or a list of values, with optional names. For example</p> <pre>list(60, 1)</pre> <p>where the order of the list follows the order of the covariates originally given in the model formula. Or more clearly, a named list,</p> <pre>list(age = 60, sex = 1)</pre> <p>If some covariates are specified but not others, the missing ones default to zero. With <code>covariates="mean"</code>, for factor / categorical variables, the mean of the 0/1 dummy variable for each factor level is used, representing an average over all values in the data, rather than a specific factor level.</p>
sojourn	Set to TRUE if the estimated sojourn times and their standard errors should also be returned.

ci	<p>If "delta" (the default) then confidence intervals are calculated by the delta method, or by simple transformation of the Hessian in the very simplest cases. Normality on the log scale is assumed.</p> <p>If "normal", then calculate a confidence interval by simulating B random vectors from the asymptotic multivariate normal distribution implied by the maximum likelihood estimates (and covariance matrix) of the log transition intensities and covariate effects, then transforming.</p> <p>If "bootstrap" then calculate a confidence interval by non-parametric bootstrap refitting. This is 1-2 orders of magnitude slower than the "normal" method, but is expected to be more accurate. See <a href="#">boot.msm</a> for more details of bootstrapping in <b>msm</b>.</p>
c1	Width of the symmetric confidence interval to present. Defaults to 0.95.
B	Number of bootstrap replicates, or number of normal simulations from the distribution of the MLEs.
cores	Number of cores to use for bootstrapping using parallel processing. See <a href="#">boot.msm</a> for more details.

### Details

Transition intensities and covariate effects are estimated on the log scale by [msm](#). A covariance matrix is estimated from the Hessian of the maximised log-likelihood.

A more practically meaningful parameterisation of a continuous-time Markov model with transition intensities  $q_{rs}$  is in terms of the mean sojourn times  $-1/q_{rr}$  in each state  $r$  and the probabilities that the next move of the process when in state  $r$  is to state  $s$ ,  $-q_{rs}/q_{rr}$ .

### Value

A list with components:

estimate	Estimated transition intensity matrix.
SE	Corresponding approximate standard errors.
L	Lower confidence limits
U	Upper confidence limits

Or if ci="none", then `qmatrix.msm` just returns the estimated transition intensity matrix.

If `sojourn` is TRUE, extra components called `sojourn`, `sojournSE`, `sojournL` and `sojournU` are included, containing the estimates, standard errors and confidence limits, respectively, of the mean sojourn times in each transient state.

The default print method for objects returned by `qmatrix.msm` presents estimates and confidence limits. To present estimates and standard errors, do something like

```
qmatrix.msm(x)[c("estimates", "SE")]
```

### Author(s)

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>

**See Also**

[pmatrx.msm](#), [sojourn.msm](#), [deltamethod](#), [ematrix.msm](#)

---

qmodel.object

*Developer documentation: transition model structure object*


---

**Description**

A list giving information about the structure of states and allowed transitions in a multi-state model, and options for likelihood calculation. Used in internal computations, and returned in a fitted [msm](#) model object.

**Value**

nstates	Number of states
iso	Label for which basic structure the model is isomorphic to in the list of structures for which analytic formulae for the transition probabilities are implemented in the source file <code>src/analyticp.c</code> . This list is given by the internal object <code>msm:::msm.graphs</code> which is defined and documented in the source file <code>R/constants.R</code> . iso is 0 if the analytic P matrix is not implemented for this structure, or if analytic P matrix calculations are disabled using <code>use.analyticp=FALSE</code> in the call to <a href="#">msm</a> .
perm	Permutation required to convert the base isomorphism into the structure of this model. A vector of integers whose $r$ th element is the state number in the base structure representing state $r$ in the current structure.
qperm	Inverse permutation: vector whose $r$ th element is the state number in the current structure representing the $r$ th state in the base structure.
npars	Number of allowed instantaneous transitions, equal to <code>sum(imatrix)</code> .
imatrix	Indicator matrix for allowed instantaneous transitions. This has $(r, s)$ entry 1 if the transition from $r$ to $s$ is permitted in continuous time, and 0 otherwise. The diagonal entries are arbitrarily set to 0.
qmatrix	Matrix of initial values for the transition intensities, supplied as the <code>qmatrix</code> argument of <a href="#">msm</a> .
inits	Vector of these initial values, reading across rows of <code>qmatrix</code> and excluding the diagonal and disallowed transitions.
constr	Indicators for equality constraints on baseline intensities, taken from the <code>qconstraint</code> argument to <a href="#">msm</a> , and mapped if necessary to the set $(1,2,3,\dots)$ .
ndpars	Number of distinct allowed instantaneous transitions, after applying equality constraints.
expm	Use <b>expm</b> package to calculate matrix exponentials for likelihoods, as supplied to the <code>use.expm</code> argument of <a href="#">msm</a> . TRUE or FALSE.

**See Also**

[msm.object](#), [emodel.object](#), [hmodel.object](#).

---

 qratio.msm

*Estimated ratio of transition intensities*


---

### Description

Compute the estimate and approximate standard error of the ratio of two estimated transition intensities from a fitted multi-state model at a given set of covariate values.

### Usage

```
qratio.msm(x, ind1, ind2, covariates = "mean",
           ci=c("delta","normal","bootstrap","none"), c1 = 0.95,
           B=1000, cores=NULL)
```

### Arguments

x	A fitted multi-state model, as returned by <a href="#">msm</a> .
ind1	Pair of numbers giving the indices in the intensity matrix of the numerator of the ratio, for example, <code>c(1, 2)</code> .
ind2	Pair of numbers giving the indices in the intensity matrix of the denominator of the ratio, for example, <code>c(2, 1)</code> .
covariates	The covariate values at which to estimate the intensities. This can either be: <ul style="list-style-type: none"> <li>the string "mean", denoting the means of the covariates in the data (this is the default),</li> <li>the number 0, indicating that all the covariates should be set to zero,</li> <li>or a list of values, with optional names. For example <code>list(60, 1)</code> where the order of the list follows the order of the covariates originally given in the model formula, or a named list, <code>list(age = 60, sex = 1)</code></li> </ul>
ci	If "delta" (the default) then confidence intervals are calculated by the delta method. If "normal", then calculate a confidence interval by simulating B random vectors from the asymptotic multivariate normal distribution implied by the maximum likelihood estimates (and covariance matrix) of the log transition intensities and covariate effects, then transforming. If "bootstrap" then calculate a confidence interval by non-parametric bootstrap refitting. This is 1-2 orders of magnitude slower than the "normal" method, but is expected to be more accurate. See <a href="#">boot.msm</a> for more details of bootstrapping in <b>msm</b> .
c1	Width of the symmetric confidence interval to present. Defaults to 0.95.

B	Number of bootstrap replicates, or number of normal simulations from the distribution of the MLEs
cores	Number of cores to use for bootstrapping using parallel processing. See <a href="#">boot.msm</a> for more details.

### Details

For example, we might want to compute the ratio of the progression rate and recovery rate for a fitted model `disease.msm` with a health state (state 1) and a disease state (state 2). In this case, the progression rate is the (1,2) entry of the intensity matrix, and the recovery rate is the (2,1) entry. Thus to compute this ratio with covariates set to their means, we call

```
qratio.msm(disease.msm, c(1, 2), c(2, 1)) .
```

Standard errors are estimated by the delta method. Confidence limits are estimated by assuming normality on the log scale.

### Value

A named vector with elements `estimate`, `se`, `L` and `U` containing the estimate, standard error, lower and upper confidence limits, respectively, of the ratio of intensities.

### Author(s)

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>

### See Also

[qmatrix.msm](#)

---

recreate.olddata	<i>Convert data stored in msm object to old format</i>
------------------	--

---

### Description

Converts the data element of `msm` objects to the old format.

### Usage

```
recreate.olddata(x)
```

### Arguments

x                    Object returned by the `msm` function, representing a fitted multi-state model.

### Details

This is just provided for convenience and to illustrate the changes. It is not guaranteed to be complete, and is liable to be withdrawn. Users who were relying on the previous undocumented format are advised to upgrade their code to use the new format, which uses model frames and model design matrices in the standard format used in version 1.4, based on [model.frame](#) and [model.matrix](#).

**Value**

A list of vectors and matrices in the undocumented ad-hoc format used for the data component of msm objects in **msm** versions 1.3.1 and earlier.

---

scorer resid.msm	<i>Score residuals</i>
------------------	------------------------

---

**Description**

Score residuals for detecting outlying subjects.

**Usage**

```
scorer resid.msm(x, plot=FALSE)
```

**Arguments**

x	A fitted multi-state model, as returned by <a href="#">msm</a> .
plot	If TRUE, display a simple plot of the residuals in subject order, labelled by subject identifiers

**Details**

The score residual for a single subject is

$$U(\theta)^T I(\theta)^{-1} U(\theta)$$

where  $U(\theta)$  is the vector of first derivatives of the log-likelihood for that subject at maximum likelihood estimates  $\theta$ , and  $I(\theta)$  is the observed Fisher information matrix, that is, the matrix of second derivatives of minus the log-likelihood for that subject at theta.

Subjects with a higher influence on the maximum likelihood estimates will have higher score residuals.

These are only available for models with analytic derivatives (which includes all non-hidden and most hidden Markov models).

**Value**

Vector of the residuals, named by subject identifiers.

**Author(s)**

Andrew Titman <a.titman@lancaster.ac.uk> (theory), Chris Jackson <chris.jackson@mrc-bsu.cam.ac.uk> (code)

---

sim.msm	<i>Simulate one individual trajectory from a continuous-time Markov model</i>
---------	---

---

### Description

Simulate one realisation from a continuous-time Markov process up to a given time.

### Usage

```
sim.msm(qmatrix, maxtime, covs=NULL, beta=NULL, obstimes=0, start=1,
mintime=0)
```

### Arguments

qmatrix	The transition intensity matrix of the Markov process. The diagonal of qmatrix is ignored, and computed as appropriate so that the rows sum to zero. For example, a possible qmatrix for a three state illness-death model with recovery is: <code>rbind( c( 0,0.1,0.02 ), c( 0.1,0,0.01 ), c( 0,0,0 ) )</code>
maxtime	Maximum time for the simulated process.
covs	Matrix of time-dependent covariates, with one row for each observation time and one column for each covariate.
beta	Matrix of linear covariate effects on log transition intensities. The rows correspond to different covariates, and the columns to the transition intensities. The intensities are ordered by reading across rows of the intensity matrix, starting with the first, counting the positive off-diagonal elements of the matrix.
obstimes	Vector of times at which the covariates are observed.
start	Starting state of the process. Defaults to 1.
mintime	Starting time of the process. Defaults to 0.

### Details

The effect of time-dependent covariates on the transition intensity matrix for an individual is determined by assuming that the covariate is a step function which remains constant in between the individual's observation times.

### Value

A list with components,

states	Simulated states through which the process moves. This ends with either an absorption before <code>obstime</code> , or a transient state at <code>obstime</code> .
times	Exact times at which the process changes to the corresponding states
qmatrix	The given transition intensity matrix

**Author(s)**

C. H. Jackson <chris.jackson@mrc-bsu.cam.ac.uk>

**See Also**

[simmulti.msm](#)

**Examples**

```
qmatrix <- rbind(
  c(-0.2, 0.1, 0.1 ),
  c(0.5, -0.6, 0.1 ),
  c(0, 0, 0)
)
sim.msm(qmatrix, 30)
```

---

simfitted.msm

*Simulate from a Markov model fitted using msm*

---

**Description**

Simulate a dataset from a Markov model fitted using [msm](#), using the maximum likelihood estimates as parameters, and the same observation times as in the original data.

**Usage**

```
simfitted.msm(x, drop.absorb=TRUE, drop.pci.imp=TRUE)
```

**Arguments**

x	A fitted multi-state model object as returned by <a href="#">msm</a> .
drop.absorb	Should repeated observations in an absorbing state be omitted. Use the default of TRUE to avoid warnings when using the simulated dataset for further <a href="#">msm</a> fits. Or set to FALSE if exactly the same number of observations as the original data are needed.
drop.pci.imp	In time-inhomogeneous models fitted using the pci option to <a href="#">msm</a> , censored observations are inserted into the data by <a href="#">msm</a> at the times where the intensity changes, but dropped by default when simulating from the fitted model using this function. Set this argument to FALSE to keep these observations and the corresponding indicator variable.



**Details**

This function is a wrapper around [simmulti.msm](#), and only simulates panel-observed data. To generate datasets with the exact times of transition, use the lower-level [sim.msm](#).

Markov models with misclassified states fitted through the `ematrix` option to [msm](#) are supported, but not general hidden Markov models with `hmodel`. For misclassification models, this function includes misclassification in the simulated states.

This function is used for parametric bootstrapping to estimate the null distribution of the test statistic in [pearson.msm](#).

**Value**

A dataset with variables as described in [simmulti.msm](#).

**Author(s)**

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>

**See Also**

[simmulti.msm](#), [sim.msm](#), [pearson.msm](#), [msm](#).

---

simmulti.msm	<i>Simulate multiple trajectories from a multi-state Markov model with arbitrary observation times</i>
--------------	--

---

**Description**

Simulate a number of individual realisations from a continuous-time Markov process. Observations of the process are made at specified arbitrary times for each individual, giving panel-observed data.

**Usage**

```
simmulti.msm(data, qmatrix, covariates=NULL, death = FALSE, start,
             ematrix=NULL, miscovariates=NULL, hmodel=NULL, hcovariates=NULL,
             censor.states=NULL, drop.absorb=TRUE)
```

**Arguments**

data	A data frame with a mandatory column named <code>time</code> , representing observation times. The optional column named <code>subject</code> , corresponds to subject identification numbers. If not given, all observations are assumed to be on the same individual. Observation times should be sorted within individuals. The optional column named <code>cens</code> indicates the times at which simulated states should be censored. If <code>cens==0</code> then the state is not censored, and if <code>cens==k</code> , say, then all simulated states at that time which are in the set <code>censor.states</code> are replaced by <code>k</code> . Other named columns of the data frame represent any covariates, which may be time-constant or time-dependent. Time-dependent covariates are assumed to be constant between the observation times.
------	---

qmatrix	The transition intensity matrix of the Markov process, with any covariates set to zero. The diagonal of qmatrix is ignored, and computed as appropriate so that the rows sum to zero. For example, a possible qmatrix for a three state illness-death model with recovery is: <code>rbind( c( 0,0.1,0.02 ), c( 0.1,0,0.01 ), c( 0,0,0 ) )</code>
covariates	List of linear covariate effects on log transition intensities. Each element is a vector of the effects of one covariate on all the transition intensities. The intensities are ordered by reading across rows of the intensity matrix, starting with the first, counting the positive off-diagonal elements of the matrix. For example, for a multi-state model with three transition intensities, and two covariates x and y on each intensity, <code>covariates=list(x = c(-0.3, -0.3, -0.3), y=c(0.1, 0.1, 0.1))</code>
death	Vector of indices of the death states. A death state is an absorbing state whose time of entry is known exactly, but the individual is assumed to be in an unknown transient state ("alive") at the previous instant. This is the usual situation for times of death in chronic disease monitoring data. For example, if you specify <code>death = c(4, 5)</code> then states 4 and 5 are assumed to be death states. <code>death = TRUE</code> indicates that the final state is a death state, and <code>death = FALSE</code> (the default) indicates that there is no death state.
start	A vector with the same number of elements as there are distinct subjects in the data, giving the states in which each corresponding individual begins. Or a single number, if all of these are the same. Defaults to state 1 for each subject.
ematrix	An optional misclassification matrix for generating observed states conditionally on the simulated true states. As defined in <a href="#">msm</a> .
miscovariates	Covariate effects on misclassification probabilities via multinomial logistic regression. Linear effects operate on the log of each probability relative to the probability of classification in the correct state. In same format as <code>covariates</code> .
hmodel	An optional hidden Markov model for generating observed outcomes conditionally on the simulated true states. As defined in <a href="#">msm</a> .
hcovariates	List of the same length as <code>hmodel</code> , defining any covariates governing the hidden Markov outcome models. Unlike in the <code>msm</code> function, this should also define the values of the covariate effects. Each element of the list is a named vector of the initial values for each set of covariates for that state. For example, for a three-state hidden Markov model with two, one and no covariates on the state 1, 2 and 3 outcome models respectively, <code>hcovariates = list( c(acute=-8, age=0), c(acute=-8), NULL)</code>
sensor.states	Set of simulated states which should be replaced by a censoring indicator at censoring times. By default this is all transient states (representing alive, with unknown state).
drop.absorb	Drop repeated observations in the absorbing state, retaining only one.

## Details

`sim.msm` is called repeatedly to produce a simulated trajectory for each individual. The state at each specified observation time is then taken to produce a new column state. The effect of time-dependent covariates on the transition intensity matrix for an individual is determined by assuming

that the covariate is a step function which remains constant in between the individual's observation times. If the subject enters an absorbing state, then only the first observation in that state is kept in the data frame. Rows corresponding to future observations are deleted. The entry times into states given in death are assumed to be known exactly.

### Value

A data frame with columns,

subject	Subject identification indicators
time	Observation times
state	Simulated (true) state at the corresponding time
obs	Observed outcome at the corresponding time, if <code>ematrix</code> or <code>hmodel</code> was supplied
keep	Row numbers of the original data. Useful when <code>drop.absorb=TRUE</code> , to show which rows were not dropped

plus any supplied covariates.

### Author(s)

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>

### See Also

[sim.msm](#)

### Examples

```
### Simulate 100 individuals with common observation times
sim.df <- data.frame(subject = rep(1:100, rep(13,100)), time = rep(seq(0, 24, 2), 100))
qmatrix <- rbind(c(-0.11, 0.1, 0.01),
                c(0.05, -0.15, 0.1),
                c(0.02, 0.07, -0.09))
simmulti.msm(sim.df, qmatrix)
```

---

sojourn.msm

*Mean sojourn times from a multi-state model*

---

### Description

Estimate the mean sojourn times in the transient states of a multi-state model and their confidence limits.

### Usage

```
sojourn.msm(x, covariates="mean", ci=c("delta","normal","bootstrap","none"),
            cl=0.95, B=1000)
```

**Arguments**

x	A fitted multi-state model, as returned by <a href="#">msm</a> .
covariates	The covariate values at which to estimate the mean sojourn times. This can either be: <ul style="list-style-type: none"> <li>the string "mean", denoting the means of the covariates in the data (this is the default),</li> <li>the number 0, indicating that all the covariates should be set to zero,</li> <li>a list of values, with optional names. For example, <code>list(60, 1)</code>, where the order of the list follows the order of the covariates originally given in the model formula, or a named list, e.g. <code>list(age = 60, sex = 1)</code></li> </ul>
ci	If "delta" (the default) then confidence intervals are calculated by the delta method, or by simple transformation of the Hessian in the very simplest cases. If "normal", then calculate a confidence interval by simulating B random vectors from the asymptotic multivariate normal distribution implied by the maximum likelihood estimates (and covariance matrix) of the log transition intensities and covariate effects, then transforming. If "bootstrap" then calculate a confidence interval by non-parametric bootstrap refitting. This is 1-2 orders of magnitude slower than the "normal" method, but is expected to be more accurate. See <a href="#">boot.msm</a> for more details of bootstrapping in <b>msm</b> .
c1	Width of the symmetric confidence interval to present. Defaults to 0.95.
B	Number of bootstrap replicates, or number of normal simulations from the distribution of the MLEs

**Details**

The mean sojourn time in a transient state  $r$  is estimated by  $-1/q_{rr}$ , where  $q_{rr}$  is the  $r$ th entry on the diagonal of the estimated transition intensity matrix.

A continuous-time Markov model is fully specified by the mean sojourn times and the probability that each state is next ([pnext.msm](#)). This is a more intuitively meaningful description of a model than the transition intensity matrix ([qmatrix.msm](#)).

Time dependent covariates, or time-inhomogeneous models, are not supported. This would require the mean of a piecewise exponential distribution, and the package author is not aware of any general analytic form for that.

**Value**

A data frame with components:

estimates	Estimated mean sojourn times in the transient states.
SE	Corresponding standard errors.
L	Lower confidence limits.
U	Upper confidence limits.

**Author(s)**

C. H. Jackson <chris.jackson@mrc-bsu.cam.ac.uk>

**See Also**

[msm](#), [qmatrix.msm](#), [deltamethod](#)

---

statetable.msm	<i>Table of transitions</i>
----------------	-----------------------------

---

**Description**

Calculates a frequency table counting the number of times each pair of states were observed in successive observation times. This can be a useful way of summarising multi-state data.

**Usage**

```
statetable.msm(state, subject, data=NULL)
```

**Arguments**

state	Observed states, assumed to be ordered by time within each subject.
subject	Subject identification numbers corresponding to state. If not given, all observations are assumed to be on the same subject.
data	An optional data frame in which the variables represented by subject and state can be found.

**Details**

If the data are intermittently observed (panel data) this table should not be used to decide what transitions should be allowed in the  $Q$  matrix, which works in continuous time. This function counts the transitions between states over a time interval, not in real time. There can be observed transitions between state  $r$  and  $s$  over an interval even if  $q_{r,s} = 0$ , because the process may have passed through one or more intermediate states in the middle of the interval.

**Value**

A frequency table with starting states as rows and finishing states as columns.

**Author(s)**

C. H. Jackson <chris.jackson@mrc-bsu.cam.ac.uk>

**See Also**

[crudeinits.msm](#)

**Examples**

```
## Heart transplant data
data(cav)

## 148 deaths from state 1, 48 from state 2 and 55 from state 3.
statetable.msm(state, PTNUM, data=cav)
```

---

surface.msm                      *Explore the likelihood surface*

---

**Description**

Plot the log-likelihood surface with respect to two parameters.

**Usage**

```
surface.msm(x, params=c(1,2), np=10, type=c("contour","filled.contour","persp","image"),
           point=NULL, xrange=NULL, yrange=NULL,...)
## S3 method for class 'msm'
contour(x, ...)
## S3 method for class 'msm'
persp(x, ...)
## S3 method for class 'msm'
image(x, ...)
```

**Arguments**

x	Output from <a href="#">msm</a> , representing a fitted msm model.
params	Integer vector with two elements, giving the indices of the parameters to vary. All other parameters will be fixed. Defaults to <code>c(1, 2)</code> , representing the first two log transition intensities. See the <code>fixedpars</code> argument to <code>msm</code> for a definition of these indices.
np	Number of grid points to use in each direction, by default 10. An <code>np x np</code> grid will be used to evaluate the likelihood surface. If 100 likelihood function evaluations is slow, then reduce this.
type	Character string specifying the type of plot to produce. <ul style="list-style-type: none"> <li>"contour"                      Contour plot, using the R function <a href="#">contour</a>.</li> <li>"filled.contour"              Solid-color contour plot, using the R function <a href="#">filled.contour</a>.</li> <li>"persp"                        Perspective plot, using the R function <a href="#">persp</a>.</li> <li>"image"                        Grid color plot, using the R function <a href="#">image</a>.</li> </ul>
point	Vector of length <code>n</code> , where <code>n</code> is the number of parameters in the model, including the parameters that will be varied here. This specifies the point at which to fix

	the likelihood. By default, this is the maximum likelihood estimates stored in the fitted model <code>x</code> , <code>x\$estimates</code> .
<code>xrange</code>	Range to plot for the first varied parameter. Defaults to plus and minus two standard errors, obtained from the Hessian at the maximum likelihood estimate.
<code>yrange</code>	Range to plot for the second varied parameter. Defaults to plus and minus two standard errors, obtained from the Hessian at the maximum likelihood estimate.
<code>...</code>	Further arguments to be passed to the plotting function.

### Details

Draws a contour or perspective plot. Useful for diagnosing irregularities in the likelihood surface. If you want to use these plots before running the maximum likelihood estimation, then just run `msm` with all estimates fixed at their initial values.

`contour.msm` just calls `surface.msm` with `type = "contour"`.

`persp.msm` just calls `surface.msm` with `type = "persp"`.

`image.msm` just calls `surface.msm` with `type = "image"`.

As these three functions are methods of the generic functions `contour`, `persp` and `image`, they can be invoked as `contour(x)`, `persp(x)` or `image(x)`, where `x` is a fitted `msm` object.

### Author(s)

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>

### See Also

[msm](#), [contour](#), [filled.contour](#), [persp](#), [image](#).

---

tnorm	<i>Truncated Normal distribution</i>
-------	--------------------------------------

---

### Description

Density, distribution function, quantile function and random generation for the truncated Normal distribution with mean equal to `mean` and standard deviation equal to `sd` before truncation, and truncated on the interval `[lower, upper]`.

### Usage

```

dtnorm(x, mean=0, sd=1, lower=-Inf, upper=Inf, log = FALSE)
ptnorm(q, mean=0, sd=1, lower=-Inf, upper=Inf,
       lower.tail = TRUE, log.p = FALSE)
qtnorm(p, mean=0, sd=1, lower=-Inf, upper=Inf,
       lower.tail = TRUE, log.p = FALSE)
rtnorm(n, mean=0, sd=1, lower=-Inf, upper=Inf)

```

**Arguments**

<code>x, q</code>	vector of quantiles.
<code>p</code>	vector of probabilities.
<code>n</code>	number of observations. If <code>length(n) &gt; 1</code> , the length is taken to be the number required.
<code>mean</code>	vector of means.
<code>sd</code>	vector of standard deviations.
<code>lower</code>	lower truncation point.
<code>upper</code>	upper truncation point.
<code>log</code>	logical; if TRUE, return log density or log hazard.
<code>log.p</code>	logical; if TRUE, probabilities <code>p</code> are given as <code>log(p)</code> .
<code>lower.tail</code>	logical; if TRUE (default), probabilities are $P[X \leq x]$ , otherwise, $P[X > x]$ .

**Details**

The truncated normal distribution has density

$$f(x, \mu, \sigma) = \phi(x, \mu, \sigma) / (\Phi(u, \mu, \sigma) - \Phi(l, \mu, \sigma))$$

for  $l \leq x \leq u$ , and 0 otherwise.

$\mu$  is the mean of the original Normal distribution before truncation,

$\sigma$  is the corresponding standard deviation,

$u$  is the upper truncation point,

$l$  is the lower truncation point,

$\phi(x)$  is the density of the corresponding normal distribution, and

$\Phi(x)$  is the distribution function of the corresponding normal distribution.

If `mean` or `sd` are not specified they assume the default values of 0 and 1, respectively.

If `lower` or `upper` are not specified they assume the default values of `-Inf` and `Inf`, respectively, corresponding to no lower or no upper truncation.

Therefore, for example, `dtnorm(x)`, with no other arguments, is simply equivalent to `dnorm(x)`.

Only `rtnorm` is used in the `msm` package, to simulate from hidden Markov models with truncated normal distributions. This uses the rejection sampling algorithms described by Robert (1995).

These functions are merely provided for completion, and are not optimized for numerical stability or speed. To fit a hidden Markov model with a truncated Normal response distribution, use a [hmmTNorm](#) constructor. See the [hmm-dists](#) help page for further details.

**Value**

`dtnorm` gives the density, `ptnorm` gives the distribution function, `qtnorm` gives the quantile function, and `rtnorm` generates random deviates.

**Author(s)**

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>



## References

Robert, C. P. Simulation of truncated normal variables. *Statistics and Computing* (1995) 5, 121–125

## See Also

[dnorm](#)

## Examples

```
x <- seq(50, 90, by=1)
plot(x, dnorm(x, 70, 10), type="l", ylim=c(0,0.06)) ## standard Normal distribution
lines(x, dtnorm(x, 70, 10, 60, 80), type="l")      ## truncated Normal distribution
```

---

totlos.msm	<i>Total length of stay, or expected number of visits</i>
------------	---

---

## Description

Estimate the expected total length of stay, or the expected number of visits, in each state, for an individual in a given period of evolution of a multi-state model.

## Usage

```
totlos.msm(x, start=1, end=NULL, fromt=0, tot=Inf, covariates="mean",
           piecewise.times=NULL, piecewise.covariates=NULL,
           num.integ=FALSE, discount=0, env=FALSE,
           ci=c("none", "normal", "bootstrap"), cl=0.95, B=1000,
           cores=NULL, ...)
envisits.msm(x, start=1, end=NULL, fromt=0, tot=Inf, covariates="mean",
             piecewise.times=NULL, piecewise.covariates=NULL,
             num.integ=FALSE, discount=0,
             ci=c("none", "normal", "bootstrap"), cl=0.95, B=1000,
             cores=NULL, ...)
```

## Arguments

x	A fitted multi-state model, as returned by <a href="#">msm</a> .
start	Either a single number giving the state at the beginning of the period, or a vector of probabilities of being in each state at this time.
end	States to estimate the total length of stay (or number of visits) in. Defaults to all states. This is deprecated, since with the analytic solution (see "Details") it doesn't save any computation to only estimate for a subset of states.
fromt	Time from which to estimate. Defaults to 0, the beginning of the process.

tot	Time up to which the estimate is made. Defaults to infinity, giving the expected time spent in or number of visits to the state until absorption. However, the calculation will be much more efficient if a finite (potentially large) time is specified: see the "Details" section. For models without an absorbing state, t must be specified.
covariates	The covariate values to estimate for. This can either be:  the string "mean", denoting the means of the covariates in the data (this is the default),  the number 0, indicating that all the covariates should be set to zero,  or a list of values, with optional names. For example list (60, 1) where the order of the list follows the order of the covariates originally given in the model formula, or a named list, list (age = 60, sex = 1)
piecewise.times	Times at which piecewise-constant intensities change. See <a href="#">pmatrix.piecewise.msm</a> for how to specify this. This is only required for time-inhomogeneous models specified using explicit time-dependent covariates, and should not be used for models specified using "pci".
piecewise.covariates	Covariates on which the piecewise-constant intensities depend. See <a href="#">pmatrix.piecewise.msm</a> for how to specify this.
num.integ	Use numerical integration instead of analytic solution (see below).
discount	Discount rate in continuous time.
env	Supplied to <a href="#">totlos.msm</a> . If TRUE, return the expected number of visits to each state. If FALSE, return the total length of stay in each state. <a href="#">envisits.msm</a> simply calls <a href="#">totlos.msm</a> with env=TRUE.
ci	If "normal", then calculate a confidence interval by simulating B random vectors from the asymptotic multivariate normal distribution implied by the maximum likelihood estimates (and covariance matrix) of the log transition intensities and covariate effects, then calculating the total length of stay for each replicate. If "bootstrap" then calculate a confidence interval by non-parametric bootstrap refitting. This is 1-2 orders of magnitude slower than the "normal" method, but is expected to be more accurate. See <a href="#">boot.msm</a> for more details of bootstrapping in <b>msm</b> . If "none" (the default) then no confidence interval is calculated.
c1	Width of the symmetric confidence interval, relative to 1
B	Number of bootstrap replicates
cores	Number of cores to use for bootstrapping using parallel processing. See <a href="#">boot.msm</a> for more details.
...	Further arguments to be passed to the <a href="#">integrate</a> function to control the numerical integration.

## Details

The expected total length of stay in state  $j$  between times  $t_1$  and  $t_2$ , from the point of view of an individual in state  $i$  at time 0, is defined by the integral from  $t_1$  to  $t_2$  of the  $i, j$  entry of the transition probability matrix  $P(t) = \text{Exp}(tQ)$ , where  $Q$  is the transition intensity matrix.

The corresponding expected number of visits to state  $j$  (excluding the stay in the current state at time 0) is  $\sum_{i \neq j} T_i Q_{i,j}$ , where  $T_i$  is the expected amount of time spent in state  $i$ .

More generally, suppose that  $\pi_0$  is the vector of probabilities of being in each state at time 0, supplied in `start`, and we want the vector  $\mathbf{x}$  giving the expected lengths of stay in each state. The corresponding integral has the following solution (van Loan 1978; van Rosmalen et al. 2013)

$$\mathbf{x} = \begin{bmatrix} 1 & \mathbf{0}_K \end{bmatrix} \text{Exp}(tQ') \begin{bmatrix} \mathbf{0}_K \\ I_K \end{bmatrix}$$

where

$$Q' = \begin{bmatrix} 0 & \pi_0 \\ \mathbf{0}_K & Q - rI_K \end{bmatrix}$$

$\pi_0$  is the row vector of initial state probabilities supplied in `start`,  $\mathbf{0}_K$  is the row vector of  $K$  zeros,  $r$  is the discount rate,  $I_K$  is the  $K \times K$  identity matrix, and  $\text{Exp}$  is the matrix exponential.

Alternatively, the integrals can be calculated numerically, using the `integrate` function. This may take a long time for models with many states where  $P(t)$  is expensive to calculate. This is required where `tot = Inf`, since the package author is not aware of any analytic expression for the limit of the above formula as  $t$  goes to infinity.

With the argument `num.integ=TRUE`, numerical integration is used even where the analytic solution is available. This facility is just provided for checking results against versions 1.2.4 and earlier, and will be removed eventually. Please let the package maintainer know if any results are different.

For a model where the individual has only one place to go from each state, and each state is visited only once, for example a progressive disease model with no recovery or death, these are equal to the mean sojourn time in each state. However, consider a three-state health-disease-death model with transitions from health to disease, health to death, and disease to death, where everybody starts healthy. In this case the mean sojourn time in the disease state will be greater than the expected length of stay in the disease state. This is because the mean sojourn time in a state is conditional on entering the state, whereas the expected total time diseased is a forecast for a healthy individual, who may die before getting the disease.

In the above formulae,  $Q$  is assumed to be constant over time, but the results generalise easily to piecewise-constant intensities. This function automatically handles models fitted using the `pci` option to `msm`. For any other inhomogeneous models, the user must specify `piecewise.times` and `piecewise.covariates` arguments to `totlos.msm`.

## Value

A vector of expected total lengths of stay (`totlos.msm`), or expected number of visits (`envisits.msm`), for each transient state.

## Author(s)

C. H. Jackson <chris.jackson@mrc-bsu.cam.ac.uk>

## References

C. van Loan (1978). Computing integrals involving the matrix exponential. IEEE Transactions on Automatic Control 23(3)395-404.

J. van Rosmalen, M. Toy and J.F. O'Mahony (2013). A mathematical approach for evaluating Markov models in continuous time without discrete-event simulation. Medical Decision Making 33:767-779.

## See Also

[sojourn.msm](#), [pmatrix.msm](#), [integrate](#), [boot.msm](#).

---

transient.msm

*Transient and absorbing states*

---

## Description

Returns the transient and absorbing states of either a fitted model or a transition intensity matrix.

## Usage

```
transient.msm(x=NULL, qmatrix=NULL)
absorbing.msm(x=NULL, qmatrix=NULL)
```

## Arguments

x	A fitted multi-state model as returned by <a href="#">msm</a> .
qmatrix	A transition intensity matrix. The diagonal is ignored and taken to be minus the sum of the rest of the row.

## Value

A vector of the ordinal indices of the transient or absorbing states.

## Author(s)

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>

---

updatepars.msm	<i>updatepars.msm</i>
----------------	-----------------------

---

**Description**

Update the maximum likelihood estimates in a fitted model object. Developer use only.

**Usage**

```
updatepars.msm(x, pars)
```

**Arguments**

x	A fitted multi-state model object, as returned by <a href="#">msm</a> .
pars	Vector of new parameters, in their untransformed real-line parameterisations, to substitute for the maximum likelihood estimates corresponding to those in the estimates component of the fitted model object ( <a href="#">msm.object</a> ). The order of the parameters is documented in <a href="#">msm</a> , argument <code>fixedpars</code> .

**Value**

An updated [msm](#) model object with the updated maximum likelihood estimates, but with the covariances / standard errors unchanged.

Point estimates from output functions such as [qmatrix.msm](#), [pmatrix.msm](#), or any related function, can then be evaluated with the new parameters, and at arbitrary covariate values.

This function is used, for example, when computing confidence intervals from [pmatrix.msm](#), and related functions, using the `ci="normal"` method.

**Author(s)**

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---

viterbi.msm	<i>Calculate the probabilities of underlying states and the most likely path through them</i>
-------------	---

---

**Description**

For a fitted hidden Markov model, or a model with censored state observations, the Viterbi algorithm recursively constructs the path with the highest probability through the underlying states. The probability of each hidden state is also computed for hidden Markov models.

**Usage**

```
viterbi.msm(x, normboot=FALSE)
```

**Arguments**

x	A fitted hidden Markov multi-state model, or a model with censored state observations, as produced by <a href="#">msm</a>
normboot	If TRUE, then before running the algorithm, the maximum likelihood estimates of the model parameters are replaced by an alternative set of parameters drawn randomly from the asymptotic multivariate normal distribution of the MLEs.

**Value**

A data frame with columns:

subject = subject identification numbers

time = times of observations

observed = corresponding observed states

fitted = corresponding fitted states found by Viterbi recursion. If the model is not a hidden Markov model and there are no censored state observations, this is just the observed states.

For hidden Markov models, an additional matrix pstate is also returned inside the data frame, giving the probability of each hidden state at each point, conditionally on all the data. This is computed by the forward/backward algorithm.

**Author(s)**

C. H. Jackson <[chris.jackson@mrc-bsu.cam.ac.uk](mailto:chris.jackson@mrc-bsu.cam.ac.uk)>

**References**

Durbin, R., Eddy, S., Krogh, A. and Mitchison, G. *Biological sequence analysis*, Cambridge University Press, 1998.

**See Also**

[msm](#)

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