

Package ‘FKF.SP’

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Title Fast Kalman Filtering Through Sequential Processing

Version 0.1.0

Description Fast and flexible Kalman filtering implementation utilizing sequential processing, designed for efficient parameter estimation through maximum likelihood estimation. 'FKF.SP' was built upon the existing 'FKF' package and was designed to generally increase the computational efficiency of Kalman filtering when independence is assumed in the measurement error of observations. Sequential processing is described in the textbook of Durbin and Koopman (2001, ISBN:978-0-19-964117-8).

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URL <https://github.com/TomAspinall/FKF.SP>

BugReports <https://github.com/TomAspinall/FKF.SP/issues>

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fkf.SP

*Fast Kalman Filtering using Sequential Processing.***Description**

The `fkf.SP` function performs fast and flexible Kalman Filtering using Sequential Processing. It is designed for efficient parameter estimation through maximum likelihood estimation. `fkf.SP` wraps the C-function `fkf_SP` which relies upon the linear algebra subroutines of BLAS. Sequential Processing is a univariate treatment of a multivariate series of observations and can benefit from computational efficiencies over traditional Kalman Filtering. `fkf.SP` is based from the `fkf` function of the FKF package, sharing identical arguments except for the `GGt` argument (see **arguments**). The `fkf.SP` function is, in general, a faster Kalman Filtering method than the `fkf` function. `fkf.SP` is compatible with missing observations.

Usage

```
fkf.SP(a0, P0, dt, ct, Tt, Zt, HHT, GGt, yt)
```

Arguments

<code>a0</code>	A vector giving the initial value/estimation of the state variable
<code>P0</code>	A matrix giving the variance of <code>a0</code>
<code>dt</code>	A matrix giving the intercept of the transition equation
<code>ct</code>	A matrix giving the intercept of the measurement equation
<code>Tt</code>	An array giving factor of the transition equation
<code>Zt</code>	An array giving the factor of the measurement equation
<code>HHT</code>	An array giving the variance of the innovations of the transition equation
<code>GGt</code>	A vector giving the diagonal elements of the matrix for the variance of the disturbances of the measurement equation. Covariance between disturbances is not supported under the Sequential Processing method.
<code>yt</code>	A matrix containing the observations. "NA"- values are allowed

Details**Parameters:**

The `fkf.SP` function builds upon the `fkf` function of the FKF package by adjusting the Kalman Filtering algorithm to utilize sequential processing. The `fkf.SP` and `fkf` functions feature highly similar arguments for compatibility purposes; only argument `GGt` has changed from an array type object to a vector or matrix type object. The `fkf.SP` function takes the additional assumption over the `fkf` function that the variance of the disturbances of the measurement equation are independent; a requirement of sequential processing (see below).

The parameters can either be constant or deterministic time-varying. Assume the number of discrete time observations is n i.e. $y = y_t$ where $t = 1, \dots, n$. Let m be the dimension of the state variable and d the dimension of the observations. Then, the parameters admit the following classes and dimensions:

<code>dt</code>	either a $m \times n$ (time-varying) or a $m \times 1$ (constant) matrix.
<code>Tt</code>	either a $m \times m \times n$ or a $m \times m \times 1$ array.

HHT either a $m \times m \times n$ or a $m \times m \times 1$ array.
 ct either a $d \times n$ or a $d \times 1$ matrix.
 Zt either a $d \times m \times n$ or a $d \times m \times 1$ array.
 GGt either a $d \times n$ (time-varying) or a $d \times 1$ matrix.
 yt a $d \times n$ matrix.

State Space Form

The following notation follows that of Koopman *et al.* (1999) and the documentation of the fkf function. The Kalman Filter is characterized by the transition and measurement equations:

$$\alpha_{t+1} = d_t + T_t \cdot \alpha_t + H_t \cdot \eta_t$$

$$y_t = c_t + Z_t \cdot \alpha_t + G_t \cdot \epsilon_t$$

where η_t and ϵ_t are iid $N(0, I_m)$ and iid $N(0, I_d)$, respectively, and α_t denotes the state vector. The parameters admit the following dimensions:

$$\begin{array}{lll}
 a_t \in R^m & d_t \in R^m & \eta_t \in R^m \\
 T_t \in R^{m \times m} & H_t \in R^{m \times m} & \\
 y_t \in R^d & c_t \in R^d & \epsilon_t \in R^d \\
 Z_t \in R^{d \times m} & G_t \in R^{d \times d} &
 \end{array}$$

Note that fkf.SP takes as input HHT and GGt which corresponds to $H_t H_t'$ and $\text{diag}(G_t)^2$ respectively.

Sequential Processing Iteration:

Traditional Kalman Filtering takes the entire observational vector y_t as the items for analysis. Sequential Processing is an alternate approach that filters the elements of y_t one at a time. Let p equal the number of observations at time t (ie. when considering possible missing observations $p \leq d$). This univariate treatment of the multivariate series has the advantage that the function of the covariance matrix, F_t , becomes 1×1 , avoiding the inversion of a $p \times p$ matrix. This can provide computational efficiencies (especially under the case of many observations, ie. p is large)

The SP iteration involves treating the vector series: y_1, \dots, y_n instead as the scalar series $y_{1,1}, \dots, y_{(1,p)}, y_{2,1}, \dots, y_{(n,p)}$. For any time point, the observation vector is given by:

$$y'_t = (y_{(t,1)}, \dots, y_{(t,p)})$$

The filtering equations are written as:

$$a_{t,i+1} = a_{t,i} + K_{t,i} v_{t,i}$$

$$P_{t,i+1} = P_{t,i} - K_{t,i} F_{t,i} K_{t,i}'$$

Where:

$$\hat{y}_{t,i} = c_t + Z_t \cdot a_{t,i}$$

$$v_{t,i} = y_{t,i} - \hat{y}_{t,i}$$

$$F_{t,i} = Z_{t,i} P_{t,i} Z_{t,i}' + G G_{t,i}$$

$$K_{t,i} = P_{t,i} Z_{t,i}' F_{t,i}^{-1}$$

$$i = 1, \dots, p$$

Transition from time t to $t + 1$ occurs through the standard transition equations.

$$\alpha_{t+1,1} = d_t + T_t \cdot \alpha_{t,p}$$

$$P_{t+1,1} = T_t \cdot P_{t,p} \cdot T_t' + H H'$$

The log-likelihood at time t is given by:

$$\log L_t = -\frac{p}{2} \log(2\pi) - \frac{1}{2} \sum_{i=1}^p (\log F_i + \frac{v_i^2}{F_i})$$

Where the log-likelihood of observations is:

$$\log L = \sum_t^n \log L_t$$

There's several distinctions between the Multivariate Kalman Filter and the Sequential processing filtered values. The elements of the innovation vector v_t are not the same as $v_{t,i}$ for $i = 1, \dots, p$. This is also true of the diagonal elements of the variance matrix F_t and the variances $F_{t,i}$, for $i = 1, \dots, p$; only the first diagonal element of F_t is equal to $F_{t,1}$.

Value

A numeric value corresponding to the log-likelihood calculated by the Kalman Filter. Ideal for maximum likelihood estimation through optimization routines such as `optim`.

fkf and fkf.SP values:

Outputs of the `fkf` and `fkf.SP` functions differ slightly in that `fkf` returns a list object of filtered values returned by the algorithm, whereas `fkf.SP` returns only a numeric value corresponding to the log-likelihood returned by the filter. `fkf` is thus appropriate when filtered values are required and `fkf.SP` is appropriate for parameter estimation through maximum likelihood methods.

When there are no missing observations (ie. "NA" values) in matrix `yt`, the return of function `fkf.SP` and the `logLik` object returned within the list of function `fkf` are identical. When there are NA values, however, the log-likelihood score returned by `fkf.SP` is always higher. The log-likelihood score of the FKF C code is instantiated through the calculation of $-n \times d \times \log(2\pi)$, where n is the number of columns of matrix `yt` and d is the number of rows of matrix `yt`. Under the assumption that there are missing observations, d would instead become d_t , where $d_t \leq d \forall t$. Whilst this doesn't influence parameter estimation, because observation matrix `yt` and thus the offset resulting from this is kept constant during maximum likelihood estimation, this does result in low bias of the log-likelihood scores output by the `fkf` function.

References

- Anderson, B. D. O. and Moore, J. B. (1979). *Optimal Filtering* Englewood Cliffs: Prentice-Hall.
- Fahrmeir, L. and tutz, G. (1994) *Multivariate Statistical Modelling Based on Generalized Linear Models*. Berlin: Springer.
- Koopman, S. J., Shephard, N., Doornik, J. A. (1999). Statistical algorithms for models in state space using SsfPack 2.2. *Econometrics Journal*, Royal Economic Society, vol. 2(1), pages 107-160.
- Durbin, James, and Siem Jan Koopman (2012). *Time series analysis by state space methods*. Oxford university press.
- David Luethi, Philipp Erb and Simon Otziger (2018). FKF: Fast Kalman Filter. R package version 0.1.5. <https://CRAN.R-project.org/package=FKF>

Examples

```
# FKF.SP is suitable for parameter estimation, and thus the following examples
# showcase how to estimate parameters of different models.

## -----
##Example 1 - ARMA(2,1) model estimation.
## -----

n <- 1000

## Set the AR parameters

ar1 <- 0.6
ar2 <- 0.2
ma1 <- -0.2
sigma <- sqrt(0.2)

## Sample from an ARMA(2, 1) process
a <- stats::arima.sim(model = list(ar = c(ar1, ar2), ma = ma1), n = n,
                        innov = rnorm(n) * sigma)

## Create a state space representation out of the four ARMA parameters
arma21ss <- function(ar1, ar2, ma1, sigma) {
  Tt <- matrix(c(ar1, ar2, 1, 0), ncol = 2)
  Zt <- matrix(c(1, 0), ncol = 2)
  ct <- matrix(0)
  dt <- matrix(0, nrow = 2)
  GGt <- matrix(0)
  H <- matrix(c(1, ma1), nrow = 2) * sigma
  HHt <- H %*% t(H)
  a0 <- c(0, 0)
  P0 <- matrix(1e6, nrow = 2, ncol = 2)
  return(list(a0 = a0, P0 = P0, ct = ct, dt = dt, Zt = Zt, Tt = Tt, GGt = GGt,
              HHt = HHt))
}

## The objective function passed to 'optim'
objective <- function(theta, yt, SP = F) {
  sp <- arma21ss(theta["ar1"], theta["ar2"], theta["ma1"], theta["sigma"])
  ans <- fkf.SP(a0 = sp$a0, P0 = sp$P0, dt = sp$dt, ct = sp$ct, Tt = sp$Tt,
                Zt = sp$Zt, HHt = sp$HHt, GGt = sp$GGt, yt = yt)
  return(-ans)
}

#Estimate Parameters using the optim function:
theta <- c(ar = c(0, 0), ma1 = 0, sigma = 1)
ARMA_MLE <- optim(theta, objective, yt = rbind(a), hessian = TRUE, SP = T)

## -----
##Example 2 - Nile Example:
## -----
#Local level model for the Nile's annual flow:

## Transition equation:
```

```

## alpha[t+1] = alpha[t] + eta[t], eta[t] ~ N(0, HHT)
## Measurement equation:
## y[t] = alpha[t] + eps[t], eps[t] ~ N(0, GGt)

yt <- Nile

##Incomplete Nile Data - two NA's are present:
yt[c(3, 10)] <- NA

## Set constant parameters:
dt <- ct <- matrix(0)
Zt <- Tt <- matrix(1)
a0 <- yt[1] # Estimation of the first year flow
P0 <- matrix(100) # Variance of 'a0'

## Parameter estimation - maximum likelihood estimation:
##Unknown parameters initial estimates:
GGt <- HHT <- var(yt, na.rm = TRUE) * .5
set.seed(1)
#Perform Maximum Likelihood Estimation
Nile_MLE <- suppressWarnings(optim(c(HHT = HHT, GGt = GGt),
                                fn = function(par, ...)
                                -fkf.SP(HHT = matrix(par[1]), GGt = matrix(par[2]), ...),
                                yt = rbind(yt), a0 = a0, P0 = P0, dt = dt, ct = ct,
                                Zt = Zt, Tt = Tt))

## <-----
#Example 3 - Dimensionless Treering Example:
## <-----

## Transition equation:
## alpha[t+1] = alpha[t] + eta[t], eta[t] ~ N(0, HHT)
## Measurement equation:
## y[t] = alpha[t] + eps[t], eps[t] ~ N(0, GGt)

## tree-ring widths in dimensionless units
y <- treering

## Set constant parameters:
dt <- ct <- matrix(0)
Zt <- Tt <- matrix(1)
a0 <- y[1] # Estimation of the first width
P0 <- matrix(100) # Variance of 'a0'

##Time comparison - Estimate parameters 10 times:
Treering_MLE <- suppressWarnings(optim(c(HHT = var(y, na.rm = TRUE) * .5,
                                GGt = var(y, na.rm = TRUE) * .5),
                                fn = function(par, ...)
                                -fkf.SP(HHT = array(par[1],c(1,1,1)), GGt = matrix(par[2]), ...),
                                yt = rbind(y), a0 = a0, P0 = P0, dt = dt, ct = ct,
                                Zt = Zt, Tt = Tt))

##Not run - Filter tree ring data with estimated parameters using the FKF package:
#fkf.obj <- fkf(a0, P0, dt, ct, Tt, Zt, HHT = array(fit.fkf$par[1],c(1,1,1)),
#              GGt = matrix(fit.fkf$par[2]), yt = rbind(y))

```

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