

Package ‘NFCP’

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Title N-Factor Commodity Pricing Through Term Structure Estimation

Version 1.0.1

Description Commodity pricing models are (systems of) stochastic differential equations that are utilized for the valuation and hedging of commodity contingent claims (i.e. derivative products on the commodity) and other commodity related investments. Commodity pricing models that capture market dynamics are of great importance to commodity market participants in order to exercise sound investment and risk-management strategies. Parameters of commodity pricing models are estimated through maximum likelihood estimation, using available term structure futures data of a commodity. 'NFCP' (n-factor commodity pricing) provides a framework for the modeling, parameter estimation, probabilistic forecasting, option valuation and simulation of commodity prices through state space and Monte Carlo methods, risk-neutral valuation and Kalman filtering. 'NFCP' allows the commodity pricing model to consist of n correlated factors, with both random walk and mean-reverting elements. The n-factor commodity pricing model framework was first presented in the work of Cortazar and Naranjo (2006) <doi:10.1002/fut.20198>. Examples presented in 'NFCP' replicate the two-factor crude oil commodity pricing model presented in the prolific work of Schwartz and Smith (2000) <doi:10.1287/mnsc.46.7.893.12034> with the approximate term structure futures data applied within this study provided in the 'NFCP' package.

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American_option_value *N-factor model American put option pricing*

Description

Value American put options under the parameters of an N-factor model through the Least-Squares Monte Carlo (LSM) Simulation Method. This function is a wrapper to the 'LSM_American_option' function of the 'LSMRealOptions' package.

Usage

```
American_option_value(
  x_0,
  parameters,
  N_simulations,
  option_maturity,
  dt,
  K,
  r,
  orthogonal = "Power",
  degree = 2,
  verbose = FALSE,
  debugging = FALSE
)
```

Arguments

<code>x_0</code>	Initial values of the state vector.
<code>parameters</code>	Named vector of parameter values of a specified N-factor model. Function NFCP_parameters is recommended.
<code>N_simulations</code>	total number of simulated price paths

option_maturity	Time to expiration of the option (in years)
dt	discrete time step of simulation
K	Strike price of the American put option
r	Risk-free interest rate.
orthogonal	The orthogonal polynomial used to approximate the continuation value of the option in the LSM simulation method. Orthogonal polynomial arguments available are: "Power", "Laguerre", "Jacobi", "Legendre", "Chebyshev", "Hermite". See help(LSM.AmericanOption)
degree	The degree of polynomials used in the least squares fit. See help(LSM.AmericanOption)
verbose	logical Should additional information be output?
debugging	logical Should additional simulation information be output?

Details

The 'American_option_value' function is a wrapper to the 'spot_price_simulate' and 'LSM_American_option' of the 'LSMRealOptions' package that returns the value of American put options under a given N-factor model.

The least-squares Monte Carlo (LSM) simulation method is an option valuation method first presented by Longstaff and Schwartz (2001) that approximates the value of American options.

Methods to solve for the value of options with early exercise opportunities include partial differential equations, lattice-based methods and Monte-Carlo simulation. LSM simulation is the optimal solution method to value American options under an N-factor model due to the multiple factors that can make up the spot price process and influence the option value. Whilst lattice and partial differential equation approaches suffer from the 'curse of dimensionality', LSM simulation may be readily applied under multi-factor settings.

Longstaff and Schwartz (2001) state that as the conditional expectation of the continuation value belongs to a Hilbert space, it can be represented by a combination of orthogonal basis functions. Increasing the number of stochastic state variables therefore increases the number of required basis functions exponentially.

Value

The 'American_option_value' function by default returns a numeric object corresponding to the calculated value of the American put option.

When verbose = T, 6 objects are returned within a list class object. The objects returned are:

Value	The calculated option value.
Standard Error	The standard error of the calculated option value.
Expected Timing	The expected time of early exercise..
Expected Timing SE	The standard error of the expected time of early exercise.
Exercise Probability	The probability of early exercise of the option being exercised.
Cumulative Exercise Probability	vector. The cumulative probability of option exercise at each discrete observation

When debugging = T, an additional 2 objects are returned within the list class object. These are the results output by both the 'Spot.Price.Simulate' and 'LSM.AmericanOption' of the 'LSMRealOptions' package respectively. The objects returned are:

State_Variables A matrix of simulated state variables for each factor is returned when verbose = T. The number of fa

Prices A matrix of simulated price paths. Each column represents one simulated price path and each row represents one time step.

References

Longstaff, F.A., and E.S. Schwartz. 2001. Valuing American Options by Simulation: A Simple Least-Squares Approach. *The Review of Financial Studies*. 14:113-147.

Schwartz, E. S., and J. E. Smith, (2000). Short-Term Variations and Long-Term Dynamics in Commodity Prices. *Manage. Sci.*, 46, 893-911.

Cortazar, G., and L. Naranjo, (2006). An N-factor Gaussian model of oil futures prices. *Journal of Futures Markets: Futures, Options, and Other Derivative Products*, 26(3), 243-268.

Aspinall, T., A. Gepp, G. Harris, S. Kelly, C. Southam, and B. Vanstone, (2021). LSMRealOptions: Value American and Real Options Through LSM Simulation. R package version 0.1.0.

Examples

```
# Example 1 - An American put option on a stock following 'GBM'
# growing at the risk-free rate:
American_option_value(x_0 = log(36),
                      parameters = c(mu_rn = (0.06 - (1/2) * 0.2^2), sigma_1 = 0.2),
                      N_simulations = 1e2,
                      option_maturity = 1,
                      dt = 1/50,
                      K = 40,
                      r = 0.05,
                      verbose = FALSE,
                      orthogonal = "Laguerre",
                      degree = 3)

# Example 2 - An American put option under a two-factor crude oil model:

## Step 1 - Obtain current (i.e. most recent) state vector by filtering the
## two-factor oil model:
Schwartz_Smith_oil <- NFCP_Kalman_filter(parameter_values = SS_oil$two_factor,
                                         parameter_names = names(SS_oil$two_factor),
                                         log_futures = log(SS_oil$stitched_futures),
                                         dt = SS_oil$dt,
                                         futures_TTM = SS_oil$stitched_TTM,
                                         verbose = TRUE)

## Step 2 - Calculate 'put' option price:
American_option_value(x_0 = Schwartz_Smith_oil$x_t,
                      parameters = SS_oil$two_factor,
                      N_simulations = 1e2,
                      option_maturity = 1,
                      dt = 1/12,
                      K = 20,
                      r = 0.05,
                      verbose = FALSE,
                      orthogonal = "Power",
                      degree = 2)
```

European_option_value *N-factor model European option pricing*

Description

Value European Option Put and Calls under the parameters of an N-factor model.

Usage

```
European_option_value(
  x_0,
  parameters,
  futures_maturity,
  option_maturity,
  K,
  r,
  call,
  verbose = FALSE
)
```

Arguments

x_0	Initial values of the state vector.
parameters	Named vector of parameter values of a specified N-factor model. Function NFCP_parameters is recommended.
futures_maturity	Time, in years, when the underlying futures contract matures.
option_maturity	Time, in years, when the European option expires.
K	Strike price of the European Option
r	Risk-free interest rate.
call	logical is the European option a call or put option?
verbose	logical. Should additional information be output? see details

Details

The European_option_value function calculates analytic expressions of the value of European call and put options on futures contracts within the N-factor model. Under the assumption that future futures prices are log-normally distributed under the risk-neutral process, there exist analytic expressions of the value of European call and put options on futures contracts. The following analytic expression follows from that presented by Schwartz and Smith (2000) extended to the N-factor framework. The value of a European option on a futures contract is given by calculating its expected future value using the risk-neutral process and subsequently discounting at the risk-free rate.

One can verify that under the risk-neutral process, the expected futures price at time t is:

$$E^*[F_{T,t}] = \exp\left(\sum_{i=1}^N e^{-\kappa_i T} x_i(0) + \mu^* t + A(T-t) + \frac{1}{2}(\sigma_1^2 t + \sum_{i,j \neq 1} e^{-(\kappa_i + \kappa_j)(T-t)} \sigma_i \sigma_j \rho_{i,j} \cdot \frac{1 - e^{-(\kappa_i + \kappa_j)t}}{\kappa_i + \kappa_j})\right) \equiv F_{T,0}$$

This follows from the derivation provided within the vignette of the NFCP package as well as the details of the futures_price_forecast package. The equality of expected futures price at time t being equal to the time- t current futures price $F_{T,0}$ is proven by Futures prices being given by expected spot prices under the risk-neutral process ($F_{T,t} = E_t^* [S_T]$) and the law of iterated expectations ($E^* [E_t^* [S_T]] = E^* [S_T]$)

Because future futures prices are log-normally distributed under the risk-neutral process, we can write a closed-form expression for valuing European put and call options on these futures. When $T = 0$ these are European options on the spot price of the commodity itself. The value of a European call option on a futures contract maturing at time T , with strike price K , and with time t until the option expires, is:

$$\begin{aligned} & e^{-rt} E^* [\max (F_{T,t} - K, 0)] \\ &= e^{-rt} (F_{T,0} N(d) - K N(d - \sigma_\phi(t, T))) \end{aligned}$$

Where:

$$d = \frac{\ln(F/K)}{\sigma_\phi(t, T)} + \frac{1}{2} \sigma_\phi(t, T)$$

and:

$$\sigma_\phi(t, T) = \sqrt{\sigma_1^2 t + \sum_{i,j \neq 1} e^{-(\kappa_i + \kappa_j)(T-t)} \sigma_i \sigma_j \rho_{i,j} \cdot \frac{1 - e^{-(\kappa_i + \kappa_j)t}}{\kappa_i + \kappa_j}}$$

Parameter $N(d)$ indicates cumulative probabilities for the standard normal distribution (i.e. $P(Z < d)$).

Similarly, the value of a European put with the same parameters is given by:

$$\begin{aligned} & e^{-rt} E^* [\max (K - F_{T,t}, 0)] \\ &= e^{-rt} (-F_{T,0} N(-d) + K N(\sigma_\phi(t, T) - d)) \end{aligned}$$

The presented option valuation formulas are analogous to the Black-Scholes formulas for valuing European options on stocks that do not pay dividends

Under this terminology, the stock price corresponds to the present value of the futures commitment ($e^{-rt} F_{T,0}$) and the equivalent annualized volatility would be $\sigma_\phi(t, T)/\sqrt{t}$

When verbose = T, the European_option_value function numerically calculates the sensitivity of option prices to the underlying parameters specified within the N-factor model, as well as some of the most common "Greeks" related to European put and call option pricing. All gradients are calculated numerically by calling the grad function from the numDeriv package.

Value

The European_option_value function returns a numeric value corresponding to the present value of an option when verbose = F. When verbose = T, European_option_value returns a list with three objects:

option value	Present value of the option.
annualized volatility	Annualized volatility of the option.
parameter sensitivity	Sensitivity of the option value to each parameter of the N-factor model.
greeks	Sensitivity of the option value to different option parameters.

References

Schwartz, E. S., and J. E. Smith, (2000). Short-Term Variations and Long-Term Dynamics in Commodity Prices. *Manage. Sci.*, 46, 893-911.

Cortazar, G., and L. Naranjo, (2006). An N-factor Gaussian model of oil futures prices. *Journal of Futures Markets: Futures, Options, and Other Derivative Products*, 26(3), 243-268.

Examples

```
##Example 1 - A European 'put' option on a stock following 'GBM'
##growing at the risk-free rate:

### Risk-free rate:
rf <- 0.05
### Stock price:
S_0 <- 20
### Stock volatility:
S_sigma <- 0.2
### Option maturity:
Tt <- 1
### Exercise price:
K <- 20
### Calculate 'put' option price:
European_option_value(x_0 = log(S_0), parameters = c(mu_rn = rf, sigma_1 = S_sigma),
                      futures_maturity = Tt, option_maturity = 1,
                      K = K, r = rf, call = FALSE, verbose = TRUE)

##Example 2 - A European call option under a two-factor crude oil model:

##Step 1 - Obtain current (i.e. most recent) state vector by filtering the
##two-factor oil model:
Schwartz_Smith_oil <- NFCP_Kalman_filter(parameter_values = SS_oil$two_factor,
                                         parameter_names = names(SS_oil$two_factor),
                                         log_futures = log(SS_oil$stitched_futures),
                                         dt = SS_oil$dt,
                                         futures_TTM = SS_oil$stitched_TTM,
                                         verbose = TRUE)

##Step 2 - Calculate 'call' option price:
European_option_value(x_0 = Schwartz_Smith_oil$x_t,
                      parameters = SS_oil$two_factor,
                      futures_maturity = 1,
                      option_maturity = 1,
                      K = 20,
                      r = 0.05,
                      call = TRUE,
                      verbose = FALSE)
```

Description

Analytically forecast future expected Futures prices under the risk-neutral version of a specified N-factor model.

Usage

```
futures_price_forecast(
  x_0,
  parameters,
  t = 0,
  futures_TTM = 1:10,
  percentiles = NULL
)
```

Arguments

x_0	Initial values of the state vector.
parameters	A named vector of parameter values of a specified N-factor model. Function NFCP_parameters is recommended.
t	a numeric specifying the time point at which to forecast futures prices
futures_TTM	a vector specifying the time to maturity of futures contracts to value.
percentiles	Optional. A vector of percentiles to include probabilistic forecasting intervals.

Details

Under the assumption of risk-neutrality, futures prices are equal to the expected future spot price. Additionally, under deterministic interest rates, forward prices are equal to futures prices. Let $F_{T,t}$ denote the market price of a futures contract at time t with time T until maturity. let $*$ denote the risk-neutral expectation and variance of futures prices. The following equations assume that the first factor follows a Brownian Motion.

$$E^*[ln(F_{T,t})] = \sum_{i=1}^N e^{-\kappa_i T} x_i(0) + \mu^* t + A(T-t)$$

Where:

$$A(T-t) = \mu^*(T-t) - \sum_{i=1}^N \frac{1 - e^{-\kappa_i(T-t)} \lambda_i}{\kappa_i} + \frac{1}{2} (\sigma_1^2(T-t) + \sum_{i,j \neq 1} \sigma_i \sigma_j \rho_{i,j} \frac{1 - e^{-(\kappa_i + \kappa_j)(T-t)}}{\kappa_i + \kappa_j})$$

The variance is given by:

$$Var^*[ln(F_{T,t})] = \sigma_1^2 t + \sum_{i,j \neq 1} e^{-(\kappa_i + \kappa_j)(T-t)} \sigma_i \sigma_j \rho_{i,j} \frac{1 - e^{-(\kappa_i + \kappa_j)t}}{\kappa_i + \kappa_j}$$

Value

futures_price_forecast returns a vector of expected Futures prices under a given N-factor model with specified time to maturities at time t . When percentiles are specified, the function returns a matrix with the corresponding confidence bands in each column of the matrix.

References

Schwartz, E. S., and J. E. Smith, (2000). Short-Term Variations and Long-Term Dynamics in Commodity Prices. *Manage. Sci.*, 46, 893-911.

Cortazar, G., and L. Naranjo, (2006). An N-factor Gaussian model of oil futures prices. *Journal of Futures Markets: Futures, Options, and Other Derivative Products*, 26(3), 243-268.

Examples

```
# Forecast futures prices of the Schwartz and Smith (2000) two-factor oil model:
## Step 1 - Run the Kalman filter for the two-factor oil model:
SS_2F_filtered <- NFCP_Kalman_filter(parameter_values = SS_oil$two_factor,
                                     parameter_names = names(SS_oil$two_factor),
                                     log_futures = log(SS_oil$stitched_futures),
                                     dt = SS_oil$dt,
                                     futures_TTM = SS_oil$stitched_TTM,
                                     verbose = TRUE)

## Step 2 - Probabilistic forecast of the risk-neutral two-factor
## stochastic differential equation (SDE):
futures_price_forecast(x_0 = SS_2F_filtered$x_t,
                      parameters = SS_oil$two_factor,
                      t = 0,
                      futures_TTM = seq(0,9,1/12),
                      percentiles = c(0.1, 0.9))
```

futures_price_simulate

Simulate futures prices of an N-factor model through Monte Carlo simulation

Description

Simulate Futures price data with dynamics that follow the parameters of an N-factor model through Monte Carlo simulation.

Usage

```
futures_price_simulate(
  x_0,
  parameters,
  dt,
  N_obs,
  futures_TTM,
  ME_TTM = NULL,
  verbose = TRUE
)
```

Arguments

x_0	Initial values of the state vector.
parameters	A named vector of parameter values of a specified N-factor model. Function NFCP_parameters is recommended.
dt	discrete time step of simulation
N_obs	The number of observations to simulate
futures_TTM	A vector or matrix of the time to maturity of futures contracts to simulate. See details
ME_TTM	vector of maturity groupings to consider for simulated futures prices. The length of ME_TTM must be equal to the number of 'ME' parameter values.
verbose	logical. Should the simulated state variables and associated prices be output?

Details

The `futures_price_simulate` function simulates futures price data using the Kalman Filter algorithm, drawing from a normal distribution for the shocks in the transition and measurement equations at each discrete time step. At each discrete time point, an observation of the state vector is generated through the transition equation, drawing from a normal distribution with a covariance equal to Q_t . Following this, simulated futures prices are generated through the measurement equation, drawing from a normal distribution with covariance matrix equal to H .

Input `futures_TTM` can be either a matrix specifying the constant time to maturity of futures contracts to simulate, or it can be a matrix where `nrow(futures_TTM) == N_obs` for the time-varying time to maturity of the futures contracts to simulate. This allows for the simulation of both aggregate stitched data and individual futures contracts.

Value

`futures_price_simulate` returns a list with three objects when `verbose = T` and a matrix of simulated futures prices when `verbose = F`. The list objects returned are:

#'

state_vector	A matrix of Simulated state variables at each discrete time point. The columns represent each factor
futures_prices	A matrix of Simulated futures prices, with each column representing a simulated futures contract.
spot_prices	A vector of simulated spot prices

References

Schwartz, E. S., and J. E. Smith, (2000). Short-Term Variations and Long-Term Dynamics in Commodity Prices. *Manage. Sci.*, 46, 893-911.

Cortazar, G., and L. Naranjo, (2006). An N-factor Gaussian model of oil futures prices. *Journal of Futures Markets: Futures, Options, and Other Derivative Products*, 26(3), 243-268.

Examples

```
##Example 1 - Simulate Crude Oil Stitched futures prices
##under a Two-Factor model, assuming a constant time to maturity:
```

```

simulated_futures <- futures_price_simulate(x_0 = c(log(SS_oil$spot[1,1]), 0),
                                           parameters = SS_oil$two_factor,
                                           dt = SS_oil$dt,
                                           N_obs = nrow(SS_oil$stitched_futures),
                                           futures_TTM = SS_oil$stitched_TTM)

##Example 2 - Simulate Crude Oil Contract Prices under a Two-Factor model,
##using a rolling-window of measurement errors:

simulated_futures_prices <- futures_price_simulate(x_0 = c(log(SS_oil$spot[1,1]), 0),
                                                  parameters = SS_oil$two_factor,
                                                  dt = SS_oil$dt,
                                                  N_obs = nrow(SS_oil$contracts),
                                                  futures_TTM = SS_oil$contract_maturities,
                                                  ME_TTM = c(1/4, 1/2, 1, 2, 5))

```

NFCP_domains

N-Factor MLE search boundaries

Description

Generate boundaries for the domain of parameters of the N-factor model for parameter estimation.

Usage

```

NFCP_domains(
  parameters,
  kappa = NULL,
  lambda = NULL,
  sigma = NULL,
  mu = NULL,
  mu_rn = NULL,
  rho = NULL,
  ME = NULL,
  x_0 = NULL,
  E = NULL
)

```

Arguments

parameters	a vector of parameter names of an N-factor model. Function NFCP_parameters is recommended.
kappa	A vector of length two specifying the lower and upper bounds for the 'kappa' parameter
lambda	A vector of length two specifying the lower and upper bounds for the 'lambda' parameter
sigma	A vector of length two specifying the lower and upper bounds for the 'sigma' parameter
mu	A vector of length two specifying the lower and upper bounds for the 'mu' parameter

mu_rn	A vector of length two specifying the lower and upper bounds for the 'mu_rn' parameter
rho	A vector of length two specifying the lower and upper bounds for the 'rho' parameter
ME	A vector of length two specifying the lower and upper bounds for the 'ME' (i.e., measurement error) parameter
x_0	A vector of length two specifying the lower and upper bounds for the 'x_0' parameter
E	A vector of length two specifying the lower and upper bounds for the 'E' parameter

Details

The NFCP_domains function generates lower and upper bounds for the parameter estimation procedure in the format required of the 'Domains' argument of the 'genoud' function. NFCP_domains allows easy setting of custom boundaries for parameter estimation, whilst also providing default domains of parameters.

Value

A matrix of defaulted domains for the given unknown parameters. The first column corresponds to the lower bound of the allowable search space for the parameter, whilst the second column corresponds to the upper bound. These values were set to allow for the 'realistic' possible values of given parameters as well as restricting some parameters (such as variance and mean-reverting terms) from taking negative values. The format of the returned matrix matches that required by the Domains argument of the Genoud function from the package RGenoud.

References

- Mebane, W. R., and J. S. Sekhon, (2011). Genetic Optimization Using Derivatives: The rgenoud Package for R. *Journal of Statistical Software*, 42(11), 1-26. URL <http://www.jstatsoft.org/v42/i11/>.
- Schwartz, E. S., and J. E. Smith, (2000). Short-Term Variations and Long-Term Dynamics in Commodity Prices. *Manage. Sci.*, 46, 893-911.

Examples

```
##Specify the Schwartz and Smith (2000) two-factor model
##with fixed measurement error:
parameters_2F <- NFCP_parameters(N_factors = 2,
                                GBM = TRUE,
                                initial_states = TRUE,
                                N_ME = 1)

###Generate the default 'domains' argument of 'NFCP_MLE' function:
NFCP_MLE_bounds <- NFCP_domains(parameters_2F)
```

NFCP_Kalman_filter	<i>Filter an N-factor commodity pricing model through the Kalman filter</i>
--------------------	---

Description

Given a set of parameters of the N-factor model, filter term structure data using the Kalman filter.

Usage

```
NFCP_Kalman_filter(
  parameter_values,
  parameter_names,
  log_futures,
  dt,
  futures_TTM,
  ME_TTM = NULL,
  verbose = FALSE,
  debugging = FALSE
)
```

Arguments

parameter_values	Vector of parameter values of an N-factor model. The NFCP_Kalman_filter function is designed for application to optim type functions, and thus parameter values and corresponding parameters different inputs within the function.
parameter_names	Vector of parameter names. Each element of parameter_names must correspond to its respective value element in object parameter_values.
log_futures	Object of class matrix corresponding to the natural logarithm of observable futures prices. NA's are allowed within the matrix. Every column of the matrix must correspond to a particular futures contract, with each row corresponding to a quoted price on a given date.
dt	Constant, discrete time step of observations, in years.
futures_TTM	Object of class 'vector' or 'matrix' that specifies the time to maturity of observed futures contracts. time to maturity can be either constant (ie. class 'vector') or time homogeneous (ie. class 'matrix'). When the time to maturity of observed futures contracts is time homogeneous, the dimensions of futures_TTM must be identical to that of log_futures. Every element of futures_TTM corresponds to the time to maturity, in years, of a futures contract at a given observation date.
ME_TTM	vector of maturity groupings to consider for observed futures prices. The length of ME_TTM must be equal to the number of 'ME' parameter values, and the maximum of ME_TTM must be greater than the maximum observed time-to-maturity of a futures contract. When the number of 'ME' parameter values is equal to one or the total number of contracts (i.e., columns of log_futures), this argument is optional and not considered. The measurement error of an observation is highly influenced by its time-to-maturity, see details .
verbose	logical. Should additional information be output? see values . When verbose = F, the NFCP_Kalman_filter function is significantly faster, see details
debugging	logical. Should additional filtering information be output? see values

Details

NFCP_Kalman_filter applies the Kalman filter algorithm for observable log_futures prices against the input parameters of an N-factor model provided through the parameter_values and parameter_names input vectors.

The NFCP_Kalman_filter function is designed for subsequent input into optimization functions and is called within the N-factor parameter estimation function NFCP_MLE. The first input to the NFCP_Kalman_filter function is a vector of parameters of an N-factor model, with elements of this vector corresponding to the parameter names within the elements of input vector parameter_names. When logical input verbose = F, the NFCP_Kalman_filter function calls the fkf_SP function of the FKF_SP package, which itself is a wrapper of a routine of the Kalman filter written in C utilizing Sequential Processing for maximum computational efficiency (see fkf_SP for more details). When verbose = T, the NFCP_Kalman_filter instead applies a Kalman filter algorithm written in base R and outputs several other list objects, including filtered values and measures for model fit and robustness (see **Returns**)

The N-factor model The N-factor model was first presented in the work of Cortazar and Naranjo (2006, equations 1-3). The N-factor framework describes the spot price process of a commodity as the correlated sum of N state variables x_t .

When GBM = TRUE:

$$\log(S_t) = \sum_{i=1}^N x_{i,t}$$

When GBM = FALSE:

$$\log(S_t) = E + \sum_{i=1}^N x_{i,t}$$

Additional factors within the spot-price process are designed to result in additional flexibility, and possibly fit to the observable term structure, in the spot price process of a commodity. The fit of different N-factor models, represented by the log-likelihood can be directly compared with statistical testing possible through a chi-squared test.

Flexibility in the spot price under the N-factor framework allows the first factor to follow a Brownian Motion or Ornstein-Uhlenbeck process to induce a unit root. In general, an N-factor model where GBM = T allows for non-reversible behaviour within the price of a commodity, whilst GBM = F assumes that there is a long-run equilibrium that the commodity price will revert to in the long-term.

State variables are thus assumed to follow the following processes:

When GBM = TRUE:

$$dx_{1,t} = \mu^* dt + \sigma_1 dw_{1,t}$$

When GBM = FALSE:

$$dx_{1,t} = -(\lambda_1 + \kappa_1 x_{1,t})dt + \sigma_1 dw_{1,t}$$

And:

$$dx_{i,t} =_{i \neq 1} -(\lambda_i + \kappa_i x_{i,t})dt + \sigma_i dw_{i,t}$$

where:

$$E(w_i)E(w_j) = \rho_{i,j}$$

The following constant parameters are defined as:

param μ : long-term growth rate of the Brownian Motion process.

param E : Constant equilibrium level.

param $\mu^* = \mu - \lambda_1$: Long-term risk-neutral growth rate

param λ_i : Risk premium of state variable i .

param κ_i : Reversion rate of state variable i .

param σ_i : Instantaneous volatility of state variable i .

param $\rho_{i,j} \in [-1, 1]$: Instantaneous correlation between state variables i and j .

Disturbances - Measurement Error:

The Kalman filtering algorithm assumes a given measure of measurement error or disturbance in the measurement equation (ie. matrix H). Measurement errors can be interpreted as error in the model's fit to observed prices, or as errors in the reporting of prices (Schwartz and Smith, 2000). These disturbances are typically assumed independent.

var ME_i measurement error of contract i .

where the measurement error of futures contracts ME_i is equal to 'ME_' [i] (i.e. 'ME_1', 'ME_2', ...) specified in arguments parameter_values and parameter_names.

There are three particular cases on how the measurement error of observations can be treated in the NFCP_Kalman_filter function:

Case 1: Only one ME is specified. The Kalman filter assumes that the measurement error of observations are independent and identical.

Case 2: One ME is specified for every observed futures contract. The Kalman filter assumes that the measurement error of observations are independent and unique.

Case 3: A series of ME's are specified for a given grouping of maturities of futures contracts. The Kalman filter assumes that the measurement error of observations are independent and unique to their respective time-to-maturity.

Grouping of maturities for case 3 is specified through the ME_TTM argument. This is a vector that specifies the maximum maturity to consider for each respective ME parameter argument.

in other words, ME_1 is considered for observations with TTM less than ME_TTM[1], ME_2 is considered for observations with TTM less than ME_TTM[2], ..., etc.

The first case is clearly the simplest to estimate, but can be a restrictive assumption. The second case is clearly the most difficult to estimate, but can be an infeasible assumption when considering all available futures contracts that make up the term structure of a commodity.

Case 3 thus serves to ease the restriction of case 1, and allow the user to make the modeling of measurement error as simple or complex as desired for a given set of maturities.

Kalman Filtering

The following section describes the Kalman filter equations used to filter the N-factor model.

The Kalman filter iteration is characterised by a transition and measurement equation. The transition equation develops the vector of state variables between discretised time steps (whilst considering a given level of covariance between state variables over time). The measurement equation relates the unobservable state vector to a vector of observable measurements (whilst also considering a given level of measurement error). The typical Kalman filter algorithm is a Gaussian process state space model.

Transition Equation:

$$\hat{x}_{t|t-1} = c_t + G_t \hat{x}_{t-1} + Q_t \eta_t$$

Measurement Equation:

$$\hat{y}_t = d_t + Z_t \hat{x}_{t|t-1} + H_t \epsilon_t$$

$$t = 1, \dots, n$$

Where η_t and ϵ_t are IID $N(0, I(m))$ and iid $N(0, I(d))$ respectively.

The state vector follows a normal distribution, $x_1 \sim N(a_1, P_1)$, with a_1 and P_1 as the mean vector and variance matrix of the initial state vector x_1 , respectively.

The Kalman filter can be used for parameter estimation through the maximization of the Log-Likelihood value. See NFCP_MLE.

Filtering the N-factor model

let m represent the number of observations at time t

let n represent the number of factors in the N-factor model

observable futures prices: $y_t = [\ln(F(t, T_1)), \ln(F(t, T_2)), \dots, \ln(F(t, T_m))]'$

State vector: $x_t = [x_{1t}, x_{2t}, \dots, x_{nt}]'$

Measurement error: $\text{diag}(H) = [ME_1^2, ME_2^2, \dots, ME_n^2]$

When the number of specified ME terms is one, $s_1 = s_2 = \dots = s_n = ME_1^2$

var Z is an $m \times n$ matrix, where each element $[i, j]$ is equal to:

$$Z_{i,j} = e^{-\kappa_i T_j}$$

var d_t is an $m \times 1$ vector:

$$d_t = [A(T_1), A(T_2), \dots, A(T_m)]'$$

Under the assumption that Factor 1 follows a Brownian Motion, $A(T)$ is given by:

$$A(T) = \mu^* T - \sum_{i=1}^N \frac{1 - e^{-\kappa_i T} \lambda_i}{\kappa_i} + \frac{1}{2} (\sigma_1^2 T + \sum_{i,j \neq 1} \sigma_i \sigma_j \rho_{i,j} \frac{1 - e^{-(\kappa_i + \kappa_j)T}}{\kappa_i + \kappa_j})$$

var v_t is a $n \times 1$ vector of serially uncorrelated Guassian disturbances with $E(V_t) = 0$ and $\text{cov}(v_t) = R^2$

Where:

$$\text{diag}(G_t) = [e^{-\kappa_1 \tau}, e^{-\kappa_2 \tau}, \dots, e^{-\kappa_n \tau}]$$

Where $\tau = T - t$

var w_t is an $n \times 1$ vector of serially uncorrelated Guassian disturbances where:

$$E(w_t) = 0$$

and $\text{cov}(w_t) = Q_t$

var $c_t = [\mu \Delta t, 0, \dots, 0]'$ is an $N \times 1$ vector of the intercept of the transition equation.

var Q_t is equal to the covariance function, given by:

$$\begin{aligned} \text{Cov}_{1,1}(x_{1,t}, x_{1,t}) &= \sigma_1^2 t \\ \text{Cov}_{i,j}(x_{i,t}, x_{j,t}) &= \sigma_i \sigma_j \rho_{i,j} \frac{1 - e^{-(\kappa_i + \kappa_j)t}}{\kappa_i + \kappa_j} \end{aligned}$$

(see also cov_func)

Penalising poorly specified models

The Kalman filter returns non-real log-likelihood scores when the function of the covariance matrix becomes singular or its determinant becomes negative. This occurs when a poorly specified parameter set is input. Non-real log-likelihood scores can break optimization algorithms. To circumvent this, the NFCP_Kalman_filter returns a heavily penalized log-likelihood score whilst also returning a warning. Penalized log-likelihood scores are calculated by:

```
stats::runif(1,-1.5e6,-1e6)
```

Diffuse Kalman filtering

If the initial values of the state vector are not supplied within the parameter_names and parameter_values vectors (ie. Initial.State = F within the NFCP_parameters function), a 'diffuse' assumption is used within the Kalman filtering algorithm. Factors that follow an Ornstein-Uhlenbeck are assumed to equal zero. The initial value of factors that follow a Brownian motion are assumed equal to the first element of log_futures. This is an assumption that the initial estimate of the spot price is equal to the closest to maturity observed futures price.

The initial covariance of the state vector for the Kalman filtering algorithm assumed to be equal to matrix Q

Initial states of factors that follow an Ornstein-Uhlenbeck have a transient effect on future observations, however the initial value of a random walk variable persists across observations and therefore influencing model fit more (see Schwartz and Smith (2000) for more details).

Value

NFCP_Kalman_filter returns a numeric object when verbose = F, which corresponds to the log-likelihood of observations. When verbose = T, the NFCP_Kalman_filter function returns a list object of length seven with the following objects:

LL	Log-Likelihood of observations
X.t	vector. The final observation of the state vector
X	matrix. All observations of the state vector, after the updating equation has been applied
Y	matrix. Estimated futures prices at each observation
V	matrix. Estimation error of each futures contracts at each observation
Filtered Error	matrix. positive mean error (high bias), negative mean error (low bias), mean error (zero bias)
Term Structure Fit	matrix. The Mean Error (Bias), Mean Absolute Error, Standard Deviation of Error and Root Mean Square Error
Term Structure Volatility Fit	matrix. The theoretical and empirical volatility of futures returns for each observed

When debugging = T, 9 objects are returned in addition to those returned when verbose = T:

P_t	array. The covariance matrix at each observation point, with the third dimension indexing across time
F_t	vector. The function of the Kalman filter covariance matrix at each observation point, with the third dimension indexing across time
K_t	matrix. The Kalman Gain at each observation point, with the third dimension indexing across time
d	matrix. d_t (see details)
Z	matrix. Z_t (see details)
G_t	matrix. G_t (see details)
c_t	vector. C_t (see details)
Q_t	matrix. Q_t (see details)
H	matrix. H (see details)

References

Anderson, B. D. O. and J. B. Moore, (1979). *Optimal filtering* Englewood Cliffs: Prentice-Hall.

Fahrmeir, L. and G. tutz,(1994) *Multivariate Statistical Modelling Based on Generalized Linear Models*. Berlin: Springer.

Schwartz, E. S., and J. E. Smith, (2000). Short-Term Variations and Long-Term Dynamics in Commodity Prices. *Manage. Sci.*, 46, 893-911.

Cortazar, G., and L. Naranjo, (2006). An N-factor Gaussian model of oil futures prices. *Journal of Futures Markets: Futures, Options, and Other Derivative Products*, 26(3), 243-268.

Durbin, J., and S. J. Koopman, (2012). *Time series analysis by state space methods*. Oxford university press.

Examples

```
##Example 1 - complete, stitched data.
##Replicating the Schwartz and Smith (2000)
##Two-Factor commodity pricing model applied to crude oil:

SS_stitched_filtered <- NFCP_Kalman_filter(
  parameter_values = SS_oil$two_factor,
  parameter_names = names(SS_oil$two_factor),
  log_futures = log(SS_oil$stitched_futures),
  futures_TTM = SS_oil$stitched_TTM,
  ## maturity groupings need not be considered here:
  ME_TTM = NULL,
  dt = SS_oil$dt,
  verbose = FALSE)

##Example 2 - incomplete, contract data.
##Replicating the Schwartz and Smith (2000)
##Two-Factor commodity pricing model applied to all available
##crude oil contracts:

SS_2F <- SS_oil$two_factor
##omit stitched contract white noise
SS_2F <- SS_2F[!grepl("ME",
  names(SS_2F))]

# Evaluate two different measurement errors
SS_2F[c("ME_1", "ME_2")] <- c(0.01, 0.04)

## Seperate measurement error into two different maturity groupings
SS_ME_TTM <- c(1,3)
## ME_1 is applied for observed contracts with less than one year
## maturity, whilst ME_2 considers contracts with maturity greater
## than one year, and less than three years

#Kalman filter
SS_contract_filtered <- NFCP_Kalman_filter(
  parameter_values = SS_2F,
  parameter_names = names(SS_2F),
  ## All available contracts are considered
  log_futures = log(SS_oil$contracts),
  ## Respective 'futures_TTM' of these contracts are input:
  futures_TTM = SS_oil$contract_maturities,
  ME_TTM = SS_ME_TTM,
```

```
dt = SS_oil$dt,
verbose = FALSE)
```

NFCP_MLE

N-factor model parameter estimation through the Kalman filter and maximum likelihood estimation

Description

the NFCP_MLE function performs parameter estimation of an n-factor model given observable term structure futures data through maximum likelihood estimation. NFCP_MLE allows for missing observations as well as constant or variable time to maturity of observed futures contracts.

Usage

```
NFCP_MLE(
  log_futures,
  dt,
  futures_TTM,
  N_factors,
  N_ME = 1,
  ME_TTM = NULL,
  GBM = TRUE,
  estimate_initial_state = FALSE,
  Richardsons_extrapolation = TRUE,
  cluster = FALSE,
  Domains = NULL,
  ...
)
```

Arguments

log_futures	Object of class <code>matrix</code> corresponding to the natural logarithm of observable futures prices. NA's are allowed within the <code>matrix</code> . Every column of the matrix must correspond to a particular futures contract, with each row corresponding to a quoted price on a given date.
dt	Constant, discrete time step of observations
futures_TTM	Object of class <code>vector</code> or <code>matrix</code> that specifies the time to maturity of observed futures contracts. time to maturity can be either constant (i.e. class <code>vector</code>) or time dependent (i.e. class <code>matrix</code>). When the time to maturity of observed futures contracts is time dependent, the dimensions of <code>futures_TTM</code> must be identical to that of <code>log_futures</code> . Every element of <code>futures_TTM</code> corresponds to the time to maturity, in years, of a futures contract at a given observation date.
N_factors	numeric. Number of state variables in the spot price process.
N_ME	numeric. The number of independent measuring errors of observable futures contracts to consider in the Kalman filter.
ME_TTM	vector of maturity groupings to consider for observed futures prices. The length of <code>ME_TTM</code> must be equal to the number of 'ME' parameter values, and the maximum of <code>ME_TTM</code> must be greater than the maximum observed time-to-maturity of a futures contract. When the number of 'ME' parameter values is equal to

	one or the total number of contracts (i.e., columns of <code>log_futures</code>), this argument is optional and not considered. The measurement error of an observation is highly influenced by its time-to-maturity, see details .
GBM	logical. When TRUE, factor 1 of the model is assumed to follow a Brownian Motion, inducing a unit-root in the spot price process.
estimate_initial_state	logical. When TRUE, the initial state vector is specified as unknown parameters of the commodity pricing model. When FALSE, a "diffuse" assumption is taken instead (see details)
Richardsons_extrapolation	logical. When TRUE, the <code>grad</code> function from the <code>numDeriv</code> package is called to approximate the gradient within the <code>genoud</code> optimization algorithm.
cluster	an optional object of the 'cluster' class returned by one of the <code>makeCluster</code> commands in the <code>parallel</code> package to allow for parameter estimation to be performed across multiple cluster nodes.
Domains	an optional matrix of the lower and upper bounds for the parameter estimation process. The <code>NFCP_domains</code> function is highly recommended. When <code>Domains</code> is not specified, the standard bounds specified within the <code>NFCP_domains</code> function are used.
...	additional arguments to be passed into the <code>genoud</code> function. See <code>help(genoud)</code>

Details

NFCP_MLE is a wrapper function that uses the genetic algorithm optimization function `genoud` from the `rgenoud` package to optimize the log-likelihood score returned from the `NFCP_Kalman_filter` function. When `Richardsons_extrapolation = TRUE`, gradients are approximated numerically within the optimization algorithm through the `grad` function from the `numDeriv` package. NFCP_MLE is designed to perform parameter estimation as efficiently as possible, ensuring a global optimum is reached even with a large number of unknown parameters and state variables. Arguments passed to the `genoud` function can greatly influence estimated parameters and must be considered when performing parameter estimation. Recommended arguments to pass into the `genoud` function are included within the vignette of NFCP. All arguments of the `genoud` function may be passed through the `NFCP_MLE` function (except for `gradient.check`, which is hard set to false).

NFCP_MLE performs boundary constrained optimization of log-likelihood scores and does not allow for out-of-bounds evaluations within the `genoud` optimization process, preventing candidates from straying beyond the bounds provided by `Domains`. When `Domains` is not specified, the default bounds specified by the `NFCP_domains` function are used.

The N-factor model The N-factor model was first presented in the work of Cortazar and Naranjo (2006, equations 1-3). The N-factor framework describes the spot price process of a commodity as the correlated sum of N state variables x_t .

When GBM = TRUE:

$$\log(S_t) = \sum_{i=1}^N x_{i,t}$$

When GBM = FALSE:

$$\log(S_t) = E + \sum_{i=1}^N x_{i,t}$$

Additional factors within the spot-price process are designed to result in additional flexibility, and possibly fit to the observable term structure, in the spot price process of a commodity. The fit of different N-factor models, represented by the log-likelihood can be directly compared with statistical testing possible through a chi-squared test.

Flexibility in the spot price under the N-factor framework allows the first factor to follow a Brownian Motion or Ornstein-Uhlenbeck process to induce a unit root. In general, an N-factor model where $GBM = T$ allows for non-reversible behaviour within the price of a commodity, whilst $GBM = F$ assumes that there is a long-run equilibrium that the commodity price will revert to in the long-term.

State variables are thus assumed to follow the following processes:

When $GBM = TRUE$:

$$dx_{1,t} = \mu^* dt + \sigma_1 dw_1 t$$

When $GBM = FALSE$:

$$dx_{1,t} = -(\lambda_1 + \kappa_1 x_{1,t})dt + \sigma_1 dw_1 t$$

And:

$$dx_{i,t} =_{i \neq 1} -(\lambda_i + \kappa_i x_{i,t})dt + \sigma_i dw_i t$$

where:

$$E(w_i)E(w_j) = \rho_{i,j}$$

The following constant parameters are defined as:

param μ : long-term growth rate of the Brownian Motion process.

param E : Constant equilibrium level.

param $\mu^* = \mu - \lambda_1$: Long-term risk-neutral growth rate

param λ_i : Risk premium of state variable i .

param κ_i : Reversion rate of state variable i .

param σ_i : Instantaneous volatility of state variable i .

param $\rho_{i,j} \in [-1, 1]$: Instantaneous correlation between state variables i and j .

Disturbances - Measurement Error:

The Kalman filtering algorithm assumes a given measure of measurement error or disturbance in the measurement equation (ie. matrix H). Measurement errors can be interpreted as error in the model's fit to observed prices, or as errors in the reporting of prices (Schwartz and Smith, 2000). These disturbances are typically assumed independent.

var ME_i measurement error of contract i .

where the measurement error of futures contracts ME_i is equal to 'ME_' [i] (i.e. 'ME_1', 'ME_2', ...) specified in arguments parameter_values and parameter_names.

There are three particular cases on how the measurement error of observations can be treated in the `NFCP_Kalman_filter` function:

Case 1: Only one ME is specified. The Kalman filter assumes that the measurement error of observations are independent and identical.

Case 2: One ME is specified for every observed futures contract. The Kalman filter assumes that the measurement error of observations are independent and unique.

Case 3: A series of ME's are specified for a given grouping of maturities of futures contracts. The Kalman filter assumes that the measurement error of observations are independent and unique to their respective time-to-maturity.

Grouping of maturities for case 3 is specified through the `ME_TTM` argument. This is a vector that specifies the maximum maturity to consider for each respective ME parameter argument.

in other words, `ME_1` is considered for observations with TTM less than `ME_TTM[1]`, `ME_2` is considered for observations with TTM less than `ME_TTM[2]`, ..., etc.

The first case is clearly the simplest to estimate, but can be a restrictive assumption. The second case is clearly the most difficult to estimate, but can be an infeasible assumption when considering all available futures contracts that make up the term structure of a commodity.

Case 3 thus serves to ease the restriction of case 1, and allow the user to make the modeling of measurement error as simple or complex as desired for a given set of maturities.

Diffuse Kalman Filtering

If `estimate_initial_state = F`, a 'diffuse' assumption is used within the Kalman filtering algorithm. Factors that follow an Ornstein-Uhlenbeck are assumed to equal zero. When `estimate_initial_state = F` and `GBM = T`, the initial value of the first state variable is assumed to equal the first element of `log_futures`. This is an assumption that the initial estimate of the spot price is equal to the closest to maturity observed futures price.

The initial covariance of the state vector for the Kalman Filtering algorithm assumed to be equal to matrix Q

Initial states of factors that follow an Ornstein-Uhlenbeck process are generally not estimated with a high level of precision, due to the transient effect of the initial state vector on future observations, however the initial value of a random walk variable persists across observations (see Schwartz and Smith (2000) for more details).

Value

NFCP_MLE returns a list with 10 objects. 9 objects are returned when the user has specified not to calculate the hessian matrix at solution.

MLE	numeric The Maximum-Likelihood-Estimate of the solution
estimated_parameters	vector. The estimated parameters
standard_errors	vector. Standard error of the estimated parameters. Returned only when <code>hessian = TRUE</code>
x_t	vector. The final observation of the state vector
X	matrix. All observations of the state vector, after the updating equation has been applied
Y	matrix. Estimated futures prices at each observation
V	matrix. Estimation error of each futures contracts at each observation
Filtered Error	matrix. The Mean Error (Bias), Mean Absolute Error, Standard Deviation of Error and
Term Structure Volatility Fit	matrix. The theoretical and empirical volatility of futures returns for each observed
proc_time	list. The real and CPU time (in seconds) the NFCP_MLE function has taken.
genoud_value	list. The output of the called genoud function.

References

- Schwartz, E. S., and J. E. Smith, (2000). Short-Term Variations and Long-Term Dynamics in Commodity Prices. *Manage. Sci.*, 46, 893-911.
- Cortazar, G., and L. Naranjo, (2006). An N-factor Gaussian model of oil futures prices. *Journal of Futures Markets: Futures, Options, and Other Derivative Products*, 26(3), 243-268.
- Mebane, W. R., and J. S. Sekhon, (2011). Genetic Optimization Using Derivatives: The rgenoud Package for R. *Journal of Statistical Software*, 42(11), 1-26. URL <http://www.jstatsoft.org/v42/i11/>.

Examples

```
##Perform One Generation of Maximum Likelihood Estimation on the
##first 20 weekly observations of the Schwartz and Smith (2000) Crude Oil Data:
SS_2F_estimated_model <- NFCP_MLE(
  ####Arguments
  log_futures = log(SS_oil$contracts)[1:20,1:5],
```

```

dt = SS_oil$dt,
futures_TTM= SS_oil$contract_maturities[1:20,1:5],
N_ME = 1,
N_factors = 1, GBM = TRUE,
####Genoud arguments:
hessian = TRUE,
Richardsons_extrapolation = FALSE,
pop.size = 4, optim.method = "L-BFGS-B", print.level = 0,
max.generations = 0, solution.tolerance = 10)

```

NFCP_parameters

Specify parameters of N-factor model

Description

the NFCP_parameters function specifies the parameters of a commodity pricing model under the N-factor framework first described by Cortazar and Naranjo (2006). This function is a recommended starting position for the application of N-factor models within the NFCP package.

Usage

```
NFCP_parameters(N_factors, GBM, initial_states, N_ME, verbose = TRUE)
```

Arguments

N_factors	numeric. Number of state variables in the spot price process.
GBM	logical. If GBM = T, factor 1 of the model is assumed to follow a Brownian Motion, inducing a unit-root in the spot price process.
initial_states	logical. If initial_states = T, the initial state vector is specified as unknown parameters of the commodity pricing model.
N_ME	numeric. The number of independent measuring errors of observable futures contracts to consider in the Kalman filter.
verbose	logical. If verbose = T, the specified N-factor model is printed when the function is called.

Details

The N-factor model The N-factor model was first presented in the work of Cortazar and Naranjo (2006, equations 1-3). The N-factor framework describes the spot price process of a commodity as the correlated sum of N state variables x_t .

When GBM = TRUE:

$$\log(S_t) = \sum_{i=1}^N x_{i,t}$$

When GBM = FALSE:

$$\log(S_t) = E + \sum_{i=1}^N x_{i,t}$$

Additional factors within the spot-price process are designed to result in additional flexibility, and possibly fit to the observable term structure, in the spot price process of a commodity. The fit of different N-factor models, represented by the log-likelihood can be directly compared with statistical testing possible through a chi-squared test.

Flexibility in the spot price under the N-factor framework allows the first factor to follow a Brownian Motion or Ornstein-Uhlenbeck process to induce a unit root. In general, an N-factor model where $GBM = T$ allows for non-reversible behaviour within the price of a commodity, whilst $GBM = F$ assumes that there is a long-run equilibrium that the commodity price will revert to in the long-term.

State variables are thus assumed to follow the following processes:

When $GBM = TRUE$:

$$dx_{1,t} = \mu^* dt + \sigma_1 dw_1 t$$

When $GBM = FALSE$:

$$dx_{1,t} = -(\lambda_1 + \kappa_1 x_{1,t})dt + \sigma_1 dw_1 t$$

And:

$$dx_{i,t} =_{i \neq 1} -(\lambda_i + \kappa_i x_{i,t})dt + \sigma_i dw_i t$$

where:

$$E(w_i)E(w_j) = \rho_{i,j}$$

The following constant parameters are defined as:

var μ : long-term growth rate of the Brownian Motion process.

var E : Constant equilibrium level.

var $\mu^* = \mu - \lambda_1$: Long-term risk-neutral growth rate

var λ_i : Risk premium of state variable i .

var κ_i : Reversion rate of state variable i .

var σ_i : Instantaneous volatility of state variable i .

var $\rho_{i,j} \in [-1, 1]$: Instantaneous correlation between state variables i and j .

Disturbances - Measurement Error:

The Kalman filtering algorithm assumes a given measure of measurement error or disturbance in the measurement equation (ie. matrix H). Measurement errors can be interpreted as error in the model's fit to observed prices, or as errors in the reporting of prices (Schwartz and Smith, 2000). These disturbances are typically assumed independent.

var ME_i measurement error of contract i .

where the measurement error of futures contracts ME_i is equal to 'ME_' [i] (i.e. 'ME_1', 'ME_2', ...) specified in arguments `parameter_values` and `parameter_names`.

There are three particular cases on how the measurement error of observations can be treated in the `NFCP_Kalman_filter` function:

Case 1: Only one ME is specified. The Kalman filter assumes that the measurement error of observations are independent and identical.

Case 2: One ME is specified for every observed futures contract. The Kalman filter assumes that the measurement error of observations are independent and unique.

Case 3: A series of ME's are specified for a given grouping of maturities of futures contracts. The Kalman filter assumes that the measurement error of observations are independent and unique to their respective time-to-maturity.

Grouping of maturities for case 3 is specified through the `ME_TTM` argument. This is a vector that specifies the maximum maturity to consider for each respective ME parameter argument.

in other words, ME_1 is considered for observations with TTM less than ME_TTM[1], ME_2 is considered for observations with TTM less than ME_TTM[2], ..., etc.

The first case is clearly the simplest to estimate, but can be a restrictive assumption. The second case is clearly the most difficult to estimate, but can be an infeasible assumption when considering all available futures contracts that make up the term structure of a commodity.

Case 3 thus serves to ease the restriction of case 1, and allow the user to make the modeling of measurement error as simple or complex as desired for a given set of maturities.

Diffuse Assumption: If `initial_states = F`, a 'diffuse' assumption is made within Kalman filtering and parameter estimation functions (See `NFCP.MLE` or `NFCP.Kalman.filter` for more information)

Value

A vector of parameter names for a specified N-factor spot price process. This vector is ideal for application within many other functions within the NFCP package

References

Schwartz, E. S., and J. E. Smith, (2000). Short-Term Variations and Long-Term Dynamics in Commodity Prices. *Manage. Sci.*, 46, 893-911.

Cortazar, G., and L. Naranjo, (2006). An N-factor Gaussian model of oil futures prices. *Journal of Futures Markets: Futures, Options, and Other Derivative Products*, 26(3), 243-268.

Examples

```
##Generate parameter of a Two-factor model Crude Oil model
##as first presented by Schwartz and Smith (2000):
two_factor_parameters <- NFCP_parameters(N_factors = 2,
                                         GBM = TRUE,
                                         initial_states = FALSE,
                                         N_ME = 5)

print(two_factor_parameters)
```

spot_price_forecast	<i>Forecast the spot prices of an N-factor model</i>
---------------------	--

Description

Analytically forecast expected spot prices following the "true" process of a given n-factor stochastic model

Usage

```
spot_price_forecast(x_0, parameters, t, percentiles = NULL)
```

Arguments

<code>x_0</code>	Initial values of the state vector.
<code>parameters</code>	A named vector of parameter values of a specified N-factor model. Function <code>NFCP_parameters</code> is recommended.
<code>t</code>	a vector of discrete time points to forecast
<code>percentiles</code>	Optional. A vector of percentiles to include probabilistic forecasting intervals.

Details

Future expected spot prices under the N-factor model can be forecasted through the analytic expression of expected future prices under the "true" N-factor process.

Given that the log of the spot price is equal to the sum of the state variables (equation 1), the spot price is log-normally distributed with the expected prices given by:

$$E[S_t] = \exp(E[\ln(S_t)] + \frac{1}{2} \text{Var}[\ln(S_t)])$$

Where:

$$E[\ln(S_t)] = \sum_{i=1}^N e^{-(\kappa_i t)} x_i(0) + \mu t$$

Where $\kappa_i = 0$ when GBM=T and $\mu = 0$ when GBM = F

$$\text{Var}[\ln(S_t)] = \sigma_1^2 t + \sum_{i,j \neq 1} \sigma_i \sigma_j \rho_{i,j} \frac{1 - e^{-(\kappa_i + \kappa_j)t}}{\kappa_i + \kappa_j}$$

and thus:

$$E[S_t] = \exp\left(\sum_{i=1}^N e^{-\kappa_i t} x_i(0) + \left(\mu + \frac{1}{2} \sigma_1^2\right)t + \frac{1}{2} \sum_{i,j \neq 1} \sigma_i \sigma_j \rho_{i,j} \frac{1 - e^{-(\kappa_i + \kappa_j)t}}{\kappa_i + \kappa_j}\right)$$

Under the assumption that the first factor follows a Brownian Motion, in the long-run expected spot prices grow over time at a constant rate of $\mu + \frac{1}{2} \sigma_1^2$ as the $e^{-\kappa_i t}$ and $e^{-(\kappa_i + \kappa_j)t}$ terms approach zero.

An important consideration when forecasting spot prices using parameters estimated through maximum likelihood estimation is that the parameter estimation process takes the assumption of risk-neutrality and thus the true process growth rate μ is not estimated with a high level of precision. This can be shown from the higher standard error for μ than other estimated parameters, such as the risk-neutral growth rate μ^* . See Schwartz and Smith (2000) for more details.

Value

spot_price_forecast returns a vector of expected future spot prices under a given N-factor model at specified discrete future time points. When percentiles are specified, the function returns a matrix with the corresponding confidence bands in each column of the matrix.

References

- Schwartz, E. S., and J. E. Smith, (2000). Short-Term Variations and Long-Term Dynamics in Commodity Prices. *Manage. Sci.*, 46, 893-911.
- Cortazar, G., and L. Naranjo, (2006). An N-factor Gaussian model of oil futures prices. *Journal of Futures Markets: Futures, Options, and Other Derivative Products*, 26(3), 243-268.

Examples

```
# Forecast the Schwartz and Smith (2000) two-factor oil model:

##Step 1 - Kalman filter of the two-factor oil model:
SS_2F_filtered <- NFCP_Kalman_filter(SS_oil$two_factor,
```

```

names(SS_oil$two_factor),
log(SS_oil$stitched_futures),
SS_oil$dt,
SS_oil$stitched_TTM,
verbose = TRUE)

##Step 2 - Probabilistic forecast of N-factor stochastic differential equation (SDE):
spot_price_forecast(x_0 = SS_2F_filtered$x_t,
                    parameters = SS_oil$two_factor,
                    t = seq(0,9,1/12),
                    percentiles = c(0.1, 0.9))

```

spot_price_simulate	<i>Simulate spot prices of an N-factor model through Monte Carlo simulation</i>
---------------------	---

Description

Simulate risk-neutral price paths of an an N-factor commodity pricing model through Monte Carlo Simulation.

Usage

```

spot_price_simulate(
  x_0,
  parameters,
  t = 1,
  dt = 1,
  N_simulations = 2,
  antithetic = TRUE,
  verbose = FALSE
)

```

Arguments

x_0	Initial values of the state vector.
parameters	A named vector of parameter values of a specified N-factor model. Function NFCP_parameters is recommended.
t	the number of years to simulate
dt	discrete time step of simulation
N_simulations	total number of simulations
antithetic	logical. Should antithetic price paths be simulated?
verbose	logical. Should simulated state variables be output? see returns

Details

The `spot_price_simulate` function is able to quickly simulate a large number of risk-neutral price paths of a commodity following the N-factor model. Simulating risk-neutral price paths of a commodity under an N-factor model through Monte Carlo simulations allows for the valuation of commodity related investments and derivatives, such as American Options and Real Options through dynamic programming methods. The `spot_price_simulate` function quickly and efficiently simulates an N-factor model over a specified number of years, simulating antithetic price paths as a simple variance reduction technique. The `spot_price_simulate` function uses the `mvrnorm` function from the MASS package to draw from a multivariate normal distribution for the simulation shocks.

The N-factor model stochastic differential equation is given by:

Brownian Motion processes (ie. factor one when GBM = T) are simulated using the following solution:

$$x_{1,t+1} = x_{1,t} + \mu^* \Delta t + \sigma_1 \Delta t Z_{t+1}$$

Where Δt is the discrete time step, μ^* is the risk-neutral growth rate and σ_1 is the instantaneous volatility. Z_t represents the independent standard normal at time t .

Ornstein-Uhlenbeck Processes are simulated using the following solution:

$$x_{i,t} = x_{i,0} e^{-\kappa_i t} - \frac{\lambda_i}{\kappa_i} (1 - e^{-\kappa_i t}) + \int_0^t \sigma_i e^{\kappa_i s} dW_s$$

Where a numerical solution is obtained by numerically discretising and approximating the integral term using the Euler-Maruyama integration scheme:

$$\int_0^t \sigma_i e^{\kappa_i s} dW_s = \sum_{j=0}^t \sigma_i e^{\kappa_i j} dW_s$$

Value

`spot_price_simulate` returns a list when `verbose = T` and a matrix of simulated price paths when `verbose = F`. The returned objects in the list are:

State_Variables	A matrix of simulated state variables for each factor is returned when <code>verbose = T</code> . The number of factors is equal to the number of state variables.
Prices	A matrix of simulated price paths. Each column represents one simulated price path and each row represents a time step.

References

- Schwartz, E. S., and J. E. Smith, (2000). Short-Term Variations and Long-Term Dynamics in Commodity Prices. *Manage. Sci.*, 46, 893-911.
- Cortazar, G., and L. Naranjo, (2006). An N-factor Gaussian model of oil futures prices. *Journal of Futures Markets: Futures, Options, and Other Derivative Products*, 26(3), 243-268.

Examples

```
# Example 1
###Simulate a geometric Brownian motion (GBM) process:
```

```

## starting price of 20, with a growth of 5% p.a. and
## volatility of 20% p.a.
simulated_spot_prices <- spot_price_simulate(
  x_0 = log(20),
  parameters = c(mu_rn = (0.05 - (1/2) * 0.2^2), sigma_1 = 0.2),
  t = 1,
  dt = 1/12,
  N_simulations = 1e3)

# Example 2
###Simulate future spot price paths under Risk-Neutrality and under the
###Schwartz - Smith two factor model:

##Step 1 - Run the Kalman Filter for the Two-Factor Oil Model:
SS_2F_filtered <- NFCP_Kalman_filter(parameter_values = SS_oil$two_factor,
                                     parameter_names = names(SS_oil$two_factor),
                                     log_futures = log(SS_oil$stitched_futures),
                                     dt = SS_oil$dt,
                                     futures_TTM = SS_oil$stitched_TTM,
                                     verbose = TRUE)

#Step 2 - Simulate spot prices:
##100 antithetic simulations of one year of monthly observations
simulated_spot_prices <- spot_price_simulate(
  x_0 = SS_2F_filtered$x_t,
  parameters = SS_oil$two_factor,
  t = 1,
  dt = 1/12,
  N_simulations = 1e3,
  antithetic = TRUE,
  verbose = TRUE)

```

SS_oil

Crude oil term structure futures data (1990 - 1995)

Description

The SS_oil list object features the approximate weekly observations of Crude Oil (WTI) futures contracts used to develop a two-factor commodity pricing model within the prominent work of Schwartz and Smith (2000) titled: "Short-Term Variations and long-Term Dynamics in Commodity Prices". The two-factor commodity pricing model presented within this study is also included. The SS_oil list object is used extensively within the NFCP package to provide working examples and showcase the features of the package.

Usage

```
data(SS_oil)
```

Format

A list Containing eight objects:

Usage

```
stitch_contracts(
  futures,
  futures_TTM = NULL,
  maturity_matrix = NULL,
  rollover_frequency = NULL,
  contract_numbers = NULL,
  verbose = FALSE
)
```

Arguments

futures	Contract futures price data. Each row of Futures should represent one observation of futures prices and each column should represent one quoted futures contract. NA's in Futures are allowed, representing missing observations.
futures_TTM	A vector of contract maturities to stitch
maturity_matrix	The time-to-maturity (in years) for each contract at each given observation point. The dimensions of maturity_matrix should match those of Futures
rollover_frequency	the frequency (in years) at which contracts should be rolled over
contract_numbers	A vector of contract numbers offset from the closest-to-maturity contract at which to stitch contracts.
verbose	logical. Should additional information be output? see details

Details

This function aggregates a set of futures contract data by stitching contract data over an observation period, resulting in a set of futures observations that is 'complete' (ie. Does not feature missing observations). Aggregated futures data benefit from several computational efficiencies compared to raw contract data, but results in the loss of futures price information.

There are two methods of the `stitch_contracts` function that can be utilized the stitch contracts:

Method 1

`stitch_contracts(futures, contract_numbers, verbose = T)` Futures data may be aggregated by stitching prices according to maturity matching. This method requires the inputs `futures_TTM`, `maturity_matrix` and `rollover_frequency`. This method stitched contracts by matching the observation prices according to which contract has the closest time-to-maturity of the desired maturity specified in `futures_TTM`. Contracts are rolled over at the frequency specified in `rollover_frequency`.

Method 2

`stitch_contracts(futures, futures_TTM, maturity_matrix, rollover_frequency, verbose = T)` Futures data may be stitched according to the contract numbers offset from the closest-to-maturity contract. This method requires only the input `contract_numbers` specifying which contracts should be included. This method is most appropriate when the maturity of available contracts are consistent (ie. contracts expire every month or three months).

Value

`stitch_contracts` returns a matrix of stitched futures prices if `verbose = T` and a list with two or three objects otherwise (see below).

prices	A data frame of Stitched futures prices. Each row represents an observation of the specified contracts.
maturities	A data frame of the time-to-maturity of observed futures prices. Each row represents an observation of the
tickers	A data frame of the named columns of observed futures prices (e.g. contract tickers). Returned only when

References

Schwartz, E. S., and J. E. Smith, (2000). Short-Term Variations and Long-Term Dynamics in Commodity Prices. *Manage. Sci.*, 46, 893-911.

Cortazar, G., and L. Naranjo, (2006). An N-factor Gaussian model of oil futures prices. *Journal of Futures Markets: Futures, Options, and Other Derivative Products*, 26(3), 243-268.

Examples

```
##These examples approximately replicate the Crude Oil data utilized within the
##prominent work of Schwartz and Smith (2000):

###Method 1 - Stitch crude oil contracts according to maturity matching:
SS_stitched_M1 <- stitch_contracts(futures = SS_oil$contracts,
                                   futures_TTM = c(1, 5, 9, 13, 17)/12,
                                   maturity_matrix = SS_oil$contract_maturities,
                                   rollover_frequency = 1/12, verbose = TRUE)

###Method 2 - Stitch crude oil contracts according to nearest contract numbers:
SS_stitched_M2 <- stitch_contracts(futures = SS_oil$contracts,
                                   contract_numbers = c(1, 5, 9, 13, 17), verbose = TRUE)
```

TSfit_volatility	<i>Calculate the volatility term structure of futures returns</i>
------------------	---

Description

Estimate the Theoretical and Empirical Volatility Term Structure of futures returns

Usage

```
TSfit_volatility(parameters, futures, futures_TTM, dt)
```

Arguments

parameters	A named vector of parameters of an N-factor model. Function NFCP.Parameters is recommended.
futures	A Matrix of futures price data. Each column corresponds to a given futures contract, and each row is an observation of the futures contracts.
futures_TTM	A vector listing the Time to Maturities of each listed futures contract from the current observation point.
dt	Numeric. The length of the discrete time step (years).

Details

The fit of the models theoretical volatility term structure of futures returns to those obtained directly from observed futures prices can be used as an additional measure of robustness for the models ability to explain the behavior of a commodities term structure. A commodity pricing model should capture all dynamics of a commodities term structure,

The theoretical model volatility term structure of futures returns is given by the following equation:

$$\sigma_F(\tau) = \sum_{i=1}^N \sum_{j=1}^N \sigma_i \sigma_j \rho_{i,j} e^{-(\kappa_i + \kappa_j)\tau}$$

Under the case that $\kappa_1 = 0$, the model volatility term structure converges to σ_1^2 as τ grows large.

The empirical volatility term structure of futures returns is given by:

$$\hat{\sigma}_F^2(\tau) = \frac{1}{\Delta t} \sum_{i=1}^N (\log(F(t_i, \tau)/F(t_i - \Delta t, \tau)) - \bar{\mu})^2$$

According to Cortazar and Naranjo (2006): "A larger number of factors gives more flexibility to adjust first and second moments simultaneously, hence explaining why (a) four-factor (may) outperform (a) three-factor one in fitting the volatility term structure."

Value

TSfit_volatility returns a matrix with the theoretical and empirical volatility term structure of futures returns, with the number of columns of this matrix coinciding with the number of input futures contracts.

References

Schwartz, E. S., and J. E. Smith, (2000). Short-Term Variations and Long-Term Dynamics in Commodity Prices. *Manage. Sci.*, 46, 893-911.

Cortazar, G., and L. Naranjo, (2006). An N-factor Gaussian model of oil futures prices. *Journal of Futures Markets: Futures, Options, and Other Derivative Products*, 26(3), 243-268.

Examples

```
# Test the volatility term structure fit of the Schwartz-Smith two-factor model on crude oil:
V_TSFit <- TSfit_volatility(
  parameters = SS_oil$two_factor,
  futures = SS_oil$stitched_futures,
  futures_TTM = SS_oil$stitched_TTM,
  dt = SS_oil$dt)
```

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