# Estimating and Testing Exponential-Affine Term Structure Models by Kalman Filter

JIN-CHUAN DUAN

Hong Kong University of Science and Technology, Clear Water Bay, Hong Kong, E-mail: jcduan@ust.hk

JEAN-GUY SIMONATO HEC (Montréal), Canada

**Abstract.** This paper proposes a unified state-space formulation for parameter estimation of exponential-affine term structure models. The proposed method uses an approximate linear Kalman filter which only requires specifying the conditional mean and variance of the system in an approximate sense. The method allows for measurement errors in the observed yields to maturity, and can simultaneously deal with many yields on bonds with different maturities. An empirical analysis of two special cases of this general class of model is carried out: the Gaussian case (Vasicek 1977) and the non-Gaussian case (Cox Ingersoll and Ross 1985 and Chen and Scott 1992). Our test results indicate a strong rejection of these two cases. A Monte Carlo study indicates that the procedure is reliable for moderate sample sizes.

Key words: term structure, Kalman filter, exponential-affine, state-space model, quasi-maximum likelihood

JEL Classification: C22.

### Introduction

The term structure of interest rates describes the relationship between the yield on a default-free debt security and its maturity. Given the high correlation among bond yields of different maturities, many theoretical models attempt to use a small number of factors to explain these joint movements. The typical macro-econometric approach is to specify a time-series model for the short-term interest rate and then employ the expectation hypothesis to derive a structural time series model for bond yields of different maturities. Examples of this approach, such as Hamilton (1987) and Hall, Anderson and Granger (1992), are abundant in the literature.

A different modelling approach, popular in the finance literature and pioneered by Vasicek (1977) and Dothan (1978), starts out by assuming a diffusion process for the instantaneous spot interest rate. Arbitrage arguments are then used to facilitate the derivation of a bond pricing formula. According to these models, the bond price is a function of the unobserved instantaneous spot interest rate and the model's parameters. A more general approach is to assume a set of unobserved state variables and proceed to derive the bond price as a function of these state variables by arbitrage and/or equilibrium arguments. Cox, Ingersoll and Ross (1985) (hereafter CIR), Richard (1978), Longstaff and Schwartz (1992) and Chen and Scott (1992) are some examples. Recently, Duffie and Kan (1996) have provided a characterization for the class of exponential-affine term structure

models, which contains most of the aforementioned term structure models as special cases. Among the existing term structure models, Vasicek (1977) and CIR (1985) have gained prominence in the literature of derivative contract pricing; for example, CIR (1985), Jamshidian (1989), Rabinovitch (1989), Hull and White (1990) and Duan, Moreau and Sealey (1995).

The empirical research focussing on some particular models of the exponential-affine family is extensive. The existent literature can be loosely classified into four categories. The first approach uses proxies for the unobserved factors; for example, Marsh and Rosenfeld (1983), Chan et al. (1992) and Daves and Ehrhardt (1993). The second approach uses a cross-sectional restriction implied by a term structure model to yield a reduced-form model. Examples of this approach are Brown and Dybvig (1986), Titman and Torous (1989), and De Munnik and Schotman (1994). The third category, such as Gibbons and Ramaswamy (1993), involves the derivation of conditional moment restrictions and the application of the generalized method of moments for estimation. The fourth category is the application of the transformed data maximum likelihood method proposed in Duan (1994). Particular cases of this approach are found in Pearson and Sun (1994) and Chen and Scott (1993a). The method proposed in this paper can be regarded as a generalization of the transformed data method. The transformed data method uses a term structure model to define a one-to-one data transformation from the unobserved state variables to the observed bond yields. This method breaks down when the number of yield series is greater than the number of state variables, unless one is willing to impose ad hoc restrictions on the structure of the measurement errors.

The purpose of this study is to develop a unified framework in which all exponential-affine term structure models can be estimated with the simultaneous use of many bond yield series. We utilize a result of Duffie and Kan (1996) which establishes the sufficient and necessary conditions for the obtention of an exponential-affine term structure model. These conditions simply require the drift and variance functions of the underlying diffusion process to be affine in state variables. In this paper, we establish that the conditional mean and variance function of the state variables, over any discrete-time interval, must also be affine in state variables. This result for the discrete-time interval makes it possible to use the Kalman filter and the prediction-error decomposition to obtain an approximate quasi-maximum likelihood solution to the estimation problem for the entire class of exponential affine term structure models.

Special cases of our approach exist in the literature. To our knowledge, Pennacchi (1991) is the first study that uses a state-space formulation for estimating a term structure model of Gaussian nature. Chen and Scott (1993b) and Lund (1994) also propose the use of a Kalman filter-based method to estimate term structure model. When our method is applied to a Gaussian model, if measurement errors are also normally distributed, our method becomes a maximum likelihood estimation. If the model is non-Gaussian as in CIR (1985) case, the method yields an approximate quasi-likelihood function. In that case, the method can be used directly to obtain the parameter estimates or indirectly as a score generator for the efficient method of moment procedure proposed in Gallant and Tauchen (1996).

The proposed method is implemented for two special cases of the exponential affine family. The first case, the Gaussian case, examines the Vasicek (1977) model in which the

unobserved factor follows a Ornstein Uhlenbeck process. The second case examines two non-Gaussian models: the CIR (1985) and Chen and Scott (1992) models where the unobserved factors follow a square-root process. The empirical study is carried out using monthly data for the period from April, 1964 to December, 1997 consisting of yields on the U.S. Treasury debt securities with maturities: 3, 6, 12 and 60 months. For the Vasicek (1977) and CIR (1985) models, the parameter estimates in most cases are statistically significant. Using an approximate robust Lagrange multiplier test, we find that all three models are strongly rejected.

To examine the statistical properties of the proposed method, a Monte Carlo study is carried out. The simulation study suggests that the method is reasonably reliable even for a sample size as small as 150.

## A state-space formulation for the exponential-affine term structure models

Vasicek (1977) and CIR (1985) bond pricing models have been widely used in the term structure literature. Both models involve specifying a stochastic process for the unobserved instantaneous spot interest rate. While Vasicek specifies a continuous time Ornstein-Uhlenbeck process, CIR use a mean-reverting square root process. Multi-factor generalizations of the CIR model have been proposed by Chen and Scott (1992) and Longstaff and Schwartz (1992). Recently Duffie and Kan (1996) have provided a characterization for the class of exponential-affine term structure model which contains, amongst others, the aforementioned models as special cases.

The exponential-affine term structure model is a class of models in which the yields to maturity are affine functions in some abstract state variable vector  $X_t$ , which is assumed to obey the following dynamic:

$$dX_t = U(X_t; \Psi)dt + \Sigma(X_t; \Psi)dW_t, \tag{1}$$

where  $W_t$  is a  $n \times 1$  vector of independent Wiener processes;  $\Psi$  denotes a  $p \times 1$  vector containing the model parameters;  $U(\cdot)$  and  $\Sigma(\cdot)$  are  $n \times 1$  and  $n \times n$  functions regular enough so that equation (1) has a unique solution. The bond pricing formula for this class of model can be generically expressed as:

$$D_t(X_t; \Psi, \tau) = A(\Psi, \tau)e^{-B(\Psi, \tau)X_t}, \tag{2}$$

where  $D_t(X_t; \Psi, \tau)$  is the price at time t of a default-free discount bond with time to maturity  $\tau$ ;  $A(\Psi, \tau)$  is a scalar function and  $B(\Psi, \tau)$  is a  $1 \times n$  vector function. The instantaneous interest rate is, as usual, defined as

$$r_t(X_t; \Psi) = -\lim_{\tau \to 0} \frac{\ln D_t(X_t; \Psi, \tau)}{\tau}.$$
 (3)

Duffie and Kan (1996) have shown that  $D(\cdot)$  is generically exponential-affine, i.e. in the form of equation (2), if and only if  $U(\cdot)$ ,  $\Sigma(\cdot)\Sigma(\cdot)'$  and  $r_t(\cdot)$  are affine in  $X_t$ . Moreover  $A(\cdot)$  and  $B(\cdot)$  in equation (2) can be obtained as the solutions to some ordinary differential equations (see Appendix A).

Let  $R_t(X_t; \Psi, \tau)$  denote the time-t continuously compounded yield on a zero-coupon bond of maturity  $\tau$ . The yield-to-maturity of this bond is given by:

$$R_t(X_t; \Psi, \tau) = -\frac{1}{\tau} \ln(D_t(X_t; \Psi, \tau)). \tag{4}$$

To deal with the estimation problem, it is reasonable to assume that the yields for different maturities are observed with errors of unknown magnitudes. Using the bond pricing formula in (2), the yield to maturity can be written, after the addition of a measurement error, as:

$$R_t(X_t; \Psi, \tau) = -\frac{1}{\tau} \ln(A(\Psi, \tau)) + \frac{1}{\tau} B(\Psi, \tau) X_t + \varepsilon_t, \tag{5}$$

where  $\varepsilon_t$  is assumed to be a normally distributed error term with zero mean and standard deviation  $\sigma_{\varepsilon}$ .

Given that *N* bonds with different maturities are observed, the *N* corresponding yields can be stacked to obtain the following representation:

$$\begin{bmatrix} R_{t}(X_{t}; \Psi, \tau_{1}) \\ R_{t}(X_{t}; \Psi, \tau_{2}) \\ \vdots \\ R_{t}(X_{t}; \Psi, \tau_{N}) \end{bmatrix} = \begin{bmatrix} -\ln(A(\Psi, \tau_{1}))/\tau_{1} \\ -\ln(A(\Psi, \tau_{2}))/\tau_{2} \\ \vdots \\ -\ln(A(\Psi, \tau_{N}))/\tau_{N} \end{bmatrix} + \begin{bmatrix} (1/\tau_{1})B(\Psi, \tau_{1}) \\ (1/\tau_{2})B(\Psi, \tau_{2}) \\ \vdots \\ (1/\tau_{N})B(\Psi, \tau_{N}) \end{bmatrix} X_{t} + \begin{bmatrix} \varepsilon_{t,1} \\ \varepsilon_{t,2} \\ \vdots \\ \varepsilon_{t,N} \end{bmatrix}.$$

$$(6)$$

In terms of the state-space model, this equation is referred to as the measurement equation. To obtain the transition equation for the state-space model, we need to derive expressions for the conditional mean and variance of the unobserved state variables process over a discrete time interval of length h. Define  $m(X_t; \Psi, h) \equiv E(X_{t+h}|X_t)$  and  $\Phi(X_t; \Psi, h) \equiv Var(X_{t+h}|X_t)$ . The transition equation over a discrete time interval, with length equal to h units of time, can be written as:

$$X_{t+1} = m(X_t; \Psi, h) + \Phi(X_t; \Psi, h)^{1/2} \eta_{t+1}$$
(7)

where  $\eta_{t+1}$  is a  $n \times 1$  vector of zero mean and unit variance error terms with  $\Phi(X_t; \Psi, h)^{1/2}$  denoting the Cholesky decomposition of  $\Phi(X_t; \Psi, h)$ . For the special cases where  $\Sigma(\cdot)$  is not a function of  $X_t$ , the transition distribution of  $X_{t+h}$  conditional on  $X_t$  will be Gaussian. For the other cases, this transition distribution will not be Gaussian. Our proposed method does not rely on a complete characterization of the transition distribution. We instead use a partial characterization focusing on the affine form of the first two moments of the

transition distribution. This enables us to obtain an exact likelihood function for the Gaussian model and an approximate one for the non-Gaussian one.

For a Gaussian state-space model, the Kalman filter provides an optimal solution to prediction, updating and evaluating the likelihood function. The Kalman filter recursion is a set of equations which allows an estimator to be updated once a new observation becomes available. The Kalman filter first forms an optimal predictor of the unobserved state variable vector given its previously estimated value. This prediction is obtained using the distribution of the unobserved state variables, conditional on the previous estimated values. These estimates for the unobserved state variables are then updated using the information provided by the observed variables. Prediction errors, obtained as a byproduct of the Kalman filter, can then be used to evaluate the likelihood function.

When the state-space model is non-Gaussian, the Kalman filter can still be applied to obtain approximate first and second moments of the model and the resulting filter is quasi-optimal. The use of this quasi-optimal filter yields an approximate quasi-likelihood function with which parameter estimation can be carried out.

Given that  $U(\cdot)$  and  $\Sigma(\cdot)\Sigma(\cdot)'$  are affine functions of  $X_t, m(X_t; \Psi, h)$  and  $\Phi(X_t; \Psi, h)$  must be affine in  $X_t$  (see Appendix B). Let  $\mathscr{F}_t$  be the information set generated by the observations up to and including time t. That is,  $\mathscr{F}_t$  is a filtration generated by  $\{R_s(X_s; \Psi, \tau_1), \ldots, R_s(X_s; \Psi, \tau_N); \text{ for } s \leq t\}$ . The conditional mean and variance are functions of the lagged unobserved state variables. This implies that  $m(X_t; \Psi, h)$  and  $\Phi(X_t; \Psi, h)$  are not measurable with respect to  $\mathscr{F}_t$ . However, their expected values conditional on  $\mathscr{F}_t$  can easily be expressed. Define  $P_{t+1|t} \equiv \operatorname{Var}(X_{t+1}|\mathscr{F}_t)$  and  $P_t \equiv \operatorname{Var}(X_t|\mathscr{F}_t)$ . Because  $m(\cdot)$  is affine in  $X_t$  we have:

$$m(\hat{X}_t; \Psi, h) = a(\Psi, h) + b(\Psi, h)\hat{X}_t, \tag{8}$$

where  $a(\cdot)$  and  $b(\cdot)$  are  $n \times 1$  and  $n \times n$  matrices and  $\hat{X}_t \equiv E(X_t | \mathscr{F}_t)$ . To implement the quasi-optimal Kalman filter, one must derive a relationship between  $P_{t+1|t}$  and  $P_t$ . By the law of iterated expectations and recognizing that  $\Phi(X_t; \Psi, h)$  is affine in  $X_t$  and  $Cov(X_t, \Phi(X_t; \Psi, h)^{1/2} \eta_{t+1} | \mathscr{F}_t) = 0$  one can derive:

$$P_{t+1|t} = b(\Psi, h)P_t b(\Psi, h)' + \Phi(\hat{X}_t; \Psi, h).$$
(9)

If it were possible to compute  $\hat{X}_t$ , the conditional mean and variance of the system could be correctly specified. Using the measurement and transition equations described in (6) and (7), the standard Kalman filter recursion<sup>1</sup> could then be used to obtain a prediction-error decomposition to evaluate the quasi-likelihood function. The estimates obtained with this quasi-likelihood function would, according to Bollerslev and Wooldridge (1992), be consistent and asymptotically normal. Unfortunately, the linear Kalman filter cannot produce  $\hat{X}_t$ ; rather it yields  $\bar{X}_t$ , that is the linear projection of  $X_t$  on the linear sub-space generated by the observed yields. The conditional mean and variance computed with the estimate  $\bar{X}_t$  should thus be different from the true conditional mean and variance of the system. Nevertheless, this approximation can be expected to work well because it is linearly optimal. This approximation is needed because of the non-Gaussian nature of the

problem, which can be likened to linearizing a non-linear function in the typical Kalman filtering applications.

Since  $\bar{X}_t$  is computable, we use it to approximate  $\hat{X}_t$  which yields a quasi-likelihood function in an approximate sense. We thus define the approximate quasi-likelihood function as the quasi-likelihood function computed with  $\bar{X}_t$  (the estimate of the conditional mean) instead of  $\hat{X}_t$  (the conditional mean). This approximate quasi-likelihood function is used as if it were the correct quasi-likelihood function. Since the statistical properties of this procedure are unknown theoretically, Monte Carlo experiments will be conducted later to assess the quality of this procedure.

Let

$$L_T(\Psi; \mathbf{R}) = \sum_{t=1}^{T} l_t(\Psi; \mathbf{R}_t), \tag{10}$$

where  $l_t(\Psi; \mathbf{R}_t)$  denotes the quasi-likelihood function at time t obtained from the approximate prediction-error decomposition with  $\mathbf{R}_t = \{R_t(X_t; \Psi, \tau_1), \dots, R_t(X_t; \Psi, \tau_N)\}'$  and  $\mathbf{R} = (\mathbf{R}_1, \dots, \mathbf{R}_T)$ . Following Bollerslev and Wooldridge (1992), it is reasonable to expect the parameter vector  $\hat{\Psi}_T$  maximizing  $L_T(\Psi; \mathbf{R})$  to be approximately consistent and asymptotically normal. This approximately consistent parameter should have the following approximate asymptotic distribution:

$$\sqrt{T}(\hat{\Psi}_T - \Psi_0) \sim N(0, \hat{F}_T^{-1} \hat{G}_T \hat{F}_T^{-1}), \tag{11}$$

where

$$\hat{F}_T = \frac{1}{T} \sum_{t=1}^{T} f_t(\hat{\Psi}_T; \mathbf{R}_t), \tag{12}$$

$$\hat{G}_T = \frac{1}{T} \sum_{t=1}^T \frac{\partial \ln l_t(\hat{\Psi}_T; \mathbf{R}_t)'}{\partial \Psi} \frac{\partial \ln l_t(\hat{\Psi}_T; \mathbf{R}_t)}{\partial \Psi},\tag{13}$$

with

$$f_t(\hat{\Psi}; \mathbf{R}_t) = \frac{\partial \mu_t'}{\partial \Psi} \Omega_t^{-1} \frac{\partial \mu_t}{\partial \Psi} + \frac{1}{2} \frac{\partial \Omega_t'}{\partial \Psi} [\Omega_t^{-1} \otimes \Omega_t^{-1}] \frac{\partial \Omega_t}{\partial \Psi},$$

where  $\mu_t$  and  $\Omega_t$  are the conditional mean and variance functions obtained from the linear Kalman filter and  $(\partial \ln l_t(\hat{\Psi}_T; \mathbf{R}_t)/\partial \Psi)$ ,  $(\partial \mu_t/\partial \Psi)$  and  $(\partial \Omega_t/\partial \Psi)$  are of dimension  $1 \times p$ ,  $N \times p$  and  $N^2 \times p$  respectively.

The asymptotic covariance matrix of the quasi-maximum likelihood estimator, commonly referred to as the robust covariance matrix, is in general not equal to the inverse of Fisher's information matrix,  $-\lim_{T\to\infty} \hat{F}_T$ . For Gaussian exponential-affine

term structure models, the approximate quasi-likelihood function becomes the exact likelihood function. This then implies that the asymptotic variance is  $(-\lim_{T\to\infty} \hat{F}_T)^{-1}$  or  $(\lim_{T\to\infty} \hat{G}_T)^{-1}$ , a standard result.

To test the over-identification restrictions imposed by the theoretical model, an approximate robust Lagrange multiplier test can be constructed. Express the unconstrained measurement equation as:

$$\begin{bmatrix} R_{t}(X_{t}; \Psi, \tau_{1}) \\ R_{t}(X_{t}; \Psi, \tau_{2}) \\ \vdots \\ R_{t}(X_{t}; \Psi, \tau_{N}) \end{bmatrix} = \begin{bmatrix} -\ln(A(\Psi, \tau_{1}))/\tau_{1} + \alpha_{1} \\ -\ln(A(\Psi, \tau_{2}))/\tau_{2} + \alpha_{2} \\ \vdots \\ -\ln(A(\Psi, \tau_{N}))/\tau_{N} + \alpha_{N} \end{bmatrix} + \begin{bmatrix} (1/\tau_{1})B(\Psi, \tau_{1}) + \beta_{1} \\ (1/\tau_{2})B(\Psi, \tau_{2}) + \beta_{2} \\ \vdots \\ (1/\tau_{N})B(\Psi, \tau_{N}) + \beta_{N} \end{bmatrix} X_{t} + \begin{bmatrix} \varepsilon_{t,1} \\ \varepsilon_{t,2} \\ \vdots \\ \varepsilon_{t,N} \end{bmatrix}$$

$$(14)$$

where  $\alpha_1$  to  $\alpha_N$  are scalar parameters;  $\beta_1$  to  $\beta_N$  are parameter vectors of dimension  $1 \times n$ . This specification for the unconstrained measurement equation is over-parameterized. First, the exponential-affine term structure model contains risk premium parameters that do not appear in the transition equation. Specifically, risk premium parameters appear in functions  $A(\cdot)$  and/or  $B(\cdot)$  of the measurement equation. These risk premium parameters complicate the matter because the unconstrained model cannot be expressed solely in terms of the parameters  $\alpha_i$ ,  $\beta_i$  and those in the transition equation. Second, since  $X_t$  is an unobserved process, its location and scale parameters are indeterminate. This specification under the alternative hypothesis in equation (14) is thus under-identified.

To perform a Lagrange multiplier test of the null hypothesis,  $H_0$ :  $\alpha_i = 0$ ,  $\beta_i = 0$  for all i, one must deal with the over-parameterization problem by appropriately removing some parameters from the parameter space. More specifically, n parameters must be removed to account for the risk premia, and  $\frac{1}{2}n(n+1) + n$  parameters removed for the alternative model's identification (see Harvey (1990), Section 8.5.1). The question about which parameters can be removed turns out to be model-specific. We thus address this issue when the specific bond pricing models are discussed.

The approximate robust Lagrange multiplier test proceeds as follows. Denote by  $\theta$  the parameter vector under the alternative, that is let  $\theta' = \{\Psi', \varphi'\}$  where  $\varphi$  is a parameter vector of dimension  $N(n+1) - \frac{1}{2}n(n+1) - 2n$ . The parameter vector  $\Psi$  is a subset of  $\{\alpha_i, \beta_i\}$  after removing an appropriate number of parameters. The null hypothesis can be stated as:

$$\dot{H}_0$$
:  $\phi = 0$ .

A result from White (1982) and Domowitz and White (1982) can be used to construct the approximate robust LM test statistic as follows:

$$\varepsilon_{LM} = \hat{S}_{1T} \hat{F}_T^{11} \hat{C}_{11}^{-1} \hat{F}_T^{11} \hat{S}_{1T}' \to \chi_{\{N(n+1) - \frac{1}{2}n(n+1) - 2n\}}^2, \tag{15}$$

where  $\hat{F}_T^{11}$  is the partitioned inverse corresponding to  $\phi$ ;  $\hat{C} = \hat{F}_T^{-1} \hat{G}_T \hat{F}_T^{-1}$ ; and  $\hat{C}_{11}$  is the

submatrix of  $\hat{C}$  corresponding to  $\phi$ ;  $\hat{S}_T = (\partial L_T(\theta; \mathbf{R})/\partial \theta)$  and  $\hat{S}_{1T}$  is the subvector of  $\hat{S}_T$  corresponding to  $\phi$ .

## **Empirical results**

The data set used in this study consists of four monthly yield series for the U.S. Treasury debt securities with maturities: 3, 6, 12 and 60 months taken from the Fama-Bliss data file. All interest rates are expressed on an annualized continuously compounded basis. These data series cover the period from April, 1964 to December, 1997, totalling 405 time series observations each. Table 1 reports summary statistics for these data series. Many empirical studies, for example, Hamilton (1988) and Spindt and Tarhan (1987), have found that the shift in the Federal Reserve monetary policy from October, 1979 to October, 1982, caused a structural break in the interest rate process. We thus report the results for the entire sample period as well as for two subperiods: April, 1964 to October, 1979 and October, 1982 to December, 1997.

The basic unit of time in this study is set equal to one year so that the maturities for all yields are stated in terms of the number of years and the parameter estimates can be interpreted as annualized values. Since the data frequency is monthly, the length of the discrete sampling interval, *h*, equals 1/12. The numerical optimization routine used in this

Table 1. Summary statistics

	April 1964 to I	December 1997	
Maturity	Mean	Std. Dev.	Auto.
3 month	0.0656	0.0265	0.9733
6 months	0.0678	0.0264	0.9754
12 months	0.0702	0.0255	0.9737
60 months	0.0758	0.0234	0.9844
	April 1964 to	October 1979	
Maturity	Mean	Std. Dev.	Auto.
3 month	0.0576	0.0173	0.9581
6 months	0.0601	0.0172	0.9640
12 months	0.0621	0.0165	0.9603
60 months	0.0642	0.0139	0.9724
	October 1982 to	December 1997	
Maturity	Mean	Std. Dev.	Auto.
3 month	0.0623	0.0195	0.9879
6 months	0.0642	0.0198	0.9870
12 months	0.0676	0.0202	0.9839
60 months	0.0784	0.0197	0.9829

study is the quadratic hill-climbing algorithm of Goldfeld, Quandt and Trotter (1966). The convergence criterion, based on the maximum absolute difference in both parameter and functional values between two successive iterations, is set to 0.0001. For each model examined below, the approximate Kalman filter recursion is initialized with the stationary mean and variance of the unobserved state variable(s).

The Gaussian case: Vasicek model

For the Vasicek (1977) model, the unobserved state variable is the instantaneous interest rate which has the following specification:

$$dr_t = \kappa(\theta - r_t)dt + \sigma dz_t,\tag{16}$$

where  $z_t$  is a Wiener process;  $\theta$  is the long-run average of the instantaneous spot interest rate;  $\kappa \ge 0$  is the mean-reverting intensity at which the process returns to its long-run mean; and  $\sigma > 0$  is the volatility parameter of the process.

In terms of this model, the functional forms for  $A(\cdot), B(\cdot), a(\cdot), b(\cdot)$  and  $\Phi(\cdot)$  are given by:

$$A(\Psi, \tau) = e^{\left[\gamma(B(\Psi, \tau) - \tau) - \frac{\sigma^2 B^2(\Psi, \tau)}{4\kappa}\right]},\tag{17}$$

$$B(\Psi, \tau) = \frac{1}{\kappa} [1 - e^{-\kappa \tau}], \tag{18}$$

$$\gamma = \theta + \frac{\sigma\lambda}{\kappa} - \frac{\sigma^2}{2\kappa^2},\tag{19}$$

$$a(\Psi, h) = \theta(1 - e^{-\kappa h}),\tag{20}$$

$$b(\Psi, h) = e^{-\kappa h},\tag{21}$$

$$\Phi(X_t; \Psi, h) = \frac{\sigma^2}{2\kappa} (1 - e^{-2\kappa h}), \tag{22}$$

where  $\lambda$  is the risk premium parameter. In this model,  $\lambda > 0$  implies a positive premium for bond prices.

We assume a diagonal covariance structure for the measurement errors. Its elements are denoted by  $\sigma_{\epsilon_3}^2$ ,  $\sigma_{\epsilon_6}^2$ ,  $\sigma_{\epsilon_6}^2$ ,  $\sigma_{\epsilon_6}^2$ . The second column of Tables 2A, B and C, present the empirical results for the Vasicek model. Parameter estimates are reported along with their robust standard errors in parentheses. In Table 2A, the results for the entire sample period are reported. The measurement error of the six month yield to maturity,  $\sigma_{\epsilon_6}$ , is not significantly different from zero at the usual significance level. All other parameters are significantly different from zero. For the two sub-periods, reported in Tables 2B and 2C, all parameters are statistically significant except for  $\lambda$  in the first sub-period and  $\sigma_{\epsilon_6}$  in the first and second sub-periods.

*Table 2A.* Estimation results for the Vasicek (1977), CIR (1985) and Chen and Scott (1992) models with monthly observations on 3, 6, 12 and 60 months yield series

	April 1964 t	o December 1997							
	Model								
Parameters	Vasicek	CIR	Chen and Scott						
$\overline{\theta_1}$	0.0585	0.0613	0.0303						
-	(0.0228)	(0.0123)	(0.0031)						
$\kappa_1$	0.1189	0.2249	1.1627						
	(0.0159)	(0.0457)	(0.1508)						
$\sigma_1$	0.0214	0.0700	0.1202						
	(0.0021)	(0.0045)	(0.0079)						
$\lambda_1$	0.3418	-0.1110	-0.3139						
•	(0.1190)	(0.0454)	(0.1222)						
$\theta_2$	_	_	4.83e-8						
-	_	_	(0.0000)						
$\kappa_2$	_	_	0.0007						
_	_	_	(0.0614)						
$\sigma_2$	_	_	0.0681						
	_	_	(0.0035)						
$\lambda_2$	_	_	-0.0266						
	_	_	(0.0636)						
$\sigma_{\epsilon_3}$	0.0028	0.0028	0.0027						
-3	(0.0002)	(0.0002)	(0.0001)						
$\sigma_{\epsilon_6}$	7.11e-7	1.88e-9	1.99e-7						
-0	(0.2485)	(7.6255)	(0.0985)						
$\sigma_{\epsilon_{12}}$	0.0030	0.0030	0.0021						
~12	(0.0002)	(0.0002)	(0.0001)						
$\sigma_{\epsilon_{60}}$	0.0099	0.0099	0.0012						
-00	(0.0003)	(0.0003)	(0.0001)						
Chi-square	187.60	201.72	104.20						
P-value	0.0000	0.0000	0.0000						
df	5	5	5						

When compared to the full sample estimates, the estimates for the pre-October, 1979 period as well as those for the post-October, 1982 period indicate a smaller instantaneous volatility,  $\sigma$ . This result supports the finding in the literature that the shift in the Federal Reserve monetary policy from October, 1979 to October, 1982 caused a structural break in the interest rate process.

The Lagrange multiplier test for the restrictions imposed by the Vasicek model on the coefficients of the measurement equation is performed using the specification described earlier. In order to make the factor model identified under the alternative hypothesis,  $\alpha_1$  and  $\beta_1$  are removed from the unconstrained parameter space. For the Vasicek model, only  $A(\cdot)$  are functions of the risk premium parameter  $\lambda$ . The parameter  $\alpha_2$  can also be conveniently regarded as an extra parameter and discarded from the unconstrained parameter space. As a result, the degrees of freedom for our test on four yield series equals

*Table 2B.* Estimation results for the Vasicek (1977), CIR (1985) and Chen and Scott (1992) models with monthly observations on 3, 6, 12 and 60 months yield series

	April 1964	to October 1979					
	Model						
Parameters	Vasicek	CIR	Chen and Scott				
$\theta_1$	0.0675	0.0690	0.0303				
	(0.0189)	(0.0155)	(0.0033)				
$\kappa_1$	0.1956	0.2279	1.3515				
•	(0.0216)	(0.0532)	(0.1899)				
$\sigma_1$	0.0170	0.0663	0.1165				
	(0.0015)	(0.0046)	(0.0086)				
$\lambda_1$	0.1581	-0.0348	-0.3578				
•	(0.2174)	(0.0516)	(0.1500)				
$\theta_2$	<u>-</u>		2.96e-10				
-	_	_	(1.87e-6)				
$\kappa_2$	_	_	1.37e-5				
-	_	_	(0.0874)				
$\sigma_2$	_	_	0.0756				
-	_	_	(0.1274)				
$\lambda_2$	_	_	0.0403				
2	_	_	(0.0513)				
$\sigma_{\epsilon_3}$	0.0028	0.0028	0.0025				
93	(0.0002)	(0.0002)	(0.0002)				
$\sigma_{\epsilon_6}$	7.21e-7	5.81e-7	2.89e-6				
~6	(0.2657)	(0.1510)	(0.0355)				
$\sigma_{\epsilon_{12}}$	0.0026	0.0026	0.0020				
012	(0.0002)	(0.0001)	(0.0001)				
$\sigma_{\epsilon_{60}}$	0.0074	0.0075	0.0010				
~00	(0.0002)	(0.0002)	(0.0001)				
Chi-square	122.95	115.31	148.40				
P-value	0.0000	0.0000	0.0000				
df	5	5	5				

five. For the whole as well as sub-samples, the null hypothesis is strongly rejected as indicated at the bottom portion of Tables 2A, B and C.

The non-Gaussian case: Chen and Scott and CIR models

The Chen and Scott (1992) model allows for several independent state variables with the following dynamic:

$$dx_{i,t} = \kappa_i(\theta_i - x_{i,t})dt + \sigma_i\sqrt{x_{i,t}}dz_{i,t}, \tag{23}$$

for i = 1, ..., k and  $z_{i,t}$  are independent Wiener processes.

Table 2C. Estimation results for the Vasicek (1977), CIR (1985) and Chen and Scott (1992) models with monthly

observations on 3, 6, 12 and 60 months yield series

October 1982 to December 1997							
Parameters	Model						
	Vasicek	CIR	Chen and Scott				
$\overline{\theta_1}$	0.0665	0.0648	0.0289				
	(0.0211)	(0.0125)	(0.0059)				
$\kappa_1$	0.0463	0.1644	1.3101				
-	(0.0152)	(0.0368)	(0.2068)				
$\sigma_1$	0.0113	0.0438	0.0718				
•	(0.0008)	(0.0029)	(0.0079)				
$\lambda_1$	0.7060	-0.1237	-0.5487				
1	(0.0875)	(0.0318)	(0.1441)				
$\theta_2$	<u> </u>	<u> </u>	2.05e-5				
. 2	_	_	(0.0115)				
$\kappa_2$	_	_	1.48e-5				
2	_	_	(0.0083)				
$\sigma_2$	_	_	0.0689				
2	_	_	(0.0425)				
$\lambda_2$	_	_	-0.0331				
2	_	_	(0.0123)				
$\sigma_{\epsilon_3}$	0.0022	0.0022	0.0017				
c <sub>3</sub>	(0.0002)	(0.0002)	(0.0001)				
$\sigma_{\epsilon_6}$	2.94e-7	4.85e-7	8.54e-8				
- 6	(0.2761)	(0.2080)	(0.1096)				
$\sigma_{\epsilon_{12}}$	0.0024	0.0024	0.0017				
812	(0.0002)	(0.0002)	(0.0001)				
$\sigma_{\epsilon_{60}}$	0.0077	0.0077	0.0013				
₹60	(0.0003)	(0.0003)	(0.0001)				
Chi-square	94.876	95.559	65.64				
P-value	0.0000	0.0000	0.0000				
df	5	5	5				

In this section, the one and two-factor version of their model are analyzed. The one-factor version of their model correspond to CIR (1985) model, where the unobserved factor is interpreted as the instantaneous interest rate.

For the two-factor model, the functional forms for  $A(\cdot)$  and  $B(\cdot)$  are given by:

$$A(\Psi, \tau) = A_1(\Psi, \tau) A_2(\Psi, \tau), \tag{24}$$

$$B(\Psi, \tau) = [B_1(\Psi, \tau), B_2(\Psi, \tau)], \tag{25}$$

where  $A_i(\Psi, \tau)$  and  $B_i(\Psi, \tau)$  for i = 1 and 2 have the following specification:

$$A_{i}(\Psi, \tau) = \left[ \frac{2\gamma_{i}e^{[(\kappa_{i} + \lambda_{i} + \gamma_{i})\tau]/2}}{(\kappa_{i} + \lambda_{i} + \gamma_{i})(e^{\gamma_{i}\tau} - 1) + 2\gamma_{i}} \right]^{2\kappa_{i}\theta_{i}/\sigma_{i}^{2}}, \tag{26}$$

$$B_i(\Psi, \tau) = \frac{2(e^{\gamma_i \tau} - 1)}{(\kappa_i + \lambda_i + \gamma_i)(e^{\gamma_i \tau} - 1) + 2\gamma_i},\tag{27}$$

with

$$\gamma_i = \sqrt{\left(\kappa_i + \lambda_i\right)^2 + 2\sigma_i^2}.$$
(28)

The functional forms for  $a(\cdot), b(\cdot)$  and  $\Phi(\cdot)$  are given by:

$$a(\Psi, h) = \begin{bmatrix} \theta_1 (1 - e^{-\kappa_1 h}) \\ \theta_2 (1 - e^{-\kappa_2 h}) \end{bmatrix}, \tag{29}$$

$$b(\Psi, h) = \begin{bmatrix} e^{-\kappa_1 h} & 0\\ 0 & e^{-\kappa_2 h} \end{bmatrix}, \tag{30}$$

$$\Phi(X_t; \Psi, h) = \begin{bmatrix} \Phi_1(x_{1,t}; \Psi, h) & 0\\ 0 & \Phi_2(x_{2,t}; \Psi, h) \end{bmatrix},$$
(31)

with

$$\Phi_{i}(x_{t}; \Psi, h) = x_{t} \frac{\sigma_{i}^{2}}{\kappa_{i}} (e^{-\kappa_{i}h} - e^{-2\kappa_{i}h}) + \theta_{i} \frac{\sigma_{i}^{2}}{2\kappa_{i}} (1 - e^{-\kappa_{i}h})^{2}$$
(32)

where  $\lambda_1$  and  $\lambda_2$  are the risk premia parameters. In this model,  $\lambda_i < 0$  implies a positive premium in bond prices for factor *i*. The CIR (1985) model is a special case of this model, and is obtained by using the top left elements of the matrices defined above.

The estimation results for the CIR (1985) model are reported in the third column of Tables 2A, B and C. These tables contain parameter estimates along with their corresponding robust standard errors in parentheses. Except for  $\lambda$  in the first sub-sample and the six month measurement error, all estimates are significantly different from zero at the 5% level in all samples. In all cases, the unobserved instantaneous interest rate is found to exhibit mean-reverting behavior, i.e.,  $\kappa > 0$ . The magnitudes of the measurement errors are generally small but significantly different from zero.

Comparing the estimates for the entire sample with the ones for the pre-October, 1979 and post-October, 1982 periods reveals that a smaller volatility scale parameter,  $\sigma$ , prevails for these two subperiods. These results are consistent with the previous findings that the shift in the Federal Reserve monetary policy from October, 1979 to October, 1982 caused a structural break in the interest rate process. The mean-reverting property of the interest rate process again prevails for these two subperiods.

The Lagrange multiplier test of the restrictions imposed by the CIR model on the

coefficients of the measurement equation is performed using the specification described earlier. For the CIR model, both  $A(\cdot)$  and  $B(\cdot)$  are functions of the risk premium parameter  $\lambda$ . As with the Vasicek model, we remove  $\alpha_1$ ,  $\alpha_2$  and  $\beta_1$  from the unconstrained parameter space to make the model identified under the alternative specification. As a result, the degrees of freedom for this test equals five. The null hypothesis is strongly rejected in the three samples as shown by the *P*-values from the Chi-square test.

The estimation results for the Chen and Scott (1992) model are reported in the fourth column of Tables 2A, B and C. These tables contain parameter estimates along with their corresponding robust standard errors (reported in parentheses). For the entire sample period, all parameters are significantly different from zero except for  $\kappa_2$  and  $\lambda_2$  and  $\sigma_{\epsilon_6}$ . For the first sub-sample,  $\theta_2, \kappa_2, \sigma_2, \lambda_2$  and  $\sigma_{\epsilon_6}$  are not significantly different from zero whereas  $\theta_2$ ,  $\kappa_2$   $\sigma_2$ , and  $\sigma_{\epsilon_6}$  are not significantly different from zero in the second sub sample.

The most striking feature of these estimates are the small values for  $\theta_2$  and  $\kappa_2$ . With these values, we  $\sigma_2^2 > 2\kappa_2\theta_2$ , which implies that the second factor can reach zero with a positive probability (see Cox, et al. 1985). A small value for  $\kappa_2$  also suggest that the second factor is close to being non-stationary. The resulting test statistics might therefore differ from the approximate asymptotic normal distribution and should be interpreted with care.

The comparison of the estimates from the entire sample with the ones from the pre-October, 1979 and post-October, 1982 periods still points to a structural instability even when the two-factor model is used. This result is once again consistent with the previous findings that the shift in the Federal Reserve monetary policy from October, 1979 to October, 1982 caused a structural break in the interest rate process.

Chen and Scott (1993b) report the estimation results for the Chen and Scott (1992) two-factor model. Their estimates are consistent with our findings. More specifically, the estimates for the long-run mean and mean reversion parameters of the second factor are small, suggesting a nearly integrated second factor.

In the Chen and Scott (1992) model, both  $A(\cdot)$  and  $B(\cdot)$  are functions of the risk premium parameters. To perform the robust Lagrange multiplier test described at the end of Section 2,  $\beta_N$ , which is a  $1 \times n$  vector, can be conveniently removed from the parameter space to account for the risk premium parameters. To ensure that the factor model under the alternative specification is identified,  $\alpha_1$ ,  $\alpha_2$ ,  $\beta_1$  and the second entry of  $\beta_2$  are removed from the unconstrained parameter space. The resulting degrees of freedom for this Chi-square test also equals five. In all samples the null hypothesis is strongly rejected as indicated by the P-values from the Chi-square test.

## A Monte Carlo analysis

To assess the statistical properties of the proposed method, we conduct Monte Carlo experiments for the sample size of 150 and 400 time-series observations with the simultaneous use of 3-, 6-, 12- and 60-month maturities. These sample sizes are chosen because they roughly correspond to the sub and whole sample size of the data sets analyzed in the preceding section.

To perform Monte Carlo experiments, the unobserved state variables of a given model

must first be simulated. For the Vasicek (1977) model the state variable follows an Ornstein-Uhlenbeck process which can be obtained using the exact conditional distribution. More formally, denote  $r_t^*$  to be the simulated value of the process at time t. Using the transition density of this process,  $r_t^*$  is obtained by:

$$r_t^* - r_{t-h}^* = \theta + (r_{t-h}^* - \theta)e^{-\kappa h} + \sqrt{\Phi}\varepsilon_t,$$

where  $\varepsilon_t$  is N(0, 1) with  $r_0^* = \theta$  and  $\phi = (\sigma^2/2\kappa)(1 - e^{-2\kappa h})$  and h set to 1/12.

For the square-root process in CIR (1985), the exact conditional distribution is also used to obtain the simulated values. Let  $x_t$  denote the value of an unobserved state variable following a square-root process. The conditional distribution function for  $x_t$  is the noncentral Chi-square,  $\chi^2[2cx_t; 2d+2, 2w]$ , with 2d+2 degrees of freedom and the parameter of non-centrality, 2w, where

$$c = \frac{2\kappa}{\sigma^2(1 - e^{-\kappa h})},$$
$$d = \frac{2\kappa\theta}{\sigma^2} - 1,$$
$$w = cx_{t-h}e^{-\kappa h}.$$

As discussed in Johnson and Kotz (1970), the non-central Chi-square distribution can be expressed as a mixture of central Chi-squares with degrees of freedom proportional to random variates from a Poisson distribution. This property is used to simulate the process. More precisely,  $x_t^*$ , the simulated value of the process at time t, is obtained by the following steps:

1. Simulate the degrees of freedom of the central Chi-square using

$$df = 2d + 2 + 2i$$
,

where j is a Poisson random variate<sup>2</sup> with mean  $w = cx_{t-h}^*e^{-\kappa h}$ . The time interval h is set to 1/12 and  $x_0^* = \theta$ .

- 2. Let *g* denote the random variate drawn from the central Chi-square with *df* degrees of freedom. This random variate is obtained from a gamma (*df*/2, 2) random variate<sup>3</sup> where *df*/2 is the shape parameter and 2 is the scale parameter.
- 3. Compute  $x_t^* = \frac{g}{2c}$ .

Using the simulated time series for either the Ornstein-Uhlenbeck or the square-root process, equation (5) is then used to compute the simulated yields with independent normally distributed measurement errors. The Monte Carlo simulation results for discretely sampled time-series are reported for the Gaussian case Vasicek (1977) and the non-Gaussian case CIR (1985). The results are obtained using 500 repetitions.

Rows one, two, three and four of Tables 3 and 4 report the true parameter values, the

*Table 3A.* Monte-Carlo experiment results for the maximum likelihood parameter estimator of the Vasicek model for 3, 6, 12 and 60-month yields (500 replications)

T = 150 observations									
	θ	κ	σ	λ	$\sigma_{\epsilon_3}$	$\sigma_{\epsilon_6}$	$\sigma_{\epsilon_1}$	$\sigma_{\epsilon_{60}}$	
True Value	0.0600	0.1000	0.0200	0.3000	1.0000e-03	1.0000e-03	1.0000e-03	1.0000e-03	
Median	0.0620	0.0998	0.0199	0.2877	1.0020e-03	9.9653e-04	9.9418e-04	9.8208e-04	
Mean	0.0617	0.1000	0.0199	0.2938	9.9979e-04	9.9927e-04	9.9575e-04	9.8599e-04	
Std. dev.	0.0214	0.0032	0.0012	0.1094	7.5826e-05	7.9260e-05	8.0301e-05	7.2084e-05	
Cov. rate									
25% Cov. rate	0.2754	0.1896	0.2475	0.2595	0.2236	0.2176	0.2515	0.2156	
50% Cov. rate	0.4810	0.3912	0.4591	0.4930	0.4631	0.4671	0.4890	0.4351	
75% Cov. rate	0.8802	0.6547	0.7006	0.8583	0.7445	0.7146	0.7325	0.7046	
95% Cov. rate	1.0000	0.8982	0.9242	1.0000	0.9621	0.9441	0.9341	0.9182	
95% cov. rate f	or Lagran	ge multipl	lier test: 0	.9321					
			7	T = 400  o	bservations				
	θ	κ	σ	λ	$\sigma_{\epsilon_3}$	$\sigma_{\epsilon_6}$	$\sigma_{\epsilon_{12}}$	$\sigma_{\epsilon_{60}}$	
True Value	0.0600	0.1000	0.0200	0.3000	1.0000e-03	1.0000e-03	1.0000e-03	1.0000e-03	
Median	0.0579	0.1001	0.0199	0.3134	1.0008e-03	1.0011e-03	9.9834e-04	9.9306e-04	
Mean	0.0587	0.1000	0.0199	0.3082	9.9902e-04	9.9984e-04	9.9942e-04	9.9431e-04	
Std. dev.	0.0227	0.0013	0.0007	0.1151	4.9212e-05	5.2000e-05	4.9237e-05	4.2409e-05	
Cov. rate									
25% Cov. rate	0.2216	0.2096	0.2834	0.2156	0.2754	0.2675	0.2435	0.2255	
50% Cov. rate	0.4072	0.4192	0.4890	0.4092	0.4611	0.4591	0.4611	0.4411	
75% Cov. rate	0.7345	0.6926	0.7345	0.7405	0.7345	0.6846	0.7265	0.7445	
95% Cov. rate	0.9960	0.9381	0.9481	0.9960	0.9601	0.9261	0.9421	0.9521	
95% cov. rate f	or Lagran	ge multipl	lier test: 0	.8683					

medians, the means and the standard deviations of the parameter estimates. The remaining four rows report the probability that the true parameter lies in the  $\alpha\%$  confidence interval. This probability is referred to as the  $\alpha\%$  coverage rate. For example, to obtain 95% coverage rate,  $\text{Prob}\{|\Psi_i-\hat{\Psi}_{iT}|<1.96\text{ s.e.}(\hat{\Psi}_{iT})\}$  is to be computed, where  $\Psi_i$  denotes the i-th parameter of the model and s.e.  $(\hat{\Psi}_{iT})$  represents the estimated standard error for the i-th parameter estimator. The bottom row of each panel reports the 95% coverage rate for the over-identification test statistic. This value indicates whether the test is biased or not.

The Monte Carlo simulation results for the Gaussian case (Vasicek (1977) model) are presented in Table 3A. The true parameter values are taken from the estimates, after rounding, reported in the empirical section. The estimated values are exceedingly close to their corresponding true values for all parameters except  $\theta$ . However, the magnitude of the bias for this parameter is small as a percentage of their corresponding true value. The coverage rates of the parameters indicate some departure from the asymptotic distribution, especially when a smaller sample size is used. This departure may be attributed to a small value for  $\kappa$ . As  $\kappa$  approaches zero, the unobserved factor and the yield series become non-stationary.

To investigate further, additional Monte Carlo experiments with 400 observations and different parameter values are reported in the third panel of Table 3B. The top panel of this

*Table 3B.* Monte-Carlo experiment results for the maximum likelihood parameter estimator of the Vasicek model for 3, 6, 12 and 60-month yields (500 replications)

T = 400 observations										
	θ	к	σ	λ	$\sigma_{\epsilon_3}$	$\sigma_{\epsilon_6}$	$\sigma_{\epsilon_1}$	$\sigma_{\varepsilon_m}$		
True Value	0.1000	0.5000	0.0500	1.0000	1.0000e-03	1.0000e-03	1.0000e-03	1.0000e-03		
Median	0.0984	0.5001	0.0499	1.0190	1.0027e-03	9.9840e-04	9.9787e-04	9.9347e-04		
Mean	0.0990	0.5000	0.0498	1.0144	9.9920e-04	9.9904e-04	9.9861e-04	9.9583e-04		
Std. dev.	0.0162	0.0021	0.0016	0.1678	5.6994e-05	5.7883e-05	4.8444e-05	3.6868e-05		
Cov. rate										
25% Cov. rate	0.2216	0.2275	0.2615	0.2216	0.2395	0.2275	0.2495	0.2255		
50% Cov. rate	0.3972	0.4491	0.4850	0.3952	0.4810	0.4631	0.4631	0.4571		
75% Cov. rate	0.7066	0.7146	0.7465	0.6906	0.7405	0.6986	0.7385	0.7365		
95% Cov. rate	0.9421	0.9281	0.9521	0.9561	0.9521	0.9281	0.9541	0.9481		
95% cov. rate f	or Lagran	ge multipl	ier test: 0	.9082						
			1	= 400 o	bservations					
	θ	κ	σ	λ	$\sigma_{\epsilon_2}$	$\sigma_{\epsilon_6}$	$\sigma_{\epsilon_{12}}$	$\sigma_{\epsilon_{60}}$		
True Value	0.1000	0.0100	0.0500	1.0000	1.0000e-03	1.0000e-03	1.0000e-03	1.0000e-03		
Median	0.0998	0.0100	0.0500	1.0018	1.0005e-03	1.0012e-03	9.9712e-04	9.9178e-04		
Mean	0.0954	0.0100	0.0499	1.0022	9.9905e-04	1.0001e-03	9.9847e-04	9.9360e-04		
Std. dev.	0.0961	0.0003	0.0011	0.0263	4.7698e-05	5.0344e-05	4.8951e-05	4.6461e-05		
Cov. rate										
25% Cov. rate	0.5200	0.1980	0.2300	0.2840	0.2580	0.2540	0.2340	0.2380		
			0.4200	0.5540	0.4720	0.4500	0.4640	0.4420		
50% Cov. rate	0.7140	0.4480	0.4380	0.5540	0.4720	0.4580	0.4640	0.4420		
50% Cov. rate 75% Cov. rate	0.7140 1.0000	0.4480 0.6920	0.4380 0.7580	0.5540 0.8980	0.4720	0.4380	0.4640	0.4420		

table reports the results for the case of a larger value for  $\kappa$ , the mean-reverting parameter. When  $\kappa=0.5$ , the biases nearly vanish, and the asymptotic distribution becomes a better approximation to the small sample distribution. The bottom panel of this table reports the results for the case of a smaller mean-reverting parameter value. When its value is set to 0.01, the coverage rates become similar to those reported in the second panel of Table 3A. This result suggests that the magnitude of the mean reversion parameter plays an important role in determining the quality of the estimator when using a finite sample size. This result is not surprising. The mean-reverting parameter  $\kappa$  is a transformation of the first order autoregressive parameter. It is well known in the time-series literature that the standard asymptotic properties of the estimators do not hold in the presence of a unit root.

The 95% coverage rate of the Lagrange multiplier test indicates a slight departure from the asymptotic distribution for all Monte Carlo experiments in Table 3. The test appears more conservative, meaning that the null hypothesis is rejected slightly more than 5% of the time.

The Monte Carlo simulation results for the non-Gaussian CIR (1985) model are presented in Table 4A. The true parameter values are taken from the estimates, after rounding, reported in the preceding empirical section. The estimates are close to their

*Table 4A.* Monte-Carlo experiment results for the approximate quasi-maximum likelihood parameter estimator of the CIR model for 3, 6, 12 and 60-month yields (500 replications)

of the CIK mod	er 101 <i>5</i> , (	), 12 and	00-111011111	yields (500	replications)				
T = 150 observations									
	θ	κ	σ	λ	$\sigma_{\epsilon_3}$	$\sigma_{\epsilon_6}$	$\sigma_{\epsilon_{12}}$	$\sigma_{\epsilon_{60}}$	
True Value	0.0600	0.2000	0.0700	-0.1000	1.0000e-03	1.0000e-03	1.0000e-03	1.0000e-03	
Median	0.0542	0.2216	0.0700	-0.1210	9.9898e-04	9.9541e-04	1.0017e-03	9.9748e-04	
Mean	0.0559	0.2230	0.0698	-0.1226	1.0002e-03	1.0003e-03	9.9575e-04	9.8918e-04	
Std. dev.	0.0111	0.0412	0.0041	0.0413	8.1982e-05	8.5580e-05	7.9028e-05	7.0259e-05	
Cov. rate									
25% Cov. rate	0.2140	0.2080	0.2440	0.2260	0.2660	0.2360	0.2440	0.2380	
50% Cov. rate	0.3940	0.4020	0.4220	0.3920	0.4560	0.4360	0.4840	0.4540	
75% Cov. rate	0.6500	0.7280	0.7160	0.7260	0.7220	0.7020	0.7400	0.7160	
95% Cov. rate	0.9800	0.9940	0.9320	0.9920	0.9420	0.9260	0.9380	0.9420	
95% cov. rate f	or Lagran	ige multip	olier test:	0.9140					
				T = 400  ob	servations				
	θ	κ	σ	λ	$\sigma_{\epsilon_2}$	$\sigma_{\epsilon_{\epsilon}}$	$\sigma_{\epsilon_{12}}$	$\sigma_{\epsilon_{60}}$	
True Value	0.0600	0.2000	0.0700	-0.1000	1.0000e-03	1.0000e-03	1.0000e-03	1.0000e-03	
Median	0.0579	0.2076	0.0699	-0.1076	9.9871e-04	9.9569e-04	9.9697e-04	9.9878e-04	
Mean	0.0583	0.2121	0.0699	-0.1120	9.9724e-04	9.9776e-04	1.0003e-03	9.9852e-04	
Std. dev.	0.0102	0.0372	0.0025	0.0370	5.0516e-05	5.0244e-05	4.8581e-05	4.1450e-05	
Cov. rate									
25% Cov. rate	0.2455	0.2375	0.2236	0.2415	0.2655	0.2415	0.2315	0.2615	
50% Cov. rate	0.4711	0.4671	0.4391	0.4691	0.4531	0.4431	0.4491	0.4711	
75% Cov. rate	0.7026	0.7405	0.7086	0.7285	0.7186	0.7146	0.7166	0.7265	
95% Cov. rate	0.9142	0.9860	0.9521	0.9880	0.9441	0.9321	0.9401	0.9561	
95% cov. rate f	or Lagran	ige multip	olier test:	0.9202					

corresponding true values for all parameters. Parameters estimates for  $\theta$  and  $\kappa$  are slightly biased. However, the magnitude of the bias is small as a percentage of the true parameter value. Overall, the coverage rates of the parameters indicate that the asymptotic distribution for the quasi-maximum likelihood estimator is a good approximation for the small sample distribution. In order to assess the impact of a nearly integrated factor, additional Monte Carlo experiments with 400 observations and different parameter values are reported in Table 4B. The top portion of this table reports the results for a larger value of  $\kappa$ . As in the Gaussian case with  $\kappa=0.5$ , the biases nearly vanish, and the approximate asymptotic distribution becomes a better approximation to the small sample distribution. The bottom panel of this table reports the results for the same parameter values with the exception of  $\kappa$ , which is now set to 0.01, a value closer to zero. With these parameter values, the estimates for  $\theta$  and  $\kappa$  are biased and their coverage rates are poorly approximated by the asymptotic distribution.

Again, as with the Gaussian case, the 95% coverage rate of the Lagrange multiplier test also indicates some departure from the asymptotic distribution in all Monte Carlo experiments in Table 4A. The test is typically more conservative, meaning that the null hypothesis is rejected more than 5% of the time.

*Table 4B.* Monte-Carlo experiment results for the approximate quasi-maximum likelihood parameter estimator of the CIR model for 3, 6, 12 and 60-month yields (500 replications)

				•	1 /			
				T = 400  ob	servations			
	θ	κ	σ	λ	$\sigma_{\epsilon_3}$	$\sigma_{\epsilon_6}$	$\sigma_{\epsilon_{12}}$	$\sigma_{\epsilon_{60}}$
True Value	0.1000	0.5000	0.0500	-1.0000	1.0000e-03	1.0000e-03	1.0000e-03	1.0000e-03
Median	0.1005	0.4975	0.0498	-0.9975	9.9887e-04	1.0003e-03	1.0007e-03	9.9103e-04
Mean	0.1007	0.4978	0.0498	-0.9977	9.9879e-04	9.9896e-04	9.9723e-04	9.6766e-04
Std. dev.	0.0050	0.0248	0.0017	0.0248	3.8536e-05	3.8484e-05	3.9517e-05	2.4056e-04
Cov. rate								
25% Cov. rate	0.2500	0.2380	0.2260	0.2460	0.2160	0.2520	0.2540	0.2600
50% Cov. rate	0.4540	0.4560	0.4460	0.4640	0.4400	0.4800	0.4840	0.4720
75% Cov. rate	0.7340	0.7220	0.7020	0.7200	0.7300	0.7400	0.7600	0.7860
95% Cov. rate	0.9560	0.9520	0.9400	0.9540	0.9420	0.9560	0.9440	0.9680
95% cov. rate f	or Lagrar	ige multip	olier test:	0.9400				
				T = 400  ob	servations			
	θ	κ	σ	λ	$\sigma_{\epsilon_3}$	$\sigma_{\epsilon_6}$	$\sigma_{\epsilon_{12}}$	$\sigma_{\epsilon_{60}}$
True Value	0.1000	0.0100	0.0500	-1.0000	1.0000e-03	1.0000e-03	1.0000e-03	1.0000e-03
Median	0.0591	0.0168	0.0500	-1.0063	9.9458e-04	9.9815e-04	9.9381e-04	9.3662e-04
Mean	0.0648	0.0161	0.0500	-1.0061	9.9469e-04	9.9682e-04	9.9359e-04	1.3205e-03
Std. dev.	0.0160	0.0035	0.0006	0.0035	3.4563e-05	3.7035e-05	3.6774e-05	1.3311e-03
Cov. rate								
25% Cov. rate	0.1140	0.1100	0.2340	0.1060	0.2660	0.2160	0.2700	0.5880
50% Cov. rate	0.1520	0.1760	0.4860	0.1660	0.4680	0.4300	0.4800	0.6820
75% Cov. rate	0.2040	0.2580	0.7480	0.2620	0.7300	0.7240	0.7660	0.7940
95% Cov. rate	0.3120	0.6760	0.9520	0.7540	0.9480	0.9460	0.9580	0.9140
95% cov. rate f	or Lagrar	ige multip	olier test:	0.8940				

Overall, these results suggest that the approximate quasi-maximum likelihood estimator in the case of the CIR model and the maximum likelihood estimator in the case of the Vasicek case tend to behave in a similar fashion. Both estimators are sensitive to a nearly integrated factor, and the Lagrange multiplier test tends to slightly over-reject the null hypothesis. The finite sample properties of the approximate quasi-maximum likelihood estimator are, by and large, reasonably described by their asymptotic counterparts.

## Conclusion

In this article, a unified state-space formulation is developed for estimating term structure models of the exponential-affine family. The method allows for measurement errors in the observed yields to maturity, and is therefore useful for simultaneous estimation based on yields on many bonds with different maturities. The Monte Carlo study indicates that the proposed method is an adequate procedure. The finite sample properties of the approximate quasi-maximum likelihood estimator are reasonably approximated by the asymptotic distribution presented in this paper.

Since our proposed filtering method provides a good approximation result, it is a good candidate for the score generator in the efficient method of moment (EMM) estimation method proposed in Gallant and Tauchen (1996). The EMM method uses the score of a density that has an analytic expression to define a GMM criterion. The method does not require the score generator to nest the structural model, but does require that the auxiliary model on which the score generator is defined closely approximates the target model. Our filtering method can thus be used as the score generator if a further refinement of the parameter estimates is preferred.

Two special cases of the exponential-affine family are used to examine the estimation method. The empirical results are, in some instances, supportive of the properties typical of the exponential-affine term structure models; for example, the mean-reverting property of the interest rate process. This, however, does not suggest that the three exponential-affine term structure models analyzed in this paper are good descriptions of the bond yield behavior. In fact the results from using four different maturities cast doubts as to whether these models can be applied to yields covering a large maturity spectrum. Using a robust Lagrange multiplier test, the three exponential affine models are strongly rejected. These results suggest that future research is needed in order to find a better specification within or beyond the exponential-affine family.

# Appendix A

Differential equations for  $A(\cdot)$  and  $B(\cdot)$  functions

As shown in Duffie and Kan (1993), the solutions for the  $A(\cdot)$  and  $B(\cdot)$  functions can be found by solving the following differential equations:

$$\begin{split} &-\frac{\partial B(\Psi,\tau)}{\partial t}+\mathscr{B}(B(\Psi,\tau))=0, \quad B(\Psi,0)=0,\\ &-\frac{\partial A(\Psi,\tau)}{\partial t}+\mathscr{A}(B(\Psi,\tau))=0, \quad A(\Psi,0)=0, \end{split}$$

with  $-\partial B_i(\Psi, \tau)/\partial t + \mathcal{B}_i(B(\Psi, \tau))$  denoting the coefficients of  $X_{i,t}$  and  $-\partial A(\Psi, \tau)/\partial t + \mathcal{B}(B(\Psi, \tau))$  denoting the term not involving  $X_t$  in the following equation:

$$\begin{split} r_t(X_t; \Psi) &= -\frac{\partial A(\Psi, \tau)}{\partial t} - \frac{\partial B(\Psi, \tau)}{\partial t} X_t + B(\Psi, \tau) U(X_t; \Psi) \\ &+ \frac{1}{2} \sum_{i=1}^n \sum_{j=1}^n B_i(\Psi, \tau) B_j(\Psi, \tau) \Sigma_i(X_t; \Psi) \Sigma_j(X_t; \Psi)', \end{split}$$

where  $\Sigma_i(\cdot)$  is the *i*-th row of matrix  $\Sigma(\cdot)$ .

## Appendix B

**Theorem.** Assume that the  $n \times 1$  vector  $X_t$  obeys the following dynamic:

$$dX_t = U(X_t; \Psi)dt + \Sigma(X_t; \Psi)dW_t$$

Suppose that  $U(X_t; \Psi)$  and  $\Sigma(X_t; \Psi)\Sigma(X_t; \Psi)'$  are affine functions of  $X_t$  so that  $U(X_t; \Psi)$  can be written as  $G + KX_t$  with G and K being matrices of dimension  $n \times 1$  and  $n \times n$ . The mean and variance of  $X_{t+h}$ , conditional on  $X_t$ , are affine functions of  $X_t$  if K is diagonable.

**Remark.** A matrix is diagonable if all of its eigenvalues are distinct. The assumption of diagonability does not involve an appreciable loss of generality. Since the eigenvalues of a matrix are continuous functions of its elements, if the matrix *K* has multiple eigenvalues, a slight alteration to any element produces a neighboring system with distinct roots that, for all practical purposes, is the same as the original system. Therefore the statement of the theorem is a generic result.

*Proof.* For notational convenience, let  $\Sigma(X_t; \Psi) = \Sigma_t$ . The integral representation for  $X_{t+h}$  can be written as:

$$X_{t+h} = X_t + \int_t^{t+h} [G + KX_s] ds + \int_t^{t+h} \Sigma_s dW_s.$$

Denote the eigenvalue decomposition of K, usually a non-symmetric matrix, by  $QkQ^{-1}$  where  $Q^{-1}Q = I$  and k a square diagonal matrix. Pre-multiplying the above equation by  $Q^{-1}$  yields:

$$Y_{t+h} = Y_t + \int_t^{t+h} [g + kY_s] ds + \int_t^{t+h} \sigma_s dW_s,$$

where  $Y_t = Q^{-1}X_t$ ,  $g = Q^{-1}G$ , and  $\sigma_t = Q^{-1}\Sigma_t$ . Taking conditional expectation gives rise to:

$$E(Y_{t+h}|Y_t) = Y_t + \int_t^{t+h} [g + kE(Y_s|Y_t)]ds.$$

The solution to this integral equation is:

$$E(Y_{t\perp h}|Y_t) = e^{kh}Y_t - (e^{kh} - I)k^{-1}g.$$

The solution for  $E(X_{t+h}|X_t)$  can be recovered from the above solution which is an affine function of  $X_t$ .

The conditional variance of  $Y_t$  can be computed using:

$$Var(Y_{t+h}|Y_t) = E(Y_{t+h}Y'_{t+h}|Y_t) - E(Y_{t+h}|Y_t)E(Y_{t+h}|Y_t)'.$$

The elements of  $E(Y_{t+h}Y'_{t+h}|Y_t)$  can be computed as follows. A  $[(n+1)n/2] \times 1$  vector  $f_t$  is formed by stacking the individual elements of the upper triangle of  $Y_{t+h}Y'_{t+h}$ . That is let:

$$f_{t} = \begin{bmatrix} Y_{1,t}Y_{1,t} \\ Y_{1,t}Y_{2,t} \\ \vdots \\ Y_{1,t}Y_{n,t} \\ Y_{2,t}Y_{2,t} \\ Y_{2,t}Y_{3,t} \\ \vdots \\ Y_{n,t}Y_{n,t} \end{bmatrix} = \begin{bmatrix} f_{t}^{1,1} \\ f_{t}^{1,2} \\ \vdots \\ f_{t}^{1,n} \\ f_{t}^{2,2} \\ f_{t}^{2,3} \\ \vdots \\ f_{t}^{n,n} \end{bmatrix}.$$

Using Ito's lemma on  $f_t$  yields:

$$f_{t+h} = f_t + \int_t^{t+h} \left[ \frac{\partial f_s}{\partial Y_s} u_s + \mathcal{D}_s ds \right] + \int_t^{t+h} \sigma_s \frac{\partial f_s}{\partial Y_s} dW_s,$$

where  $u_t = g + kY_t$  with k being the diagonal matrix defined earlier. Moreover,

$$\frac{\partial f_t}{\partial Y_t} = \begin{bmatrix} \frac{\partial f_t^{1,1}}{\partial Y_{1,t}} & \frac{\partial f_t^{1,1}}{\partial Y_{2,t}} & \dots & \frac{\partial f_t^{1,1}}{\partial Y_{n,t}} \\ \vdots & \vdots & \ddots & \vdots \\ \frac{\partial f_t^{n,n}}{\partial Y_{1,t}} & \frac{\partial f_t^{n,n}}{\partial Y_{2,t}} & \dots & \frac{\partial f_t^{n,n}}{\partial Y_{n,t}} \end{bmatrix},$$

$$\mathcal{D}_{t} = \frac{1}{2} \begin{bmatrix} \operatorname{tr}(\sigma_{t}\sigma_{t}^{'}\frac{\partial^{2}f_{t}^{1,1}}{\partial Y_{t}\partial Y_{t}^{\prime}}) \\ \vdots \\ \operatorname{tr}(\sigma_{t}\sigma_{t}^{'}\frac{\partial^{2}f_{t}^{n,n}}{\partial Y_{t}\partial Y_{t}^{\prime}}) \end{bmatrix},$$

and

$$\frac{\partial^2 f_t^{i,j}}{\partial Y_t \partial Y_t'} = \begin{bmatrix} \frac{\partial^2 f_t^{i,j}}{\partial Y_{1,i} \partial Y_{1,i}} & \cdots & \frac{\partial^2 f_t^{i,j}}{\partial Y_{1,i} \partial Y_{n,i}} \\ \vdots & \ddots & \vdots \\ \frac{\partial^2 f_t^{i,j}}{\partial Y_{n,i} \partial Y_{1,i}} & \cdots & \frac{\partial^2 f_t^{i,j}}{\partial Y_{n,i} \partial Y_{n,i}} \end{bmatrix}.$$

Since  $u_t$  is affine in  $Y_t$ , we can write:

$$\frac{\partial f_t}{\partial Y_t} u_t = pY_t + qf_t,$$

where p is a matrix of dimension  $[(n+1)n/2] \times n$  and q a diagonal matrix of dimension  $[(n+1)n/2] \times [(n+1)n/2]$  given by:

$$q = \begin{bmatrix} k_{1,1} + k_{1,1} & 0 & \dots & 0 \\ 0 & k_{1,1} + k_{2,2} & \dots & 0 \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \dots & k_{n,n} + k_{n,n} \end{bmatrix},$$

where  $k_{i,j}$  is element i,j of matrix k. Since  $\sigma_t \sigma_t'$  is affine in  $Y_t$ , it follows that:

$$\mathscr{D}_t = w + vY_t,$$

where w and v are matrices of dimension  $[(n+1)n/2] \times 1$  and  $[(n+1)n/2] \times n$ , respectively. The stochastic integral equation can thus be rewritten as:

$$f_{t+h} = f_t + \int_t^{t+h} [pY_s + qf_s + w + vY_s] ds + \int_t^{t+h} \sigma_s \frac{\partial f_s}{\partial Y_s} dW_s.$$

Taking conditional expectation yields:

$$E(f_{t+h}|Y_t) = f_t + \int_t^{t+h} [z_s + qE(f_s|Y_t)]ds,$$

where  $z_s = w + (p + v)E(Y_s|Y_t)$ . Clearly,  $z_s$  is a function of s and is an affine function of  $Y_t$ . The solution to this integral equation is

$$E(f_{t+h}|Y_t) = e^{qh}f_t + \int_t^{t+h} e^{q(t+h-s)}z_s ds.$$

As this solution shows, the only non-affine element, in terms of  $Y_t$ , for  $E(Y_{t+h}Y'_{t+h}|Y_t)$  is given by

$$e^{qh}f_t$$
.

Straightforward computations for  $E(Y_{t+h}|Y_t)E(Y_{t+h}|Y_t)'$ , shows that the only non-affine element is given by

$$e^{kh}Y_{t}Y_{t}'e^{kh}$$
.

The upper triangular elements of this matrix, when properly staked up, are precisely the

elements of  $e^{qh}f_t$ . Therefore  $Var(Y_{t+h}|Y_t)$  is affine in  $Y_t$ . Converting  $Y_t$  back to  $X_t$  gives rise to the result that  $Var(X_{t+h}|X_t)$  is affine in  $X_t$ .

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#### **Notes**

- 1. A complete description of the Kalman filter recursion can be found in Harvey (1990).
- 2. The algorithm used to simulate the Poisson random variate is taken from Devroye (1981).
- 3. The algorithm used to simulate the gamma random variate is taken from Devroye (1986).

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