

Estimation of Dynamic Term Structure Models

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ABSTRACT

We study the finite-sample properties of some of the standard techniques used to estimate modern term structure models. For sample sizes and models similar to those used in most empirical work, we reach three surprising conclusions. First, while maximum likelihood works well for simple models, it produces strongly biased parameter estimates when the model includes a flexible specification of the dynamics of interest rate risk. Second, despite having the same asymptotic efficiency as maximum likelihood, the small-sample performance of Efficient Method of Moments (a commonly used method for estimating complicated models) is unacceptable even in the simplest term structure settings. Third, the linearized Kalman filter is a tractable and reasonably accurate estimation technique, which we recommend in settings where maximum likelihood is impractical.

JEL classification: C14, C52, G12, G13.

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1 Introduction

Starting with Vasicek (1977) and Cox, Ingersoll, and Ross (1985), an enormous literature has focused on building and estimating dynamic models of the term structure. By specifying particular functional forms for both the risk-neutral dynamics of short-term interest rates and the compensation investors require to bear interest rate risk, these models describe the evolution of yields at all maturities. Much of the literature focuses on the affine class characterized by Duffie and Kan (1996). This class allows multiple state variables to drive interest rates and has the computationally convenient feature that bond yields are linear functions of these variables. The first generation of affine models, including multivariate generalizations of Vasicek and Cox et al., imposed two specializing assumptions: the state variables are independent and the price of risk is a multiple of interest rate volatility. Given these restrictions, estimation of the models' parameters is reasonably simple.

The estimation results revealed limitations in the models. For example, Dai and Singleton (2000) find strong evidence of nonzero correlations among the state variables. Duffee (2002) finds that the restriction on the price of risk implies unrealistic behavior for bonds' excess returns. Moreover, there is evidence of nonlinearity in expected interest rate movements that is inconsistent with these models.¹ In response to these limitations, researchers have introduced more flexible, "second generation" models. For example, Dai and Singleton (2000) estimate affine models in which the state variables are allowed to be correlated, while retaining the assumption that the price of risk is proportional to volatility. Duffee (2002) constructs a multifactor affine model with a more general specification for the dynamics of the price of risk than in Dai and Singleton (2000). Duarte (2004), Ahn, Dittmar, and Gallant (2002), and Leippold and Wu (2002) construct models with fairly general specifications of the price of risk that produce nonlinear dynamics.²

Although these new models are a significant improvement over earlier models, their ability to capture the observed dynamics of bond yields is not yet clear because the corresponding empirical literature is relatively immature. This is due to both the quick pace of the modeling advances and the difficulty of estimating some of these more complex models. Maximum likelihood is asymptotically efficient, but its finite-sample properties in the context of these models are not clear. Moreover, for many of these models the probability distribution of discretely sampled bond yields is unknown or intractable. Alternative techniques include

¹Nonlinearities are documented in Pfann, Schotman, and Tschernig (1996), Ait-Sahalia (1996), Conley, Hansen, Luttmer, and Scheinkman (1997), and Stanton (1997).

²Other nonlinear models include Longstaff (1989), Beaglehole and Tenney (1992), Constantinides (1992), and Ahn and Gao (1999). The latter two have prices of risk that are more flexible than those in Dai and Singleton (2000). Stanton (1997) and Boudoukh, Downing, Richardson, Stanton, and Whitelaw (2010) are nonparametric (and therefore nonlinear) models of both physical drifts and prices of risk.

moment-based methods and simulation methods. The optimal technique is difficult to determine due to our limited understanding of the properties of these techniques when applied to sophisticated term structure models.

In this paper we study the finite-sample properties of three prominent techniques used to estimate second-generation term structure models, in settings close to those facing researchers estimating these models. The first technique we study is maximum likelihood (the method of choice when the state variables can be precisely inferred by the econometrician), used, for example, by Pearson and Sun (1994), Chen and Scott (1993), Aït-Sahalia and Kimmel (2010), and Brandt and He (2002). We employ exact maximum likelihood when the likelihood function is known, and simulated maximum likelihood when it is not, using the method of Pedersen (1995) and Santa-Clara (1995). The second technique we study is the Efficient Method of Moments of Gallant and Tauchen (1996) (used, for example, by Dai and Singleton, 2000; Ahn et al., 2002; Andersen and Lund, 1997; Gallant and Tauchen, 1997). It is the most widely used alternative when maximum likelihood is infeasible, because it is tractable and can attain the same asymptotic efficiency as maximum likelihood. This technique uses an auxiliary model. We follow common practice by choosing a semi-nonparametric model; thus we refer to this technique as EMM/SNP. The final method is a variant of the Kalman filter. This may or may not be maximum likelihood, depending on the setting. For each technique, we use Monte Carlo simulations to determine the behavior of the estimators for sample sizes and models similar to those used in most empirical work. To keep the size of the paper manageable, we restrict our attention to models in the affine class with independent factors. We examine models in both the “essentially affine” class of Duffee (2002) and the “semi-affine” class of Duarte (2004).

Though we consider only a few of the infinite number of possible models, our results allow us to draw three main conclusions that apply much more generally. Each stands in surprising contrast to what we would expect based solely on asymptotic considerations. The first conclusion is that when the term structure model does not include a highly restrictive form of the price of risk, maximum likelihood does a poor job of estimating the parameters that determine expected changes in interest rates. The estimates are strongly biased and estimated with little precision. One implication is that conditional expectations of bond returns implied by maximum likelihood parameter estimates differ substantially from their true values. This behavior is related to the well-known downward bias in estimates of the speed of mean reversion of highly persistent processes such as bond yields. While Ball and Torous (1996) find that high persistence in bond yields does not pose a serious problem in estimating first-generation models, we find that it causes major problems in estimating second-generation models due to their more flexible specification of the dynamics of the price

of risk.

Our second conclusion is that the performance of EMM/SNP is unacceptable in even the simplest first-generation term structure settings, where maximum likelihood methods work well. This conclusion is particularly surprising because EMM/SNP attains the same asymptotic efficiency as maximum likelihood. The explanation for this poor performance involves the asymmetry of the EMM criterion function and is thus fairly technical. Section 8 describes the basic intuition, and a more complete discussion of the problem is in Duffee and Stanton (2008). The main result is that while the high persistence of bond yields drives a wedge between the finite-sample and asymptotic properties of all estimation techniques, the consequences of this wedge are much more dramatic for EMM.

This is a discouraging conclusion because EMM/SNP is a tractable method for estimating models for which maximum likelihood is infeasible. However, our third conclusion is more positive. We find that the Kalman filter is a reasonable choice even when it does not correspond to maximum likelihood. In many second-generation term structure models the standard Kalman filter cannot be implemented because the first and second moments of discretely observed bond yields are unknown. For these models we advocate the use of a modified Kalman filter that uses linearized instantaneous term structure dynamics. Although this method is inconsistent, its finite-sample biases are similar to the biases associated with maximum likelihood.

In the next section we describe the estimation techniques that we examine in the remainder of the paper. The specific term structure models that we consider are discussed in Section 3, and Section 4 gives details of the simulation procedure. Sections 5 and 6 present results for one-factor term structure models (Gaussian and square root respectively), Section 7 presents results for two factor models, Section 8 investigates in more detail the performance of EMM/SNP, and concluding comments are offered in Section 9.

2 Estimation techniques

This section outlines the three estimation techniques whose small-sample properties we study in the rest of the paper: maximum likelihood (ML), EMM/SNP, and a variant of the Kalman filter. Although we perform our Monte Carlo analysis using specific models presented in Section 3, our discussion here is more general.

The data are a panel of bond yields. They are equally spaced in the time series, at intervals $t = 1, \dots, T$. The random vector y_t represents a length- m vector of bond yields. Denote the history of yields through t as $Y_t = (y'_1, \dots, y'_t)'$. Yields are a function of a length- n

latent state vector x_t and (perhaps) a latent noise vector w_t :

$$y_t = y(x_t, w_t, \rho_1). \quad (1)$$

The noise may represent market microstructure effects or measurement error. Although these are certainly plausible features of the data, the main role played in the literature by this noise is to give the model flexibility to fit high-dimensional data with a low-dimensional state vector. The vector ρ_1 contains the parameters of this function. The state vector follows a diffusion process

$$dx_t = \mu(x_t, \rho_2) dt + \sigma(x_t, \rho_2) dz_t, \quad (2)$$

where ρ_2 is the parameter vector. The density function associated with the noise is

$$g_w(w_1, \dots, w_T). \quad (3)$$

For simplicity, we assume that the distribution of the noise is independent of x_t .

A term structure model (including a description of noise in bond yields) implies functional forms for (1), (2), and (3). We are interested in the resulting probability distribution of yields. Stack the parameter vectors ρ_1 and ρ_2 into ρ . Then we can always write the log density function of the data as

$$\log g_{Y_T}(Y_T) = \sum_{t=1}^T \log g_t(y_t | Y_{t-1}; \rho),$$

where $g_1(y_1 | Y_0; \rho)$ is interpreted as the unconditional distribution of y_1 . The true parameter vector is denoted ρ_0 . The primary difficulty in estimating ρ_0 with this structure is that the functional form for $g_t(\cdot)$ is often unknown or intractable.

2.1 Maximum Likelihood

The maximum likelihood estimator is the value that maximizes the (log) likelihood. Due to its asymptotic efficiency, maximum likelihood is the estimation method of choice in almost any econometric setting where we can evaluate the likelihood function. In many dynamic term structure models there is a one-to-one mapping between a length- n x_t and n bond yields. Thus we can pick any n points on the date- t yield curve, assume these yields have no noise, and invert the appropriate pricing equations to infer x_t .³ This commonly adopted

³When the true parameters are used, this inversion always produces an admissible state vector. However, for an arbitrary parameter vector the resulting state vector may be inadmissible. For example, observed bond yields might imply negative values for state variables that ought never to be negative.

approach was first used by Pearson and Sun (1994) and Chen and Scott (1993). In principle, when we can identify the state we can estimate the model with maximum likelihood because the likelihood function can be expressed as the solution to a partial differential equation involving the functions μ and σ in (2).⁴ However, this equation can be solved in closed form for only for a few special cases such as the square-root diffusion model of Cox et al. (1985) model (hereafter CIR) and the Gaussian diffusion model of Vasicek (1977). In many other models the equation can only be solved numerically, making direct maximum likelihood via this approach infeasible.

Another potential problem with the use of maximum likelihood is that when we observe $m > n$ bond yields, there is in general no set of values of x_t that exactly matches the m bond yields every period. One way to circumvent this difficulty is to assume that only n of the yields are measured without error and allow for noise in the remaining yields. This does not add to the difficulty of using maximum likelihood, but has the disadvantage that the choice of the yields estimated without error is, necessarily, somewhat ad hoc. If we are unwilling to accept this assumption, or if the term structure model does not imply a one-to-one mapping between x_t and bond yields even in the absence of noise as in Ahn, Dittmar, and Gallant (2002), our inability to infer x_t exactly will make maximum likelihood estimation even more difficult.

Much progress has been made recently in expanding the settings where maximum likelihood is possible. Pedersen (1995) and Santa-Clara (1995) develop a simulation-based approach that allows the approximation of the likelihood function when the state is observable and the likelihood function is intractable. The idea is to split each observation interval into small subintervals. The conditional distribution of the state approaches the normal distribution as the length of the subintervals shrinks towards zero. If a particular observation interval is split into k pieces, the method involves simulating a large number of paths for the first $k - 1$ of the subintervals, then for each path calculating the likelihood of jumping from the value at subperiod $n - 1$ to the (next observed) value at subperiod k . As both the number of simulated paths and the number of subintervals per observation become large, the average of these normal likelihoods converges to the true likelihood of moving from one observed value to the next. Aït-Sahalia (1999, 2008) proposes an alternative estimation procedure for this case. He develops a series of approximations to the likelihood function that are tractable to estimate and converge to the true likelihood function. Aït-Sahalia and Kimmel (2010) apply this technique to term structure modeling. The econometrician controls the accuracy of the approximation by choosing the order of the approximating series. Finally, when the under-

⁴This equation is known as the Kolmogorov forward equation (see Øksendal, 2002, for further information). Lo (1988) describes the use of this equation to implement maximum likelihood.

lying state cannot be exactly observed due to measurement error, Brandt and He (2002) and Bates (2006) have recently developed techniques that allow calculation of an approximate likelihood function for certain classes of term structure models.

We estimate a variety of models with maximum likelihood. The implementation differs depending on the model. If the model is Gaussian, we assume that bond yields are all observed with error and use the Kalman filter, as discussed in Section 2.3. If the model is not Gaussian we assume that $m = n$ bond yields are observed without error. In this case we use the exact likelihood function if it is known. If not, we use the Pedersen/Santa-Clara technique, following the implementation in Brandt and Santa-Clara (2002). Because the technique is designed to simulate conditional densities rather than unconditional densities, we condition the likelihood of the data on the first observation.

2.2 Efficient Method of Moments

When maximum likelihood is infeasible, the most commonly used method for estimating term structure models is the Efficient Method of Moments (EMM), a path simulation method. Simulations produced with the dynamic model are used to draw indirect inferences about the density function $g_{Y_T}(Y_T)$. These simulations can be used to calculate arbitrary population moments as functions of the parameters of the process being estimated, which can be compared with sample moments estimated from the data.⁵ Since it is never evaluated, the true density function $g_{Y_T}(Y_T)$ can be intractable or even unknown, and the data can be observed with or without noise, since adding noise to simulated data is trivial. The defining characteristic of EMM is the choice of moments to simulate. Following Gallant and Tauchen (1996), EMM uses the score vector from some tractable auxiliary model. Although the technique is well-known, we go through the details here to motivate our later discussion of the finite-sample behavior of EMM.

Let f be some auxiliary function that (perhaps approximately) expresses the log density of y_t as a function of Y_{t-1} and a parameter vector γ_0 :

$$f(y_t | Y_{t-1}; \gamma_0).$$

The first step in EMM is to calculate the parameters of the auxiliary function that maximize the (pseudo) log likelihood. Equivalently, the parameter vector γ_T is the vector that sets the

⁵Duffie and Singleton (1993) discuss the properties of simulation estimators in general.

sample mean of the derivative of the log-likelihood function to zero.

$$\frac{1}{T} \sum_{t=1}^T \left(\frac{\partial f(y_t | Y_{t-1}; \gamma)}{\partial \gamma} \Big|_{\gamma=\gamma_T} \right) = 0.$$

The Central Limit Theorem implies that

$$\sqrt{T}(\gamma_T - \gamma_0) \xrightarrow{d} N(0, d^{-1} S d^{-1}) \quad (4)$$

where the convergence is in distribution. The matrices S and d are defined as

$$S = E \left[\begin{pmatrix} \frac{\partial f}{\partial \gamma} \\ \frac{\partial f}{\partial \gamma'} \end{pmatrix} \Big|_{\gamma=\gamma_0} \right] \quad (5)$$

and

$$d = E \left(\frac{\partial f}{\partial \gamma \partial \gamma'} \Big|_{\gamma=\gamma_0} \right).$$

Intuitively, d transforms the variability of the moment vector S into the variability of the auxiliary parameters.

The second step in EMM is to simulate a long time series $\hat{Y}_N(\rho) = (\hat{y}_1(\rho)', \dots, \hat{y}_N(\rho)')'$ using the true model (1), (2) and (3). If the discrete density of bond yields conditional on ρ is known, yields can be generated from this density. Otherwise the continuous process (2) is discretized.⁶ The simulated time series is used to calculate the expectation of the score vector associated with the auxiliary model:

$$m_T(\rho, \gamma_T) = \frac{1}{N} \sum_{\tau=1}^N \frac{\partial f(\hat{y}_\tau(\rho) | \hat{Y}_{\tau-1}(\rho); \gamma)}{\partial \gamma} \Big|_{\gamma=\gamma_T}. \quad (6)$$

The arguments of f in (6) are explicit to show that the score vector is calculated using the combination of simulated yields and parameters from the original data Y_T . As N approaches infinity, this sample mean approaches the expectation of the score vector evaluated at γ_T :

$$\lim_{N \rightarrow \infty} m_T(\rho, \gamma_T) = E \left(\frac{\partial f(y_t(\rho) | Y_{t-1}(\rho); \gamma)}{\partial \gamma} \Big|_{\gamma=\gamma_T} \right),$$

where Y_{t-1} and y_t are drawn from the distribution of bond yields as determined by ρ .

⁶Discretization techniques are discussed in Kloeden and Platen (1992).

The Central Limit Theorem determines the asymptotic distribution of m_T :

$$\sqrt{T}m_T(\rho_0, \gamma_T) \xrightarrow{d} N(0, C(\rho_0)d^{-1}Sd^{-1}C(\rho_0)) \quad (7)$$

where

$$C(\rho) = \lim_{T \rightarrow \infty} \left(\frac{\partial m_T(\rho, \gamma)}{\partial \gamma'} \Big|_{\gamma=\gamma_T} \right) = \frac{\partial m_T(\rho, \gamma)}{\partial \gamma} \Big|_{\gamma=\gamma_0}.$$

The inner part of the variance-covariance matrix in (7) is the variance-covariance matrix of the auxiliary parameters from (4). This inner matrix is pre- and post-multiplied by the sensitivity of m_T to the auxiliary parameters.

The distribution in (7) can be simplified by recognizing that $C(\rho_0) = d$, so that

$$\sqrt{T}m_T(\rho_0, \gamma_T) \xrightarrow{d} N(0, S).$$

This asymptotic result leads to the EMM estimator

$$\rho_T = \underset{\rho}{\operatorname{argmin}} m_T(\rho, \gamma_T)' S_T^{-1} m_T(\rho, \gamma_T). \quad (8)$$

where S_T is the sample counterpart to (5):

$$S_T = \frac{1}{T} \sum_{t=1}^T \left[\left(\frac{\partial f}{\partial \gamma} \right) \left(\frac{\partial f}{\partial \gamma'} \right) \Big|_{\gamma=\gamma_T} \right].$$

An estimate of the asymptotic variance-covariance matrix of $\hat{\rho}_T$ is

$$\Sigma_T = \frac{1}{T} [(M_T)' S_T^{-1} (M_T)]^{-1},$$

where

$$M_T = \frac{\partial m_T(\rho, \gamma_T)}{\partial \rho'} \Big|_{\rho=\rho_T}.$$

If there are more moment conditions (length of γ) than parameters (length of ρ), then under the null hypothesis,

$$J = T m_T(\rho_T, \gamma_T)' S_T^{-1} m_T(\rho_T, \gamma_T)$$

is asymptotically distributed as a $\chi^2(q)$ random variable, where q is the number of over-identifying moment conditions.

This estimation procedure does not specify which auxiliary log-likelihood function to use. Gallant and Tauchen (1996) note that if the distribution implied by the auxiliary model is

close to that implied by the true underlying model, then the estimates obtained should be close to those obtained using maximum likelihood. A common choice of auxiliary model is a semi-nonparametric (SNP) description of the data (outlined in Appendix A), in large part motivated by its asymptotic properties. Gallant and Long (1997) show that with this choice of auxiliary model, EMM asymptotically attains the efficiency of maximum likelihood.

Although it is desirable to use estimation techniques that have good asymptotic properties, their finite-sample properties are more important in practice. The finite-sample properties of EMM/SNP have been studied in several contexts.⁷ However, this earlier work has not examined settings that contain the salient features of bond yield data: highly persistent and highly correlated multivariate time series. The most relevant work is Zhou (2001), who studies methods to estimate the parameters of a square-root diffusion model of the instantaneous interest rate. In this univariate setting he finds that when the data are highly persistent, the performance of EMM/SNP is mixed. There is good reason to suspect that the performance of EMM/SNP will deteriorate in a multivariate setting. Efficient Method of Moments is a GMM estimator, and it is well-known that the finite-sample properties of GMM can deteriorate seriously as the number of over-identifying restrictions increases (see, for example, Tauchen, 1986; Kocherlakota, 1990; Ferson and Foerster, 1994; Hansen et al., 1996). Because SNP puts little structure on data, the number of SNP parameters that are used to summarize a multivariate time series can be large. An SNP specification uses a minimum of $m(m+1)(3/2)$ parameters to fit an m -dimensional time series, and usually many more than this.

2.3 The Kalman Filter

Filtering is a natural approach when the underlying state is unobserved. The Kalman filter corresponds to ML when the state vector dynamics are Gaussian and the noise is also normally distributed. In non-Gaussian settings, given an analytic conditional density for the state vector, exact nonlinear filtering is possible but numerically demanding, especially for nonscalar x_t . We are unaware of any empirical term-structure implementations of exact filtering when the dynamics of x_t are nonlinear.⁸

Approximate linear filtering is easier to implement. The Kalman filter has been applied to term structure models in which x_t has affine dynamics and thus analytic expressions of

⁷See, for example, Chumacero (1997), Andersen, Chung, and Sørensen (1999), and Andersen and Sørensen (1996).

⁸The exact filter of Kitagawa (1987) is implemented by Lu (1999) for a Constantinides (1992) model, which has Gaussian dynamics for the state vector. For a discussion of the high computational cost of Kitagawa's filter, see the comments by Kohn and Ansley (1987) and Martin and Raftery (1987).

the first two moments of the conditional density are available.⁹ Outside the Gaussian class of term structure models, parameter estimates obtained directly from Kalman filter estimation are inconsistent. There is Monte Carlo evidence that when the underlying model is linear but heteroskedastic, the inconsistency may be of limited importance in practice.¹⁰

In this paper we examine the empirical performance of a variant of the Kalman filter. To introduce this variant we first review the extended Kalman filter. The observation equation expresses observed yields, y_t , as a linear function of the unobservable state, x_t , plus measurement error ϵ_t . The transition equation expresses the discrete-time evolution of x_t as linear in x_t . These equations are determined by the parameters of the term structure model ρ . The term “extended” means that the parameters of the linear functions may depend on the underlying state x_t . The structure is

$$y_t = H_0(\rho) + H_1(\rho)'x_t + \epsilon_t; \tag{9}$$

$$x_{t+1} = F_0(x_t, \rho) + F_1(x_t, \rho)x_t + v_{t+1}; \tag{10}$$

$$E(\epsilon_t) = 0; E(v_{t+1}) = 0; E(\epsilon_t \epsilon_t') = R(\rho); E(v_{t+1} v_{t+1}') = Q(x_t, \rho).$$

In (9), H_0 and H_1 are not functions of x_t . Although there are no additional complications introduced by allowing for such dependence, the term structure models we examine have pricing formulas that satisfy (9). The contemporaneous prediction of the state vector and its associated variance-covariance matrix are denoted $x_{t|t}^p$ and $P_{t|t}$ respectively. One-step-ahead forecasts of the state vector and observable vector, are denoted $x_{t+1|t}^p$ and $y_{t+1|t}^p$ and the variance-covariance matrices of these forecasts are denoted $P_{t+1|t}$ and $V_{t+1|t}$ respectively.

The extended Kalman filter is estimated using the standard Kalman filter recursion, though the resulting parameter estimates are generally inconsistent. The recursion begins with a candidate parameter vector ρ . This vector is used to calculate an unconditional expectation and variance-covariance matrix for x_1 , which we can denote $x_{0|0}^p$ and $P_{0|0}$. (If closed-form expressions for these moments are unavailable, the moments can be produced with simulations.) The steps in the recursion are

1. Use $x_{t|t}^p$ and ρ to evaluate the matrices $F_0(x_t, \rho)$, $F_1(x_t, \rho)$, and $Q(x_t, \rho)$. Denote these values by F_{0t} , F_{1t} , and Q_t . This is the step that creates inconsistency in the estimates, because the filtered x_t is used instead of the (unknown) true x_t .¹¹

⁹Applications include Pennacchi (1991), Chen and Scott (2003), Duan and Simonato (1999), Lund (1997), de Jong (2000), Geyer and Pichler (1999), and Jegadeesh and Pennacchi (1996).

¹⁰Some results are in de Jong (2000) and Duan and Simonato (1999). In addition, in certain cases, as in Lund (1997), the approximation error can be reduced using iterative techniques or numerical integration.

¹¹The distribution of the shock also affects the consistency of the estimates. See Duan and Simonato (1999) for a detailed discussion.

2. Compute the one-period-ahead prediction and variance of x_{t+1} , $x_{t+1|t}^p = F_{0t} + F_{1t}x_{t|t}^p$ and $P_{t+1|t} = F_{1t}P_{t|t}F_{1t}' + Q_t$.
3. Compute the one-period-ahead prediction and variance of y_{t+1} , $y_{t+1|t}^p = H_0 + H_1'x_{t+1|t}^p$ and $V_{t+1|t} = H_1'P_{t+1|t}H_1 + R$.
4. Compute the forecast error in y_{t+1} , $e_{t+1} = y_{t+1} - y_{t+1|t}^p$.
5. Update the prediction of x_{t+1} , $x_{t+1|t+1}^p = x_{t+1|t}^p + P_{t+1|t}H_1V_{t+1|t}^{-1}e_{t+1}$ and $P_{t+1|t+1} = P_{t+1|t} - P_{t+1|t}H_1V_{t+1|t}^{-1}H_1'P_{t+1|t}$.

The estimated parameter vector ρ_T solves

$$\rho_T = \operatorname{argmax}_{\rho} \sum_{t=1}^T f(e_t, V_{t|t-1}),$$

where the period- t approximate log-likelihood is

$$f(e_t, V_{t|t-1}) = -\frac{1}{2}[m \log(2\pi) + \log |V_{t|t-1}| + e_t' V_{t|t-1}^{-1} e_t].$$

This filter requires a closed-form expression for the discrete-time dynamics of x_t . In many term structure settings there is no such expression. In such cases we advocate the use of a variant of the Kalman filter, where (10) is replaced with a linearization of the instantaneous dynamics of x_t in (2). The linearization is taken in the neighborhood of $x_{t|t}^p$. The time between discrete observations is denoted Δt . The linearization is (suppressing the dependence on parameters)

$$x_{t+1} = F_{0t} + F_{1t}x_t + v_{t+1}; \tag{11}$$

$$F_{0t} = \left(\mu(x_{t|t}^p) - \frac{\partial \mu(x_t)}{\partial x_t'} \bigg|_{x_t=x_{t|t}^p} x_{t|t}^p \right) \Delta t; \tag{12}$$

$$F_{1t} = I + \frac{\partial \mu(x_t)}{\partial x_t'} \bigg|_{x_t=x_{t|t}^p} \Delta t; \tag{13}$$

$$Q_t = \sigma(x_{t|t}^p)\sigma(x_{t|t}^p)' \Delta t. \tag{14}$$

In (11) through (14), two new approximation errors are added to that caused by evaluating dynamics at the filtered x_t instead of at the true x_t . The first is the use of the instantaneous dynamics of x_t as a proxy for the discrete-time dynamics of x_t . The second is the linearization of these dynamics.

An estimate of the asymptotic variance-covariance matrix of the estimated parameters

ρ_T is based on the outer product of first derivatives of the log likelihood function,

$$\Sigma_T = \frac{1}{T^2} \sum_{t=1}^T \left[\left(\frac{\partial f}{\partial \rho} \right) \left(\frac{\partial f}{\partial \rho'} \right) \Big|_{\rho=\rho_T} \right].$$

Because the log-likelihood function is misspecified for non-Gaussian models, a theoretically more robust estimate of the variance-covariance matrix uses both first and second derivatives of the log likelihood function. In practice, however, we have found that when estimating term structure models, numerical difficulties in the calculation of second derivatives outweigh the value of using this more robust estimator.

3 Model description

The focus of this paper, as of most of the literature in this area, is on estimating models within the affine framework of Duffie and Kan (1996). There are n state variables, denoted $x_t \equiv (x_{t,1}, \dots, x_{t,n})'$. Uncertainty is generated by n independent Brownian motions. Under the equivalent martingale measure these are denoted $\tilde{z}_t \equiv (\tilde{z}_{t,1}, \dots, \tilde{z}_{t,n})'$; corresponding Brownian motions under the physical measure are represented without the tildes. The instantaneous nominal interest rate, denoted r_t , is affine in the state:

$$r_t = \delta_0 + \delta x_t.$$

Here, δ_0 is a scalar and δ is an n -vector. The equivalent-martingale dynamics of the vector x_t determine bond prices. We consider two special cases of this framework. In both of them the individual elements of x_t are independent. The first is when these elements follow Gaussian processes:

$$dx_{it} = (k\theta_i - k_i x_{it})dt + \sigma_i d\tilde{z}_{it}.$$

The second is when these elements follow square-root diffusion processes:

$$dx_{it} = (k\theta_i - k_i x_{it})dt + \sigma_i \sqrt{x_{it}} d\tilde{z}_{it}.$$

Duffie and Kan show that we can write the price and yield of a zero-coupon bond that matures at time $t + \tau$ in the form

$$P(x_t, \tau) = \exp[A(\tau) - B(\tau)'x_t], \tag{15}$$

$$Y(x_t, \tau) = (1/\tau)[-A(\tau) + B(\tau)'x_t]. \tag{16}$$

In (15) and (16), $A(\tau)$ is a scalar function and $B(\tau)$ is an n -valued function. Explicit solutions can be calculated for the special cases we consider, given the results of Vasicek (1977) and Cox et al. (1985).

3.1 The price of risk

The dynamics of x_t under the physical measure are determined by specifying the dynamics of the market price of risk. Defining π_s/π_t as the state price deflator for time- t pricing of time- s payoffs, we can write

$$\frac{d\pi_t}{\pi_t} = -r_t dt - \Lambda_t' dz_t. \quad (17)$$

The element i of the n -vector Λ_t represents the price of risk associated with the Brownian motion z_{it} . When the equivalent-martingale dynamics are pure Gaussian diffusions we use the following form for the price of risk:

$$\Lambda_{it} = \sigma_i^{-1}(\lambda_{i1} + \lambda_{i2}x_{it}).$$

This form is a special case of the “essentially affine” models in Duffee (2002). When the equivalent-martingale dynamics are pure square-root diffusions we use the following form:

$$\Lambda_{it} = \sigma_i^{-1}(\lambda_{i1} + \lambda_{i2}\sqrt{x_{it}}).$$

This “semi-affine” specification is introduced in Duarte (2004). In either case, when $\lambda_{i2} = 0$ we are in the “completely affine” world of Dai and Singleton (2000). When $\lambda_{i2} \neq 0$, the individual elements of Λ_t can change sign depending on the shape of the term structure (i.e., depending on the elements of x_t). Thus investors’ willingness to face certain types of interest-rate risk can switch sign in a way that is not possible when $\lambda_{i2} = 0$.

3.2 Interest rate dynamics under the physical measure

The general representation of the state price deflator’s dynamics in (17) allow us to write the dynamics of x_t under the physical measure. For the Gaussian case the physical dynamics are

$$dx_{it} = (k\theta_i + \lambda_{i1} - (k_i - \lambda_{i2})x_{it})dt + \sigma_i dz_{it}.$$

Note that under the physical measure the dynamics of x_{it} are also Gaussian. This equivalence between equivalent martingale and physical dynamics does not carry over to the case of square-root diffusions under the equivalent martingale measure. In this case the physical

dynamics are

$$dx_{it} = (k\theta_i + \lambda_{i1}\sqrt{x_{it}} - (k - \lambda_{i2})x_{it})dt + \sigma_i\sqrt{x_{it}}dz_{it}. \quad (18)$$

These dynamics are nonlinear if $\lambda_{i1} \neq 0$. Stationarity of x_t in the Gaussian case is equivalent to $k_i - \lambda_{i2} > 0$. The same condition ensures stationarity of x_{it} in (18). Stationarity is also ensured in this case if $k_i - \lambda_{i2} = 0$ and $\lambda_{i1} < 0$.

4 Details of the simulation procedure

We study one-factor and two-factor affine term structure models. For each choice of n , we consider both Gaussian dynamics and square-root diffusion dynamics. We further break down these models into the first-generation version (Vasicek for Gaussian and CIR for square-root diffusion) and a version that generalizes the price of risk. For Gaussian dynamics this generalization is an affine price of risk and for the square-root diffusion this is Duarte’s semi-affine price of risk.

The “true” parameters for all of these processes are based on the results of fitting the models to Treasury yields. We used implied zero-coupon month-end bond yields computed by Rob Bliss from coupon bond yields, and are indebted to him for sharing the data. For the two-factor semi-affine model the parameters were estimated using data from 1971 through 1998. For all other models the parameter estimates are based on data from 1970 through 2001. (The use of two samples is accidental.)

Most of the simulated data samples we examine contain 1000 weeks (a little more than 19 years) of bond yields. Typically the samples include yields on bonds with maturities of three months, one year, and ten years. Occasionally we consider only two yields. In this case we drop the one-year yield. We also occasionally consider five yields. In this case we add yields for maturities of six months and five years. In most of the simulations the bond yields are given by the sum of the model-implied bond yields and normally distributed noise. The noise in each bond’s yield is independent across other bonds and across time and has a maturity-independent variance V . When we use estimation techniques that rely on exact identification of the state, the ten-year bond yield is observed without error. With $n = 2$, the three-month bond yield is also observed without error. The simulated data are generated by discretizing the instantaneous dynamics (ten subperiods per week) using the Milstein Fortran code of Gallant and Tauchen.

For each set of simulated data we estimate the parameters of the term structure model using various techniques. Three of the techniques require simulations. For the Pedersen/Santa-Clara simulated ML procedure we divide each weekly interval into five subperiods and es-

timate the conditional distribution with 2500 Monte Carlo simulations. For the Kalman filter that uses linearized instantaneous dynamics, we estimate the unconditional first two moments of the state with a single path simulation of 5000 years. We set the length of the path simulation in EMM at 50,000 weeks. We experimented with longer simulations, but increasing the simulation length had no appreciable effect on the results. For each technique we compute parameter estimates and their associated standard errors. We use Gallant and Tauchen’s Fortran 77 code to estimate parameters using EMM/SNP. Our code to estimate parameters using the other techniques is written in Fortran 90 and calls IMSL optimization routines (Simplex and a quasi-Newton optimizer). We conduct an extensive search over the parameter space to find the optimal parameters.

Five hundred Monte Carlo simulations are produced for each model. The results from these simulations are discussed in the next three sections. The one-factor Gaussian model has a simple structure that allows us to pinpoint what is driving the finite-sample behavior of its parameter estimates. Therefore we devote the next section to these models. We then briefly examine the results of finite-sample estimation of one-factor square-root diffusion models, and finally consider two-factor models.

5 One-factor Gaussian models

In this section we make two major points. First, ML estimation of term structure models that allow for a general specification of the price of risk—more precisely, models that allow the drift of the state under the physical measure to be unrelated to the drift under the equivalent martingale measure—produces estimates of the dynamics of the price of risk that are strongly biased and imprecise. Second, the performance of EMM/SNP is substantially inferior to that of ML. In particular, the estimated asymptotic standard errors of the parameter estimates differ substantially from the finite-sample standard deviations of the parameter estimates.

For the one-factor Gaussian model the equivalent martingale dynamics of the instantaneous interest rate are

$$dr = (k\theta - kr_t)dt + \sigma d\tilde{z}_t \tag{19}$$

and the physical dynamics are

$$dr = (k\theta + \lambda_1 - (k - \lambda_2)r_t)dt + \sigma dz_t. \tag{20}$$

The Vasicek model sets $\lambda_2 = 0$. The more general form is an affine price of risk. As discussed in the previous section, the parameters of the model are estimated from the behavior of Treasury data over the past thirty years. For the Vasicek model the parameters are $k\theta =$

0.0084, $\lambda_1 = -0.005$, $k = 0.065$, and $\sigma = 0.0175$. The mean interest rate is 0.052. The half-life of a shock to r_t is almost eleven years; r_t is close to a random walk. For the more general Gaussian model we retain the same parameters identified under the equivalent martingale measure and set the price of risk parameters to $\lambda_1 = 0.005$ and $\lambda_2 = -0.14$.

Before we get into the details of term structure estimation, we take a look at the finite-sample properties of the instantaneous interest rate itself. Without cross-sectional information, we can only identify parameters that are identified under the physical measure: $k\theta + \lambda_1$, $k - \lambda_2$, and σ . For the purposes of this exercise we use the “true” parameters for the Vasicek model. Table 1 displays results from Monte Carlo simulation of ML estimation of the physical dynamics of r_t , given 1000 weeks of observations of r_t . The estimates of both $k\theta + \lambda_1$ and k are strongly biased. The bias in k is the standard finite-sample bias in estimates of the autoregressive parameter of an AR process that has a near unit root, as noted by Ball and Torous (1996). The bias in $k\theta + \lambda_1$ is created by the same bias, since the mean interest rate is $(k\theta + \lambda_1)/k$. To fit the sample mean with a biased estimate of k , the estimate of $k\theta + \lambda_1$ must also be biased. Also note that the sample standard deviations of these two parameter estimates are substantially larger (by 40 to 80 percent) than their corresponding mean standard errors. Below, we contrast these results with those from estimation of the complete dynamic term structure model.

5.1 ML estimation of Vasicek model

We estimate the model using a panel of bond yields. The panel gives us information about both the physical and equivalent-martingale dynamics of r_t , allowing us to estimate all of the model’s parameters. Recall that all bond yields are measured with normally distributed error with standard deviation \sqrt{V} . In this case the standard Kalman filter produces maximum likelihood estimates of the model’s parameters. The left side of Table 2 displays the results of 500 Monte Carlo simulations of Kalman filter estimation.

In contrast to ML estimation using only observations of r_t , here all of the estimated parameters are now unbiased (or, more precisely, there is no statistically significant bias given 500 simulations). In addition, except for λ_1 the mean standard errors are close to the sample standard deviations of the parameter estimates. The reason the bias disappears when using panel data is that the drift of r_t under the physical dynamics shares the parameter k with the drift of r_t under the risk-neutral dynamics. The cross-section contains precise information about the risk-neutral drift, thus for the purposes of estimating k the ML estimation procedure essentially discounts the imprecise information contained in the physical drift. This is why the standard deviation of the estimates of k using panel data is

less than two percent of the corresponding standard deviation using only time series data. Here we are simply restating the point made by Ball and Torous (1996) in the context of CIR-type models. Note that the panel contains much more information about the speed of mean reversion than does the time series of r_t even though the bond yields are observed with error. (No measurement error was introduced in the simulated time series of r_t .)

5.2 ML estimation of the general Gaussian model

We now extend the analysis to the affine specification for the price of interest rate risk. The price of risk is linear in r_t such that the speed of mean reversion under the physical measure $k - \lambda_2 = 0.205$. Therefore r_t is not as persistent as in the Vasicek model estimated above. As in that model, the Kalman filter corresponds to maximum likelihood.

The left side of Table 3 displays the results of 500 Monte Carlo simulations of ML estimation. Estimates of the parameters identified under the equivalent martingale measure are close to unbiased and, as in the Vasicek case, are estimated with high precision. (The variability in the estimates is a little higher here than in the Vasicek case.) However, estimates of the parameters identified only under the physical measure are strongly biased and highly variable. The bias in the estimate of λ_1 is about one standard deviation and the bias in the estimate of λ_2 is about minus one standard deviation.

These results for the one-factor general Gaussian model are characteristic of all of the models we examine that have a general specification of the price of risk. Because of the relative simplicity of this model, we can see clearly the source of the bias in the risk premium parameters. It is created by the decoupling of the risk-neutral dynamics from the physical dynamics. Recall that in the Vasicek model, the parameters identified under the risk-neutral measure help to pin down the drift of r_t under the physical measure. When the dynamics of the risk premia are more flexible, the physical and risk-neutral drifts have no common parameters. Therefore the parameters $(k\theta + \lambda_1)$ and $(k - \lambda_2)$ are determined exclusively by the time-series properties of bond yields. The near unit-root bias produces an upward-biased speed of mean reversion, and thus a downward-biased estimate of λ_2 (since k is determined by the cross section). Note that the bias in the speed of mean reversion, $0.374 - 0.140 = 0.234$, is almost identical to the bias in the speed of mean reversion we saw in Table 1, where only time-series information is used. In addition, the variability in the estimate of mean reversion from Table 1 is almost identical to the variability in the estimate of λ_2 here.

Figure 1 displays the true and estimated drift functions for both the Vasicek and generalized Vasicek model. Panel A corresponds to the Vasicek model and Panel B corresponds to the Gaussian model with an affine price of risk. The solid lines represent true drift func-

tions and the dashed lines represent drift functions implied by the mean parameter estimates from Kalman filter estimation. In Panel A these lines are indistinguishable. In Panel B, we see that the estimated drift function implies a too-high drift when r_t is below average and a too-low drift when r_t is above average. To give some intuition for the results, the true half-life of an interest rate shock in the affine-risk model is 3.38 years. The implied half-life is less than half as long, 1.58 years.

Another way to interpret the magnitude of this bias is to ask what implications it has for the dynamics of expected excess returns to bonds. In this model the instantaneous expected excess return to a bond with maturity τ is

$$\text{excess ret}_t = -\frac{1 - e^{-k\tau}}{k}(\lambda_1 + \lambda_2 r_t).$$

The intuition behind this expression is straightforward. The fraction in this expression is the sensitivity of a bond's log price to instantaneous interest rates. (For k close to zero it is approximately the maturity of the bond.) The term in parentheses is the difference between the true drift of r_t and the risk-neutral drift. The higher is the true drift relative to the risk-neutral drift, the lower is the expected excess return to the bond: investors are pricing bonds as if interest rates will rise less quickly (or fall more rapidly) than they are expected to under the physical measure.

The estimated model implies a much wider range of expected excess bond returns than does the true model for a given range of r_t . According to the true model, the expected excess return to a ten-year bond typically ranges from around zero to around six percent a year. Consider, for example, three values of r_t : one standard deviation below the mean, the mean, and one standard deviation above the mean. These values are 3.80 percent, 6.54 percent, and 9.27 percent respectively. The corresponding expected excess returns to a ten-year bond are 0.24, 3.05, and 5.87 percent per year. If we use the mean parameter estimates from the model to predict instantaneous expected excess returns at these various values of r_t , the range of expected excess returns would roughly triple. The expected excess returns are -2.04 , 5.48 , and 13.0 percent per year respectively.

The expectations hypothesis of interest rates is hard to reconcile with the behavior of Treasury yields. For example, Campbell and Shiller (1991) find that when the slope of the term structure is steep, expected excess returns to bonds are high. Backus, Foresi, Mozumdar, and Wu (2001) conclude that the evidence against the expectations hypothesis of interest rates is strengthened when we take into account finite-sample biases in tests of the hypothesis. Although their analysis focused on estimating simple regressions, the result carries over to estimating dynamic term structure models. The finite-sample bias we docu-

ment works against rejecting the expectations hypothesis. The negative bias in λ_2 implies that when r_t is low (and therefore the slope is steep), the expected excess returns implied by the model's parameter estimates are lower than those implied by the true model. This bias pushes the implied behavior of returns closer to that consistent with the expectations hypothesis.

The bias in the price of risk dynamics is an important feature of term structure estimation, thus it is worth taking a closer look at its determinants. We conduct a variety of experiments to determine the sensitivity of the bias to variations in the amount of information in the data sample. We vary the number of points along the yield curve that are observed, the amount of measurement error in bond yields, the frequency of observation, and the length of the sample period.

Given the nature of the bias, none of the results from these experiments are particularly surprising. More cross-sectional information (either more points on the yield curve or less measurement error) increases the accuracy of estimates of the parameters identified under the equivalent martingale measure but has little effect on the accuracy of estimates of the parameters identified under the physical measure. The relevant evidence is in Tables 4 and 5. When five points on the yield curve are used (the first set of results in Table 4), the standard deviations of the estimates of $k\theta$, k , σ , and \sqrt{V} drop to about 80 percent of the corresponding standard deviations when three points are used. However, this change has a minimal effect on the means and the standard deviations of λ_1 and λ_2 . The same pattern appears when only two points on the yield curve are used (the middle set of results in the table). Finally, when the standard deviation of measurement error in bond yields is set to 10 basis points instead of 60 (the final set of results in the table), the standard deviations of estimates $k\theta$, k , σ , and \sqrt{V} drop by about 3/4, but again the effect on estimates of the price of risk parameters is minimal.

Table 5 displays results based on varying the amount of information in the time series. In Panel A we consider both cutting the sample size in half and doubling the sample size. The results are easy to summarize. The standard deviations of all the parameter estimates decrease with the sample size. The bias in the estimates of the price of risk also decreases with the sample size. For example, doubling the sample to 2000 weeks (more than 38 years) cuts the bias in the estimates of both λ_1 and λ_2 in half. Nonetheless, the mean parameter estimates of k and λ_2 imply that the half-life of interest rate shocks (2.16 years) is less than 2/3 the actual half-life of interest rate shocks.

Monthly data are often used to estimate dynamic term structure models. Panel B reports results at this frequency. The first set of results is for 240 months, which is about the length of a sample of 1000 weeks. Thus a comparison of these results with those of Kalman filter

estimation in Table 3 reveals the effects of a decrease in the frequency of observation while holding the length of the sample period constant. This change decreases the precision of the parameter estimates identified under the equivalent martingale measure. The standard deviations of $k\theta$, k , σ , and \sqrt{V} roughly double. It has a much smaller effect on the estimates of λ_1 and λ_2 . Their standard deviations rise by between 10 and 20 percent. Moreover, the bias in these estimates is unaffected. Doubling the length of the data sample has the same effect with monthly data as with weekly: standard deviations of parameter estimates fall and the bias in the estimates of the price of risk parameters decreases.

We now shift our attention to estimation of the Gaussian model with EMM/SNP. Since this technique is not as efficient at exploiting information in the sample as is maximum likelihood, we expect that the performance of this estimator will not match that of ML. But the actual performance of EMM/SNP is nonetheless surprising.

5.3 EMM/SNP Estimation

Model estimation with EMM/SNP is a two-step procedure. First, an SNP specification is chosen to summarize the features of the data. Second, parameters of the term structure model are chosen to minimize a quadratic form in the score vector from this chosen specification. We follow Gallant and Tauchen by choosing the SNP specification through a sequential search process using the Schwarz Bayes criterion. This search usually resulted in a VAR(2) specification with no higher-order terms for the variance-covariance matrix of innovations. Occasionally the search resulted in a VAR(3) specification. These SNP specifications should be good auxiliary functions because they capture the relevant features of the data. In particular, no higher-order terms for variance-covariance matrix are necessary because the models are Gaussian. Denoting the number of points on the yield curve by m and the lag length as p , there are $m(mp + 1) + m(m + 1)/2$ SNP parameters and therefore an equivalent number of moment conditions.

The right side of Table 2 summarizes the results for estimation of the Vasicek model. The results differ from those produced by ML estimation in two important ways. First, the estimate of σ is strongly biased and estimated with low precision. The mean estimate is about 70 percent of its true value, while its standard deviation is almost four times the standard deviation of the ML estimate. Second, EMM/SNP's estimates of the uncertainty in the parameter estimates are much too small. For example, the sample standard deviations of the estimates of both σ and λ_1 are more than three times the corresponding mean standard errors. Moreover, the χ^2 test of the adequacy of the model, based on the over-identifying moment conditions, rejects the model at the five percent level in more than forty percent

of the simulations. The rejection rate at the one percent level is nearly thirty percent. Put differently, if we use EMM/SNP to judge whether these data are generated by the Vasicek term structure model, we frequently will conclude that they are not.

In Section 8 we discuss in detail the reasons for the poor performance of EMM/SNP. Here we simply preview some of the points we make later. The problem is with EMM, not SNP. Estimation with EMM relies on the asymptotic equivalence between the curvature of the auxiliary function given the sample data and the curvature given an infinite sample of data generated by the true parameters. When the underlying data are highly persistent, these curvatures are often quite different from each other in finite samples. An implication of this divergence is that the weighting matrix used in EMM estimation is the wrong weighting matrix. The use of a wrong weighting matrix results in inefficient parameter estimates and improper statistical inference.

The finite-sample performance of EMM/SNP does not improve when the underlying model is the more general one-factor Gaussian model. The right side of Table 3 summarizes results for estimation of the one-factor model with an affine price of risk. The divergence between EMM/SNP and ML is concentrated in the estimates of the price of risk parameters. The biases in the estimates of λ_1 and λ_2 are about 30 percent larger with EMM/SNP than with ML. The standard deviations of these EMM/SNP estimates are more than twice the corresponding ML standard deviations. They are also more than twice the corresponding mean standard errors. Thus statistical inference is again problematic, although here the χ^2 test of the over-identifying restrictions does not tend to reject the model too often.

In this one-factor Gaussian setting—the simplest possible term structure model—the finite-sample performance of EMM/SNP diverges dramatically from ML. In particular, statistical inference with EMM/SNP is, to put it mildly, problematic, while inference with ML is well-behaved. Given its performance in this simple setting, it is not too difficult to predict how EMM/SNP will perform in estimating the more complicated models that we examine later in the paper.

6 One-factor square-root models

In this section we estimate models in which the instantaneous interest rate follows a non-Gaussian process. The instantaneous interest rate is

$$r_t = \delta_0 + x_t, \tag{21}$$

where the state variable x_t follows a square-root diffusion process under the equivalent martingale measure.

$$dx_t = (k\theta - kx_t)dt + \sigma\sqrt{x_t}d\tilde{z}_t \quad (22)$$

Under the physical measure the dynamics of x_t are

$$dx_t = (k\theta + \lambda_1\sqrt{x_t} - (k - \lambda_2)x_t)dt + \sigma\sqrt{x_t}dz_t \quad (23)$$

This is Duarte’s semiaffine extension of the translated CIR model. The translated CIR model sets λ_1 to zero. The “true” parameters of the process are set to the values reported in the second column of Table 6. The true value of λ_1 is set to zero, although we fit data simulated using the true parameters to both the translated CIR model and to the more general semiaffine extension.

We examine three estimation techniques. The first is maximum likelihood using exact identification of the state vector. For this estimation technique we do not add noise to the ten-year bond yield, so that it can be inverted to infer the state. The three-month and one-year yields are contaminated with noise with a standard deviation of 70 basis points. The second is the Kalman filter and the third is EMM/SNP. For the latter two techniques, all bond yields are observed with measurement error with a standard deviation of 60 basis points. Note that because of the different structures of measurement error, the data used with ML estimation differs from that used with Kalman filter and EMM/SNP estimation.

We make three main points in the following discussion. First, generalizing the price of risk has the same qualitative consequences for finite-sample estimation as it does in the Gaussian model. Second, Kalman filter estimation does not use information from the data as efficiently as does ML estimation. The main consequence is that estimates of parameters identified under the equivalent martingale measure are estimated by the Kalman filter with less precision than they are estimated by ML. Third, EMM/SNP is an unacceptable method for estimating these models.

6.1 ML estimation

We first examine results of estimating the model with the restriction $\lambda_1 = 0$. This restriction allows computation of the probability distribution of discretely observed values of x_t . We therefore implement maximum likelihood using the exact likelihood function. The simulation results are summarized in the first set of columns in Table 6. (Unlike previous tables, this table reports medians as well as means. We refer to the median values when discussing the EMM/SNP results.) Only the estimate of λ_2 is biased, and its bias is not economically

large. The mean estimate of λ_2 , combined with the value of k , implies somewhat faster mean reversion of r_t . The implied half-life of a shock with the mean estimate is 4.5 years, while the true half-life is 5.3 years.

More troubling are the large biases in the estimates of the standard errors of the parameters. Aside from the standard deviation of measurement error, the standard deviations of the parameter estimates are all much larger (on average, twice as large) than the corresponding mean standard errors. Thus with this model, unlike the purely Gaussian model considered in the previous section, finite-sample statistical inference with ML is unreliable.

We next examine results of estimating the more general model (allowing $\lambda_1 \neq 0$) on the same data. Because the probability distribution of discretely observed values of x_t is not known, we use simulated ML in place of exact ML. The parameter space is restricted to ensure stationarity. The restriction is $k - \lambda_2 \geq 0$, with the additional requirement that $\lambda_1 < 0$ if $k - \lambda_2 = 0$.¹² This restriction is occasionally binding. Of the 500 simulations, 13 produced parameter estimates $k - \lambda_2 = 0$, $\lambda_1 < 0$. (Although this is a boundary of the parameter space, the process is strictly stationary.) Results from these 13 simulations are included in the summary statistics for the parameter estimates. They are not included in the summary statistics for standard errors.

A summary of the results is displayed in the first set of columns of Table 7. The main point to take from the table is that the estimates of the price of risk parameters are strongly biased. As with the general Gaussian model, the bias in λ_1 is about one standard deviation and the bias in λ_2 is about minus one standard deviation. Although the implied drift of r_t is nonlinear in r_t , the bias is not really a consequence of nonlinear finite-sample behavior. Instead, it is simply reflecting the same finite-sample bias in the speed of mean reversion that affected estimates of the general Gaussian model discussed in the previous section. The introduction of λ_1 in the model breaks the link between physical and equivalent martingale drifts. Therefore the physical drift is determined exclusively by the time-series properties of the data. Because of the finite-sample bias in the speed of mean reversion, the parameter estimates of the physical drift imply faster mean reversion than is implied by the true parameters.

Figure 2 illustrates the different drifts. The solid lines in both panels display the true drift of r_t as a function of r_t . In Panel A, the dashed line is the drift implied by the mean parameter estimates from exact ML estimation of the restricted model (the mean estimates reported in Table 6). In Panel B, the dashed line is the drift implied by the mean parameter estimates from simulated ML estimation of the more general model. The combination of a

¹²Because we condition on the first observation, stationarity is not necessary. Nonetheless we explicitly impose the stationarity restriction on the estimated parameters.

positive bias in the estimate of λ_1 and a negative bias in the estimate of λ_2 produces a drift function that intersects the x-axis at approximately the same point as the true drift function (thus producing the correct mean of r_t), while implying faster reversion to this mean at all other r_t .

Not all of the features of ML estimation of the more general model compare unfavorably to ML estimation of the restricted model. There is a closer correspondence between standard deviations of the parameter estimates and mean standard errors. Thus statistical inferences concerning parameters identified under the equivalent martingale measure are well-behaved.

6.2 Linearized Kalman filter estimation

First consider the simulation results for Kalman filter estimation that imposes the restriction $\lambda_1 = 0$. We implement the Kalman filter using the correct functional forms for the first and second moments of the discretely observed data and evaluate these functions at the filtered value of the state. The Kalman filter is not ML in this non-Gaussian setting. The simulation results are summarized in the second set of columns in Table 6. The parameter estimates are more biased than the ML estimates, although the differences are not large. For example, the mean Kalman filter estimates imply an unconditional mean interest rate about 80 basis points below the mean ML estimates, which in turn is about 90 basis points below the true unconditional mean. The half-life of an interest rate shock implied by the mean Kalman filter estimates is four years.

The biggest difference between ML and Kalman filter estimation is that the latter's estimates of the equivalent-martingale parameters are much less precise. The extreme cases are δ_0 and $k\theta$. The standard deviations of Kalman filter estimates of these parameters are four times the standard deviations of ML estimates. The precision of parameters not identified under the equivalent martingale measure (λ_2 and \sqrt{V}) are about the same across the two estimation techniques.

Similar patterns are associated with Kalman filter estimation of the more general model, where λ_1 is free. With this model we do not have functional forms for the the discrete-time first and second moments. Therefore the filter is implemented using linearized instantaneous dynamics. For 24 of the 500 simulations, the parameter estimates are on the boundary of the parameter space. They are treated in the same way that such simulations are treated with ML estimation. The estimation results are summarized in the second set of columns in Table 7. The mean parameter estimates produced by the Kalman filter are close to those produced with ML. (The biases in the price of risk parameters are actually slightly less with the Kalman filter than with ML.) Standard deviations of the equivalent-martingale

parameter estimates are much larger than the corresponding standard deviations produced with ML estimation. The precision of estimates of λ_1 , λ_2 , and \sqrt{V} are roughly the same across the two estimation techniques.

On balance, parameter estimates produced with Kalman filter estimation have biases similar to those produced with ML estimation but are less efficient. Thus not surprisingly ML estimation is preferable if it is feasible.¹³ If we are willing to assume exact identification of the state, feasibility effectively depends on computer run time. With the computer resources available to us, a single simulation of this one-factor square-root model (i.e., finding the optimal parameter estimates by maximizing the likelihood function, including an extensive search over the parameter space) takes a few hours. Of course, more complex models require more time; for example, the square-root model with two independent factors discussed in the next section takes about eleven hours to estimate. The current practice in term structure estimation is to use at least three correlated factors. Although such a model could be estimated using a few days of computing time, Monte Carlo analysis of the estimation properties is probably infeasible with current technology. For the models we examine, estimation with the Kalman filter is about 25 to 60 times faster than estimation with simulated ML.

6.3 EMM/SNP estimation

The final sets of results in the two tables summarize EMM/SNP estimation of the models. They do not require a detailed analysis. There are three points to note. First, distributions of parameter estimates and standard errors are strongly skewed. The mean parameter estimates are typically nowhere near the true parameters, while the median estimates are closer. Second, the distributions of parameter estimates have extremely high standard deviations, probably driven by the tail observations. For example, in Table 7, the standard deviations of SNP/EMM estimates are typically about ten times larger than the standard deviation of Kalman filter estimates. Third, tests of the over-identifying restrictions often reject the model. The χ^2 test rejects the model at the five percent level in more than one quarter of the simulations. In a nutshell, EMM/SNP is a failure at estimating the parameters of the one-factor square-root diffusion model.

¹³There is a caveat to this conclusion. The Kalman filter produces standard errors that are closer to the standard deviations of the parameter estimates. Thus inference with the Kalman filter is less prone to false rejections of the true parameters.

7 Estimation of two-factor models

In this section we briefly consider results from estimation of two-factor versions of the Gaussian and square-root diffusion models with generalized prices of risk. We do this primarily to confirm that our conclusions based on one-factor models carry over to two-factor models. We also use a two-factor setting to evaluate the performance of the Kalman filter when the true model exhibits nonlinear dynamics.

7.1 A two-factor Gaussian model with affine risk premia

The model is

$$r_t = \delta_0 + x_{1t} + x_{2t}, \quad (24)$$

where the equivalent martingale dynamics of the state variables are

$$dx_{it} = -k_i x_{it} dt + \sigma_i d\tilde{z}_{it} \quad (25)$$

and the physical dynamics are

$$dx_{it} = (\lambda_{i1} - (k_i - \lambda_{i2})x_{it})dt + \sigma_i dz_{it}. \quad (26)$$

The measurement error in bond yields has a standard deviation of 20 basis points. The true parameters of the process and a summary of the results of Kalman filter (ML) estimation of the model are displayed in Table 8. The table also summarizes the results of estimation with EMM/SNP.

There are two main points to take from these results. The first is that the properties of the Kalman filter estimates are similar to those in the one-factor case. The price of risk parameters are biased and estimated imprecisely. The true speeds of mean reversion of the factors, $k_i - \lambda_{i2}$, are 0.60 and 0.15. The mean estimated speeds of mean reversion are 0.86 and 0.36, and the resulting biases of 0.26 and 0.21 are of the same order of magnitude as the standard deviations of the estimated parameters. To put these biases in a more intuitive perspective, the true half-lives of interest rate shocks are 1.2 years and 4.7 years for the two processes. The half-lives corresponding to the mean parameter estimates are 0.9 years and 1.9 years, respectively. From this perspective, the bias is larger for the more persistent process, as we expect.

The second main point is that EMM/SNP performs much worse than the Kalman filter. The distribution of the parameter estimates of the price of risk are highly skewed. The standard deviations of these estimates are around five to ten times the standard deviations

of the corresponding Kalman filter estimates. The distributions of standard errors are also strongly biased and skewed.

The main conclusion we draw from these results is that the finite-sample estimation properties that we examined extensively in the one-factor Gaussian setting carry over to this two-factor setting. We now consider whether the same conclusion can be drawn for square-root diffusion models.

7.2 A two-factor square-root model with nonlinear risk premia

The model is

$$r_t = \delta_0 + x_{1t} + x_{2t}, \quad (27)$$

where the equivalent martingale dynamics of the state variables are

$$dx_{it} = ((k\theta)_i - k_i)x_{it}dt + \sigma_i\sqrt{x_{it}}d\tilde{z}_{it} \quad (28)$$

and the physical dynamics are

$$dx_{it} = ((k\theta)_i + \lambda_{i1}\sqrt{x_{it}} - (k_i - \lambda_{i2})x_{it})dt + \sigma_i\sqrt{x_{it}}dz_{it}. \quad (29)$$

We estimate the model with simulated ML and with the Kalman filter. The measurement error in bond yields has a standard deviation of 20 basis points. For estimation with simulated ML, we observe the three-month and ten-year yields without error. For estimation with the Kalman filter, all yields are measured with error. Estimation results with EMM/SNP are not reported because of the failure of this technique on the simpler one-factor square-root model.

Unlike the one-factor square-root model we examined, here the true values of λ_{i2} differ from zero. Another difference is that here we do not impose stationarity on the parameter estimates produced with simulated ML. (The true model exhibits stationarity.) Since simulated ML conditions on the first observation, we can estimate the model without imposing stationarity and then take a separate look at the sets of parameter estimates that imply non-stationary rates. Stationarity is imposed on parameter estimates produced with the Kalman filter.

The true parameters of the process and a summary of the estimation results are displayed in Table 9. We first examine the results for ML estimation. Estimates of the parameters identified under the equivalent martingale measure are unbiased, while estimates of the price of risk parameters are strongly biased. As we have seen elsewhere in this paper, the biases are in the direction of faster mean reversion of the states. The drift functions of both state

variables are displayed in Figure 3. The solid lines are the true drift functions.¹⁴ The dashed lines are the drift functions implied by the mean parameter estimates.

Nonstationary term structure behavior is implied by parameter estimates of 71 of the 500 Monte Carlo simulations. These simulations are characterized by negative estimates of both $k_i - \lambda_{i2}$ and λ_{i1} . (In almost all of these simulations x_{1t} is nonstationary and x_{2t} is stationary.) The negative λ_{i1} induces mean-reverting behavior for small x_{it} and the negative $k_i - \lambda_{i2}$ induces mean-averting behavior for large x_{it} . For the purposes of stationarity the latter term dominates, but in any finite sample the former term may dominate. Therefore parameter estimates that imply nonstationarity do not necessarily correspond to mean-averting behavior in the sample.

We now turn to the Kalman filter estimates. Of the 500 Monte Carlo simulations, 126 produced a set of parameter estimates that are on the boundary of the parameter space. As in one-factor square-root model, Kalman filter estimates of equivalent-martingale parameters are not strongly biased, but are much estimated with much less precision than ML estimates. Biases in estimates of the price of risk parameters are similar to the biases in the ML estimates, and the precision of these parameter estimates is close to (but somewhat less than) the precision of the corresponding ML estimates. As with the case of the two-factor Gaussian model, the main conclusion we draw from these results is that the finite-sample estimation properties that we examined extensively in the one-factor square-root setting carry over to this two-factor setting.

8 Interpreting the finite-sample behavior of EMM

One of the more surprising results documented in the previous sections is the poor finite-sample performance of EMM/SNP. Because EMM is a GMM estimator, a natural reaction is to blame the usual GMM suspects: too many moments or poorly chosen moments. In this section we argue that a more subtle effect is at work. The problem is with the matrix used to weight the moments, not with the moments themselves. When the data are highly persistent, the finite-sample variance-covariance matrix of the EMM moments is unlikely to look anything like the finite-sample estimate of this matrix.

For clarity, we illustrate this problem in the starkest possible setting: when the auxiliary likelihood function is the true likelihood function. Thus the moments used by EMM are optimal, in the sense of asymptotic efficiency. To clearly highlight the issues, we do not

¹⁴Note that they do not display much nonlinearity. Recall from Section 4 that the parameters are based on results of fitting the term structure model to Treasury data.

examine a term structure model here. Instead, we examine a Gaussian AR(1) process,

$$y_t = \alpha_0 + \beta_0 y_{t-1} + \epsilon_t, \quad \epsilon_t \sim N(0, \nu_0).$$

Stack the parameters into the vector $\rho_0 = (\alpha_0 \ \beta_0 \ \nu_0)'$. We can estimate ρ_0 given a time series of data y_0, \dots, y_T . We first discuss some features of ML estimation, then turn to EMM estimation.

8.1 ML estimation of an AR(1)

Denote the true log-likelihood function conditional on an unknown parameter vector ρ as $f(y_t|y_{t-1}; \rho)$. Its derivative with respect to ρ is

$$h(y_t|y_{t-1}; \rho) \equiv \frac{\partial f(y_t|y_{t-1}; \rho)}{\partial \rho} = \nu^{-1} \begin{pmatrix} e_t \\ e_t y_{t-1} \\ -(1/2)(1 - e_t^2/\nu) \end{pmatrix},$$

where the fitted residual is

$$e_t = y_t - \alpha - \beta y_{t-1}.$$

Denote the ML parameter estimate (conditioned on observation y_0) as ρ_T . This estimate sets the mean derivative equal to zero:

$$h_T(\rho_T, Y_T) = 0, \quad h_T(\rho, Y_T) \equiv \frac{1}{T} \sum_{t=1}^T h(y_t|y_{t-1}; \rho).$$

We can think of ML as a GMM estimator where the moment vector is $h(y_t|y_{t-1}; \rho)$. The estimates of α and β are the usual ordinary least-squares parameter estimates and the estimate of ν is the mean of the squared fitted residuals. This parameter vector is biased but consistent, the bias arising because the sample means of ϵ_t and y_{t-1} are correlated.

The asymptotic variance of the moment vector h evaluated at ρ_0 is S/T , where

$$S = \nu^{-1} \begin{pmatrix} 1 & E(y) & 0 \\ E(y) & E(y^2) & 0 \\ 0 & 0 & 1/(2\nu) \end{pmatrix}.$$

Here, $E(y) = \alpha_0/(1 - \beta_0)$ and $E(y^2) = E(y)^2 + \nu_0/(1 - \beta_0^2)$. An estimate of S can be

calculated from the outer product of h :

$$S_T = \frac{1}{T} \sum_{t=1}^T h(y_t|y_{t-1}; \rho_T) h(y_t|y_{t-1}; \rho_T)'$$

One way to judge the accuracy of this variance estimate is to use it to test the hypothesis that ρ_0 is the true parameter vector. (Although the moment vector is identically zero when evaluated at the ML estimate, it is a random variable when evaluated at some other parameter vector such as ρ_0 .) If the finite-sample variance of $h_T(\rho_0, Y_T)$ is close to S_T , then the following likelihood ratio test statistic should have a distribution close to a $\chi^2(3)$:

$$LR = T h_T(\rho_0, Y_T)' S_T^{-1} h_T(\rho_0, Y_T). \quad (30)$$

The difference between the finite-sample and asymptotic distributions of this statistic is a common metric for evaluating GMM estimation techniques. Below we use Monte Carlo simulations to compute the true distribution of this statistic.

8.2 EMM estimation of an AR(1)

Now consider estimating the same model with EMM, using the true likelihood function as the auxiliary model. Because EMM requires two sets of parameter estimates, denote the parameter vector for the auxiliary model as $p = (a \ b \ v)'$. The ML estimate of this vector is denoted p_T . The model's simplicity allows us to express the EMM moment vector analytically. It is

$$m_T(\rho, p_T) = v_T^{-1} \begin{pmatrix} \hat{a}_T + \hat{b}_T E(y; \rho) \\ \hat{a}_T E(y; \rho) + \hat{b}_T E(y^2; \rho) \\ -(1/2) \left(1 - v_T^{-1} (\hat{a}_T^2 + \hat{b}_T^2 E(y^2; \rho) + \nu + 2\hat{a}_T \hat{b}_T E(y; \rho)) \right) \end{pmatrix}, \quad (31)$$

where

$$\begin{aligned} \hat{a}_T &\equiv \alpha - a_T, & \hat{b}_T &\equiv \beta - b_T, \\ E(y; \rho) &\equiv \alpha / (1 - \beta), & E(y^2; \rho) &= E(y; \rho)^2 + \nu / (1 - \beta^2). \end{aligned}$$

It is well-known that when the true log-likelihood function is used as the auxiliary model, the EMM parameter estimates are identical to ML estimates. This is easily seen in (31). At $\rho = p_T$, this moment vector is identically zero.

Recall the general formula for the variance of the EMM moment vector in (7). Although $C(\rho_0)$ and d cancel each other asymptotically, it is helpful to write out finite-sample estimates

of $C(\rho)$ and the auxiliary model's Hessian d . The estimate of $C(\rho)$ is

$$C(\rho) = \left(\frac{\partial m_T(\rho, \gamma_T)}{\partial \gamma'} \Big|_{\gamma=\gamma_T} \right).$$

The explicit representation of this matrix is

$$C(\rho) = -v_T^{-1} \begin{pmatrix} 1 & E(y; \rho) & \frac{\hat{a}_T + \hat{b}_T E(y; \rho)}{v_T} \\ E(y; \rho) & E(y^2; \rho) & \frac{\hat{a}_T E(y; \rho) + \hat{b}_T E(y^2; \rho)}{v_T} \\ \frac{\hat{a}_T + \hat{b}_T E(y; \rho)}{v_T} & \frac{\hat{a}_T E(y; \rho) + \hat{b}_T E(y^2; \rho)}{v_T} & \frac{2v_T^{-1}(\hat{a}_T^2 + \hat{b}_T^2 E(y^2; \rho) + 2\hat{a}_T \hat{b}_T E(y; \rho) + \nu) - 1}{2v_T} \end{pmatrix}.$$

This estimate is the curvature of the auxiliary function evaluated at p_T , using an infinite sample generated by the parameter vector ρ . The estimate of d is

$$d_T = -\nu_T^{-1} \begin{pmatrix} 1 & \overline{y_{t-1}} & 0 \\ \overline{y_{t-1}} & \overline{y_{t-1}^2} & 0 \\ 0 & 0 & 1/(2\nu_T) \end{pmatrix}.$$

This estimate is the curvature of the auxiliary function evaluated at p_T , using the sample data. Note that if we evaluate $C(\rho)$ at $\rho = p_T$, it is (almost) identical to d_T .¹⁵ Asymptotically both matrices converge in absolute value to S , as does the variance of m_T . This asymptotic behavior motivates the use of S_T as an estimate of the variance of m_T .

As with ML, we can judge the accuracy of this estimate by using it to test the hypothesis that ρ_0 is the true parameter vector. If the finite-sample variance of $m_T(\rho_0, p_T)$ is close to S_T , then the following J statistic should have a distribution close to a $\chi^2(3)$.

$$J = T m_T(\rho_0; p_T)' S_T^{-1} m_T(\rho_0; p_T). \quad (32)$$

Below we use Monte Carlo simulations to compute the true distributions of these statistics.

Even if S_T is a good approximation to the finite-sample variance of $h_T(\rho_0, Y_T)$, it will not be a good approximation to the variance of $m_T(\rho_0, p_T)$ unless $C(\rho_0)$ is close to d_T . The greater the difference between ρ_0 and p_T , the greater the difference between these two curvatures. Thus it is instructive to consider an alternative statistic.

$$J_{alt} = T m_T(\rho_0; p_T)' C(\rho_0)^{-1} d_T S_T^{-1} d_T C(\rho_0)^{-1} m_T(\rho_0; p_T). \quad (33)$$

We refer to (33) as a ‘‘robust EMM’’ test statistic. It is not a statistic usually associated

¹⁵They are not identical because the sample first and second moments of y_t are not exactly equal to the sample first and second moments of y_{t-1} .

with EMM because it relies on the ability to compute $C(\rho)$. We can compute it analytically in this simple case, but when the true likelihood function is unknown or intractable this matrix is also either unknown or intractable.

8.3 Monte Carlo Results

We use Monte Carlo simulations to study the finite-sample properties of these estimation techniques. We set $\alpha_0 = 0$ and $\nu_0 = 1$, and focus on values of β_0 that are very close to one. Given ρ_0 , we use the AR model to simulate a set of data Y_T . We then calculate the ML estimate ρ_T , the estimated asymptotic variance of the moments S_T , the GMM moment vector $h_T(\rho_0, Y_T)$, and the EMM moment vector $m_T(\rho_0, p_T)$. We also calculate the tests of the true model (30), (32), and (33).

We summarize the behavior of the simulated ML and EMM moment vectors with their sample variances across the Monte Carlo simulations. We then compare these empirical variances to the mean, across the simulations, of S_T . The results are displayed in Table 10. To simplify the interpretation of the results, the empirical variances of the individual elements of these moment vectors are divided by the mean of the corresponding diagonal elements of S_T . To conserve space the table does not report the ratio of the variance of the EMM moment vector to the more robust estimate of its variance, $C(\rho_0)d_T^{-1}S_Td_T^{-1}C(\rho_0)$. We discuss this ratio in the text.

The first two rows of the table consider the case of $\beta_0 = 0.99875$, which is the AR(1) coefficient for weekly observations of the instantaneous interest rate from a Vasicek model with a speed of mean reversion of 0.065. With 1000 observations (the first row), the mean estimate of β is biased down. If the data are weekly, the annual speed of mean reversion implied by this mean coefficient is 0.11.

Although the estimate of β is biased, the finite-sample variance of the ML moment is close to its estimated asymptotic value. The ratios of individual variances lie between 0.98 and 1.1. By contrast, the finite-sample variance of the EMM moment vector differ dramatically from its estimated asymptotic value. The ratios of individual variances range from 0.01 to 12. The main reason for this large divergence between finite-sample and estimated asymptotic behavior is that the curvature of the likelihood function in the neighborhood of truth differs from the curvature in the neighborhood of the ML estimates. The robust estimate of the variance of the EMM moment vector corrects for this difference in curvature. With this correction, the ratios of the empirical variances of the elements of the EMM moment vector to the corresponding diagonal elements of this robust variance estimate (not reported in the table) range from 1.12 to 2.20.

Given these results concerning the accuracy of the asymptotic variance estimates, it is not surprising that the EMM test (32) of the hypothesis $\rho = \rho_0$ is rejected at the five percent asymptotic level in roughly three-quarters of the simulations. The ML test also over-rejects truth, but the empirical rejection rate at the five percent level is a more modest eight percent. The robust EMM test rejects the hypothesis $\rho = \rho_0$ in twelve percent of the simulations. Doubling the amount of data reduces the difference between the ML and EMM results, but as the second row of the table shows, the behavior of the EMM moment vector still differs substantially from its asymptotic behavior.

The third and fourth rows of the table reports results for a less persistent process: $\beta_0 = 0.95$. For weekly data, this corresponds to speed of mean reversion of 2.67. The lower persistence produces finite-sample behavior of the EMM moment vector that is much closer to its asymptotic behavior. For example, with 1000 observations, the ratios of true variances of the three EMM moments to their estimated asymptotic variances range from 0.96 to 1.13. Nonetheless, with 1000 observations the true model is over-rejected; the rejection rate is thirteen percent using the five percent asymptotic test statistic.

Zhou (2001) also shows that EMM over-rejects the true model when the data exhibit high persistence. He simulates a square-root diffusion model of the instantaneous interest rate and calls the empirical rejection rate with EMM/SNP “unacceptably large.” Our contribution is to explain the source of this behavior. It is a consequence of the fact that when the data are highly persistent, the curvature of the log-likelihood function evaluated using the sample data diverges from the curvature evaluated using the true parameters. More generally, convergence of the variability of the EMM moment vector to the variability of the auxiliary score vector requires longer samples of data than are usually available to the econometrician.

In practice, we are seldom confronted with a problem in which the divergence between the curvature of f using the sample data and the curvature using data generated by the true model is relevant: we usually do not know the true model. But the main message here carries over to a much broader class of problems. Anytime there is a divergence between the properties of the sample data and the simulated data, there will also be a divergence in the curvature of the log-likelihood function. This situation arises whenever the number of moment conditions exceeds the number of free parameters. In this case, the moments cannot be set exactly to zero, and instead a quadratic form in those moments is minimized (see Equation (8)). The moments are weighted by the inverse of their estimated asymptotic variance. If this matrix differs substantially from the true variance of the EMM moment vector, it will result in both inefficient parameter estimates and incorrect statistical inference.

As the results in this paper have shown, this is a particular problem in a term structure setting when SNP is used as the auxiliary model. When panel data are used to estimate the

model, the VAR specification of SNP results in many over-identifying restrictions. Of course, the problem is with EMM, not SNP. The use of SNP simply starkly reveals the problems with EMM. In this section we illustrate the inefficiency and incorrect inference in the context of the AR(1) model. Consider estimation of α in the AR(1) model when we are given the true values β_0 and ν_0 . Starting with the unrestricted model, a natural way to extend the ML estimation procedure is to use GMM to solve

$$\operatorname{argmin}_{\alpha} J = T h_T((\alpha, \beta_0, \nu_0)', Y_T)' S_T^{-1} h_T((\alpha, \beta_0, \nu_0)', Y_T) \quad (34)$$

Similarly, we can estimate α using EMM by solving

$$\operatorname{argmin}_{\alpha} J = T m_T((\alpha, \beta_0, \nu_0)', p_T)' S_T^{-1} m_T((\alpha, \beta_0, \nu_0)', p_T) \quad (35)$$

The latter problem is equivalent to estimating the restricted model with EMM using SNP as the auxiliary model, where the SNP specification is a Gaussian AR(1). The solutions to these minimization problems will differ. In our Monte Carlo simulations we also calculated these two estimates of α . In the final two columns of Table 10 we report the standard deviations, across the simulations, of the estimates. (We do not report the estimated means in the table. In every case they are close to α_0 .)

Consider the first row in Table 10. When the data are highly persistent, with 1000 observations the standard deviation of the estimates based on the EMM solution (35) is more than twice the standard deviation of the estimates based on (34). In addition, with EMM, 40 percent of the simulations produce J statistics that imply rejection of the model at the 5 percent level. The model is also rejected too frequently when it is estimated with (34), but the over-rejection rate is quite modest. The inefficiency and poor inference associated with EMM remains when the amount of data is doubled (the second row). With less persistent data (the third and fourth rows), the relative efficiency of EMM improves substantially.

The main message to draw from this extended discussion is that the finite-sample behavior of the EMM moment vector has little in common with its estimated asymptotic behavior unless the distribution of the data implied by the EMM parameter estimates is close to the distribution of the data sample. The distributions need to be sufficiently close so that the curvature of the auxiliary function is similar across the two distributions. If the true model is sufficiently tractable, we can adjust for differences in curvature to produce more reliable estimates. However, the usual application of EMM is to estimate models that are intractable. Thus as a practical matter there appears to be little we can do to shore up the performance of EMM.

9 Conclusions

The analysis in this paper leads to several important conclusions. First, despite its well-known asymptotic optimality properties, maximum likelihood yields highly biased parameter estimates in settings representative of those used in the estimation of modern term structure models. This differs from the results of Ball and Torous (1996). They find that when cross-sectional information is used in estimation, the bias in the estimated speed of mean reversion and price of risk disappears. However, they only consider very simple forms for the market price of interest rate risk. Their conclusion does not hold when we allow the price of risk to be more flexible, such as the forms specified by Duarte (2004) and Duffee (2002). Once the physical and equivalent-martingale drifts are decoupled, the use of cross-sectional information can no longer eliminate the time-series related bias in the estimated speed of mean reversion.

Our second conclusion is that, despite its asymptotic equivalence to maximum likelihood, the Efficient Method of Moments is an unacceptable alternative to maximum likelihood in finite samples. Most surprisingly, this is true even in settings where maximum likelihood performs well. Much larger samples are required for the small-sample performance of EMM to approach its asymptotic behavior than is the case for ML.

Our third conclusion is that the Kalman filter is a reasonable alternative to maximum likelihood, even in non-Gaussian settings where the two are not equivalent. Unsurprisingly, when feasible, ML usually makes more efficient use of the information in the sample than does the Kalman filter, but the Kalman filter is vastly superior to EMM.

These results underscore the importance of performing detailed Monte Carlo analysis to study the small-sample properties of new estimators, rather than merely relying on their asymptotic properties. As we have shown in this paper, even techniques that are asymptotically equivalent can have very different properties when used on finite samples.

A SNP as the Auxiliary Model for EMM

The standard auxiliary model for use with EMM has become the SNP (for **SemiNonParametric**) model of Gallant and Tauchen (1992) (see also the user's guide, Gallant and Tauchen, 1998). It consists of writing the conditional density of the dataset under analysis in the form of a Hermite polynomial multiplied by a normal density, i.e.,

$$f(y_t | x_{t-1}, \rho) = c(x_{t-1}) [h(z_t | x_{t-1})]^2 \phi(z_t),$$

where

- $\phi(\cdot)$ represents the standard normal p.d.f.,
- $h(z_t | x_{t-1})$ is a Hermite polynomial in z_t ,
- $c(x_{t-1})$ is a normalization constant (equal to $1 / \int [h(s | x_{t-1})]^2 \phi(s) ds$), and
- z_t is a normalized version of y_t , defined by

$$z_t = R_{x,t-1}^{-1} (y_t - \mu_{x,t-1}),$$

where $\mu_{x,t-1}$ is the conditional mean, and $R_{x,t-1}^{-1}$ the Cholesky decomposition of the conditional variance of y_t . This specification allows great flexibility in fitting the conditional distribution. In particular, we are free to choose:

- the dimensionality of the Hermite polynomial in z , K_z (allows for non-Gaussian behavior).
- the degree of the polynomial in x that makes up each of the coefficients in the Hermite polynomial, K_x (another way to allow for conditional heterogeneity).
- the number of lags of x in the Hermite polynomial, K_p .
- the number of lags in a VAR specification for $\mu_{x,t-1}$, L_μ .
- the degree of an ARCH (or GARCH, setting $L_g > 0$) specification for the scale transformation $R_{x,t-1}$, L_r .

Choice of an appropriate specification is performed by using a model selection criterion, such as the Schwarz Bayes information criterion (see Schwarz, 1978), which rewards good fit while penalizing over-parametrization. Gallant and Tauchen (1998) discuss a search strategy for finding an appropriate parametrization for a given problem.

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	True Value	Mean estimate	Std dev	Mean std error
$(k\theta + \lambda_1) \times 10^3$	3.4	14.5	18.3	9.8
$k \times 10^2$	6.5	30.4	23.9	16.7
$\sigma \times 10^3$	17.5	17.5	0.4	0.4

Table 1: Maximum likelihood parameter estimates of a Gaussian model of the short rate.

The interest rate model is $dr = (k\theta + \lambda_1 - kr_t)dt + \sigma dz$. This table summarizes the results of 500 Monte Carlo simulations. Each simulation consists of 1000 weekly observations of the instantaneous interest rate. The parameters are estimated with maximum likelihood.

Parameter	True Value	Kalman filter (ML)			— EMM/SNP —		
		Mean	Std dev	Mean std err	Mean	Std dev	Mean std err
$k\theta \times 10^3$	8.4	8.4	0.24	0.23	8.0	0.34	0.20
$k \times 10^2$	6.5	6.5	0.31	0.31	6.5	0.31	0.25
$\sigma \times 10^3$	17.5	17.5	1.02	0.99	12.2	3.76	1.23
$\lambda_1 \times 10^3$	-5.0	-5.0	2.33	3.35	-4.2	2.55	0.66
$\sqrt{V} \times 10^3$	6.0	6.0	0.09	0.09	6.0	0.11	0.08

Table 2: Parameter estimates of a one-factor Gaussian model with constant price of risk

The model is given by equations (19) and (20) in the text, with $\lambda_2 = 0$. This table summarizes the results of 500 Monte Carlo simulations. With each simulation, 1000 weeks of yields on bonds with maturities of three months, one year, and ten years are observed with iid measurement error. The standard deviation of the error is \sqrt{V} . The parameters of the model are estimated with the Kalman filter (which is ML) and with EMM/SNP. The mean number of overidentifying moments with EMM/SNP is 23.6. The EMM χ^2 test rejects the model at the 5% level in 43.2% of the simulations and rejects at the 1% level in 28.6% of the simulations.

Parameter	True Value	- Kalman filter (ML) -			- EMM/SNP -		
		Mean	Std dev	Mean std err	Mean	Std dev	Mean std err
$k\theta \times 10^3$	8.4	8.4	0.31	0.30	8.4	0.47	0.37
$k \times 10^2$	6.5	6.5	0.39	0.37	6.5	0.60	0.45
$\sigma \times 10^3$	17.5	17.7	1.04	1.02	12.2	1.58	1.57
$\lambda_1 \times 10^3$	5.0	20.0	16.98	14.12	25.0	46.78	17.33
$\lambda_2 \times 10^2$	-14.0	-37.4	24.14	20.96	-43.7	53.37	25.06
$\sqrt{V} \times 10^3$	6.0	6.0	0.09	0.09	6.0	0.10	0.09

Table 3: Parameter estimates of a one-factor Gaussian model with affine price of risk

The model is given by equations (19) and (20) in the text. This table summarizes the results of 500 Monte Carlo simulations. With each simulation, 1000 weeks of yields on bonds with maturities of three months, one year, and ten years are observed with iid measurement error. The standard deviation of the error is \sqrt{V} . The parameters of the model are estimated with the Kalman filter (which is ML) and with EMM/SNP. The mean number of overidentifying moments with EMM/SNP is 22.5. The EMM χ^2 test rejects the model at the 5% level in 5.6% of the simulations and rejects at the 1% level in 1.4% of the simulations.

Parameter	True Value	5 bonds			2 bonds			10 b.p. measurement error		
		Mean	Std dev	Mean std err	Mean	Std dev	Mean std err	Mean	Std dev	Mean std err
$k\theta \times 10^3$	8.4	8.4	0.27	0.26	8.4	0.34	0.32	8.4	0.06	0.06
$k \times 10^2$	6.5	6.5	0.34	0.33	6.5	0.41	0.39	6.5	0.06	0.06
$\sigma \times 10^3$	17.5	17.6	0.83	0.83	17.6	1.22	1.19	17.5	0.42	0.43
$\lambda_1 \times 10^3$	5.0	19.8	16.74	13.95	20.0	16.92	14.22	19.5	16.24	13.73
$\lambda_2 \times 10^2$	-14.0	-37.0	23.50	20.67	-37.3	24.13	21.12	-36.7	23.04	20.29
$\sqrt{V} \times 10^3$	Varies	6.0	0.07	0.06	6.0	0.11	0.11	1.0	0.02	0.02

Table 4: Maximum likelihood parameter estimates of a one-factor Gaussian model with affine price of risk: varying assumptions

The model is given by equations (19) and (20) in the text. This table summarizes the results of three sets of 500 Monte Carlo simulations. With each simulation, 1000 weeks of yields on bonds with maturities of ranging from three months to ten years are observed with iid measurement error. The standard deviation of the error is \sqrt{V} , set to 60 basis points in the first two sets of simulations and 10 basis points in the third set. The third set also assumes that three bond yields are observed.

Panel A. Weekly observations

Parameter	True Value	500 obs			2000 obs		
		Mean	Std dev	Mean std err	Mean	Std dev	Mean std err
$k\theta \times 10^3$	8.4	8.5	0.57	0.53	8.4	0.19	0.19
$k \times 10^2$	6.5	6.5	0.71	0.66	6.5	0.23	0.22
$\sigma \times 10^3$	17.5	17.7	1.37	1.48	17.6	0.70	0.71
$\lambda_1 \times 10^3$	5.0	36.9	34.99	25.57	12.4	10.27	8.70
$\lambda_2 \times 10^2$	-14.0	-62.7	48.44	38.03	-25.6	15.10	12.70
$\sqrt{V} \times 10^3$	6.0	6.0	0.12	0.12	6.0	0.06	0.06

Panel B. Monthly observations

Parameter	True Value	240 obs			480 obs		
		Mean	Std dev	Mean std err	Mean	Std dev	Mean std err
$k\theta \times 10^3$	8.4	8.4	0.67	0.61	8.4	0.37	0.37
$k \times 10^2$	6.5	6.5	0.82	0.77	6.5	0.47	0.46
$\sigma \times 10^3$	17.5	17.7	1.52	1.50	17.6	0.99	1.03
$\lambda_1 \times 10^3$	5.0	20.2	20.25	14.63	12.2	9.95	8.71
$\lambda_2 \times 10^2$	-14.0	-37.3	26.94	21.75	-25.2	13.95	12.83
$\sqrt{V} \times 10^3$	6.0	6.0	0.20	0.19	6.0	0.13	0.13

Table 5: Maximum likelihood parameter estimates of a one-factor Gaussian model with affine price of risk: varying assumptions

The model is given by equations (19) and (20) in the text. This table summarizes the results of four sets of 500 Monte Carlo simulations. With each simulation, yields on bonds with maturities of ranging from three months to ten years are observed with iid measurement error. The standard deviation of the error is \sqrt{V} .

Parameter	True Value	Exact ML			Kalman filter			EMM/SNP								
		Mean	Median	Std dev	Mean	Median	Std dev	Mean	Median	Std dev						
$\delta_0 \times 10^2$	1.0	1.0	1.0	0.67	0.28	0.23	0.6	1.2	2.83	1.91	0.94	-32.8	2.7	190.70	265.75	1.76
$k\theta \times 10^3$	7.5	7.5	7.5	0.56	0.27	0.23	7.8	7.4	2.27	1.59	0.79	34.2	5.8	154.12	209.51	1.46
$k \times 10^2$	6.3	6.3	6.3	0.26	0.21	0.20	6.3	6.2	0.67	0.62	0.59	4.2	4.7	3.08	2.73	2.07
$\sigma \times 10^2$	8.0	8.0	8.1	0.44	0.28	0.28	8.1	8.1	1.23	1.06	0.97	9.6	10.1	5.07	3.83	3.57
$\lambda_2 \times 10^2$	-6.8	-9.2	-7.6	7.04	2.23	1.77	-11.2	-9.5	8.71	7.52	7.06	-19.4	-13.5	22.42	25.86	9.05
$\sqrt{V} \times 10^3$	Varies	7.0	7.0	0.11	0.11	0.11	6.0	6.0	0.09	0.09	0.09	5.9	5.9	0.32	0.09	0.09

Table 6: Parameter estimates of a one-factor square-root model with linear risk premium

The model is given by equations (21) through (23) in the text, with $\lambda_1 = 0$. This table summarizes the results of 500 Monte Carlo simulations. With each simulation, 1000 weeks of yields on three bonds with maturities ranging from three months, one year, and ten years are observed. For “Exact ML”, the ten-year yield is observed without error and the three-month and one-year yields are observed with normally distributed, iid measurement error. The standard deviation of the error, \sqrt{V} , is 70 basis points. For Kalman filter estimation and EMM/SNP estimation, all three yields are observed with measurement error. The standard deviation of the error is 60 basis points. The mean number of overidentifying restrictions with EMM/SNP is 23.0. The EMM χ^2 test rejects the model at the five percent level in 26.4 percent of the simulations and at the one percent level in 17.4 percent of the simulations.

Parameter	True Value	Simulated ML			Linearized Kalman filter			EMM/SNP								
		Mean	Median	Std dev	Mean	Median	Std dev	Mean	Median	Std dev						
$\delta_0 \times 10^2$	1.0	1.0	1.0	0.61	0.49	0.34	0.5	1.1	3.12	2.24	1.00	-12.7	1.2	33.55	32.22	14.22
$k\theta \times 10^3$	7.5	7.5	7.5	0.51	0.44	0.33	7.9	7.5	2.50	1.85	0.87	18.4	7.4	27.30	25.11	11.52
$k \times 10^2$	6.3	6.3	6.3	0.27	0.28	0.26	6.3	6.2	0.68	0.63	0.60	5.5	6.3	2.65	3.53	3.44
$\sigma \times 10^2$	8.0	8.0	8.1	0.44	0.42	0.39	8.0	8.0	1.27	1.12	1.02	8.2	7.9	4.30	7.46	6.55
$\lambda_1 \times 10^2$	0.0	11.1	8.5	11.05	9.02	8.40	10.3	7.2	13.20	9.91	8.35	30.7	9.4	145.77	75.17	20.54
$\lambda_2 \times 10^2$	-6.8	-56.7	-46.3	46.81	39.64	37.99	-48.7	-37.9	45.07	37.93	35.91	-91.2	-41.9	435.65	79.09	40.65
$\sqrt{V} \times 10^3$	Varies	7.0	7.0	0.11	0.11	0.11	6.0	6.0	0.09	0.09	0.09	5.9	5.9	0.10	0.09	0.09

Table 7: Parameters estimates of a one-factor square-root model that allows for a nonlinear risk premium

The model is given by equations (21) through (23) in the text. This table summarizes results from 500 Monte Carlo simulations. With each simulation, 1000 weeks of yields of bonds with maturities of three months, one year, and ten years are observed. The true model has a linear risk premium, while the estimated models allow for a nonlinear risk premium. For “Simulated ML”, the ten-year yield is observed without error and the three-month and one-year yields are observed with normally distributed, iid measurement error. The standard deviation of the error, \sqrt{V} , is 70 basis points. For linearized Kalman filter estimation and EMM/SNP estimation, all three yields are observed with measurement error with a standard deviation of 60 basis points. The mean number of overidentifying restrictions with SNP/EMM is 22. The EMM χ^2 test rejects the model at the 5% level in 9.4 percent of the simulations and at the 1% level in 4.0 percent of the simulations.

Parameter	True Value	Kalman filter (exact ML)				EMM/SNP					
		Mean	Median	Std dev	Mean std err	Mean	Median	Std dev	Mean std err		
$\delta_0 \times 10$	1.0	1.1	1.0	1.05	0.69	0.08	1.6	1.1	5.04	$> 10^3$	0.12
$k_1 \times 10^2$	70.0	70.4	70.4	3.09	3.04	2.98	70.4	70.0	5.19	3.94	3.48
$k_2 \times 10^2$	2.7	2.7	2.7	0.64	0.63	0.63	2.7	2.6	1.44	1.14	0.92
$\sigma_1 \times 10^2$	2.0	2.0	2.0	0.09	0.09	0.09	2.0	2.0	0.13	0.11	0.10
$\sigma_2 \times 10^2$	1.3	1.3	1.3	0.07	0.07	0.07	1.3	1.3	0.11	0.09	0.08
$\lambda_{11} \times 10^2$	-1.0	-1.4	-1.3	0.84	0.70	0.66	-2.2	-1.8	3.89	1.01	0.70
$\lambda_{21} \times 10^2$	-0.4	-1.4	-0.8	7.84	5.15	0.62	-14.8	-1.19	68.11	91.12	1.06
$\lambda_{12} \times 10^2$	10.0	-15.2	-8.8	34.68	30.34	30.04	-61.5	-37.7	149.06	45.45	31.25
$\lambda_{22} \times 10^2$	-12.0	-33.6	-27.9	23.57	18.96	18.27	-120.6	-50.6	267.34	92.04	25.09
$\sqrt{V} \times 10^3$	2.0	2.0	2.0	0.04	0.04	0.04	2.0	2.0	0.04	0.03	0.03

Table 8: Parameter estimates of a two-factor Gaussian model with affine price of risk

The model is given by equations (24) through (26) in the text. This table summarizes the results of 500 Monte Carlo simulations. With each simulation, 1000 weeks of yields on three bonds with maturities of three months, one year, and ten years are observed with iid measurement error. The standard deviation of the error is \sqrt{V} . The parameters of the model are estimated with the Kalman filter (which is ML) and with EMM/SNP. The mean number of overidentifying restrictions with SNP/EMM is 23.1. The EMM χ^2 test rejects the model at the 5% level in 7.4 percent of the simulations and at the 1% level in 3.2 percent of the simulations.

Parameter	True Value	— Simulated ML —			Linearized Kalman filter		
		Mean	Std dev	Mean std err	Mean	Std dev	Mean std err
$\delta_0 \times 10^2$	-3.6	-3.7	0.93	0.88	-5.3	9.72	15.96
$k\theta_1 \times 10^2$	4.0	4.0	0.26	0.28	4.4	2.40	2.02
$k\theta_2 \times 10^2$	0.3	0.3	0.04	0.04	0.4	0.42	0.52
$k_1 \times 10^2$	70.0	69.9	2.18	2.25	70.7	3.67	3.77
$k_2 \times 10^2$	3.0	3.0	0.40	0.41	3.0	0.93	0.87
$\sigma_1 \times 10^2$	8.9	9.0	0.49	0.54	8.8	1.41	1.28
$\sigma_2 \times 10^2$	5.2	5.2	0.32	0.36	5.1	0.97	0.91
$\lambda_{11} \times 10^2$	-20.0	-14.3	10.65	11.44	-9.1	17.64	16.17
$\lambda_{21} \times 10^2$	0.5	8.1	10.98	7.63	9.8	14.05	15.51
$\lambda_{12} \times 10^2$	66.0	35.4	56.19	59.11	13.5	65.86	65.61
$\lambda_{22} \times 10^2$	-5.0	-40.6	47.28	36.16	-46.5	48.13	35.62
$\sqrt{V} \times 10^3$	2.0	2.0	0.04	0.05	2.0	0.03	0.03

Table 9: Parameter estimates of a two-factor square-root model with nonlinear risk premium

The model is given by equations (27) through (29) in the text. This table summarizes the results of 500 Monte Carlo simulations. With each simulation, 1000 weeks of yields on three bonds with maturities of three months, one year, and ten years are observed. For “Simulated ML”, the three-month and ten-year yields are observed without error and the one-year yield is observed with normally distributed, iid measurement error. The standard deviation of the error, \sqrt{V} , is 20 basis points.

Num obs (T)	β_0	Mean				Ratio of true var to estimated asymptotic var				Empirical 5% rejection freq			Std dev of restricted estimates of α	
		β_T	α	β	ν	α	β	ν	ML	EMM	Robust EMM	ML	EMM	
1000	0.99875	0.99794	0.977	1.095	1.013	12.192	0.012	1.901	0.078	0.732	0.116	0.032	0.065	
												[.074]	[.400]	
2000	0.99875	0.99829	1.037	1.077	1.009	5.449	0.012	1.388	0.070	0.454	0.072	0.023	0.034	
												[.068]	[.226]	
1000	0.95000	0.94702	0.977	1.021	1.014	1.133	0.960	1.076	0.052	0.130	0.060	0.032	0.030	
												[.058]	[.116]	
2000	0.95000	0.94838	1.037	1.011	1.009	1.103	0.917	1.036	0.052	0.074	0.054	0.023	0.023	
												[.060]	[.066]	

Table 10: Results of Monte Carlo simulation of ML and EMM estimation of an AR(1) model

This table reports results from 500 Monte Carlo simulations. The simulated model is

$$x_t = \alpha_0 + \beta_0 x_{t-1} + \epsilon_t, \quad \epsilon_t \sim N(0, \nu_0), \quad \alpha_0 = 0, \nu_0 = 1$$

For each simulation a sample of length T is generated and the parameters are estimated with both ML (using the score vector of the likelihood function as the moment vector in GMM estimation) and with EMM (using the true likelihood function as the auxiliary likelihood function). The asymptotic variance of the moment vectors is also estimated, and the ML and EMM moment vectors are calculated at the model's true parameters. The table reports the variances, across the simulations, of these moments divided by the corresponding means (across the simulations) of the estimated asymptotic variances of these moments. For example, a value of 10 for α implies that the empirical variance of the partial derivative of the likelihood function with respect to α is 10 times the mean estimated asymptotic variance. The same moment vectors and estimated asymptotic variances are used to construct LR tests of the hypothesis that the model's parameters equal the true parameters. The table reports the frequency with which the hypothesis is rejected at the 5% level. In addition, a restricted version of the model is estimated where β_0 and ν_0 are known and α is unknown. Standard deviations of the estimates of α are displayed in the final two columns along with the frequency with which the hypothesis that $\alpha = \alpha_0$ is rejected at the 5% level.

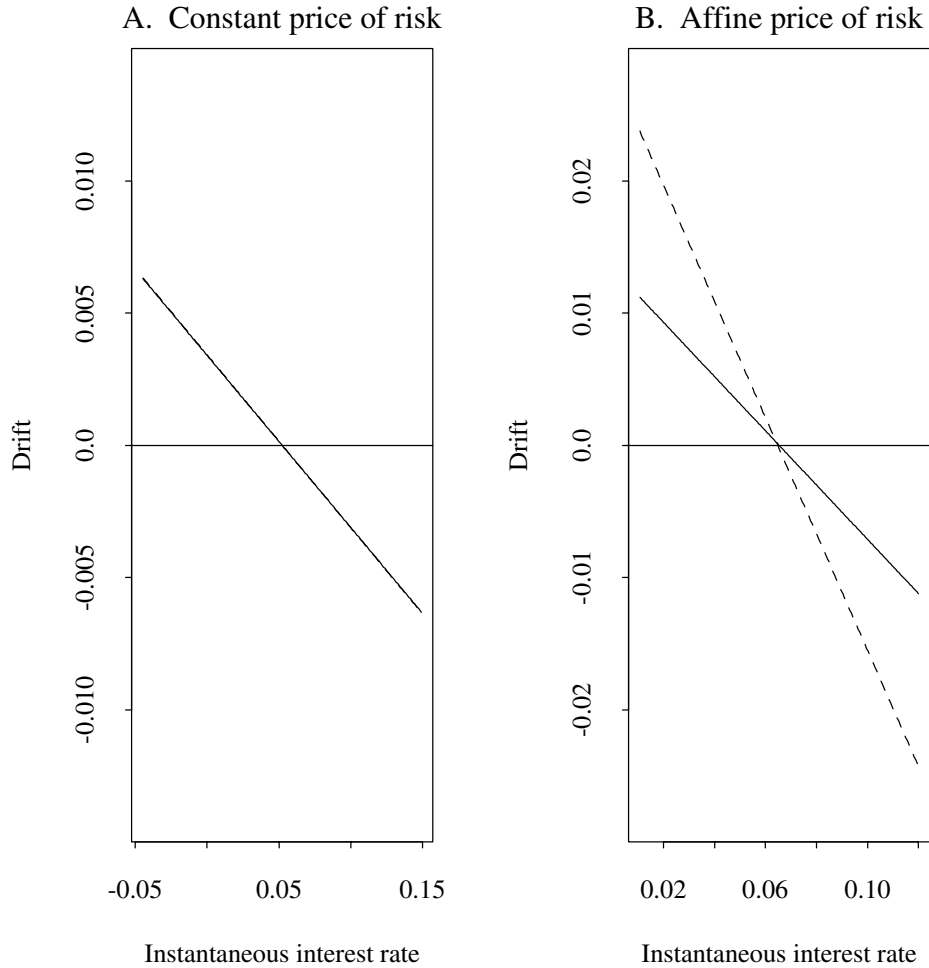


Figure 1: True and estimated drift functions for one-factor Gaussian models

Panels A and B display drift functions for r_t from two Gaussian term structure models. The model for Panel A specifies a constant price of interest rate risk and the model for Panel B specifies an affine price of interest rate risk. True drift functions are illustrated with solid lines. The dashed lines are drift functions implied by the mean parameter estimates from 500 Monte Carlo simulations of maximum likelihood estimation. The simulated data consists of a panel of 1000 weeks of yields on three bonds. The yields are all observed with normally distributed, iid measurement error of 60 basis points. The maturities of the bonds range from three months to ten years. The ranges of the x axes are approximately two unconditional standard deviations of r_t (which differ across the models).

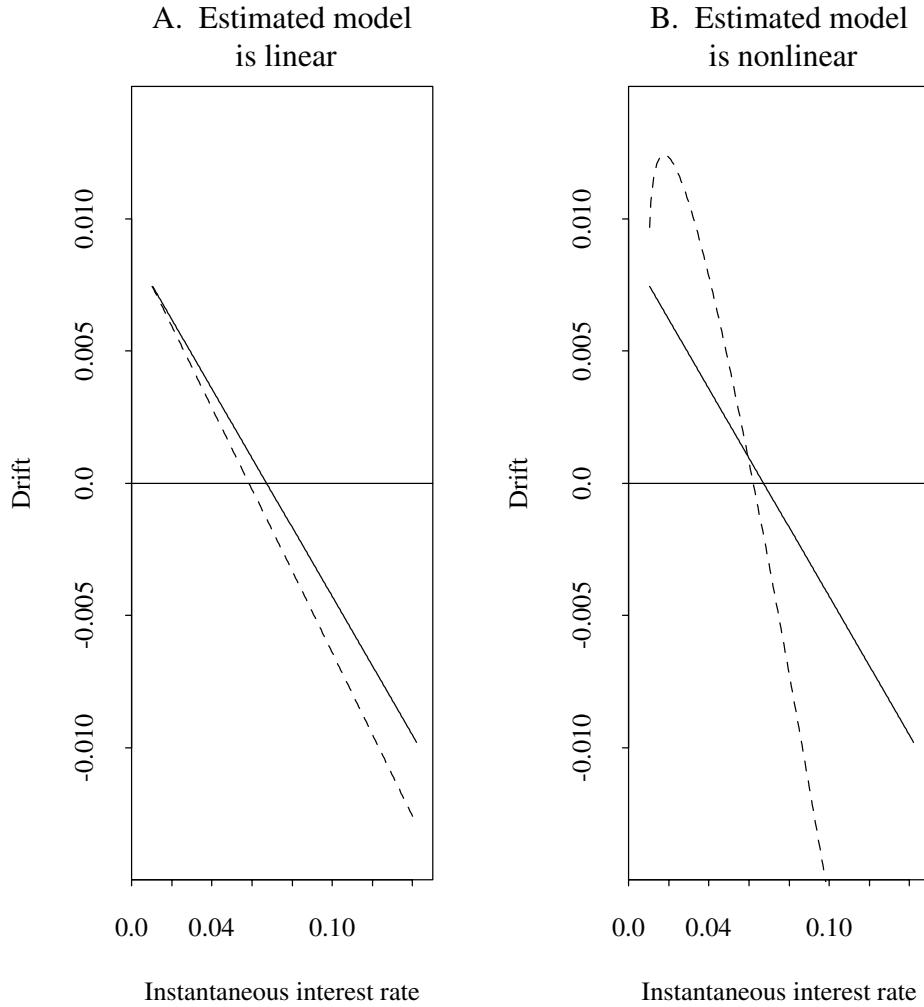


Figure 2: True and estimated drift functions for a one-factor square-root model

Simulated data of 1000 weeks of yields on three bonds are generated by a one-factor translated CIR model. The maturities of the bonds range from three months to ten years. One bond's yield is observed without error and the other bonds' yields are observed with normally distributed, iid measurement error of 60 basis points. The true model's drift function for r_t is indicated by the identical solid lines. The dashed lines are drift functions implied by mean parameter estimates from 500 Monte Carlo simulations of linearized instantaneous Kalman filter estimation. The estimated model in Panel A is a translated CIR model. The estimated model in Panel B nests the translated CIR model by allowing for a nonlinear drift under the physical measure.

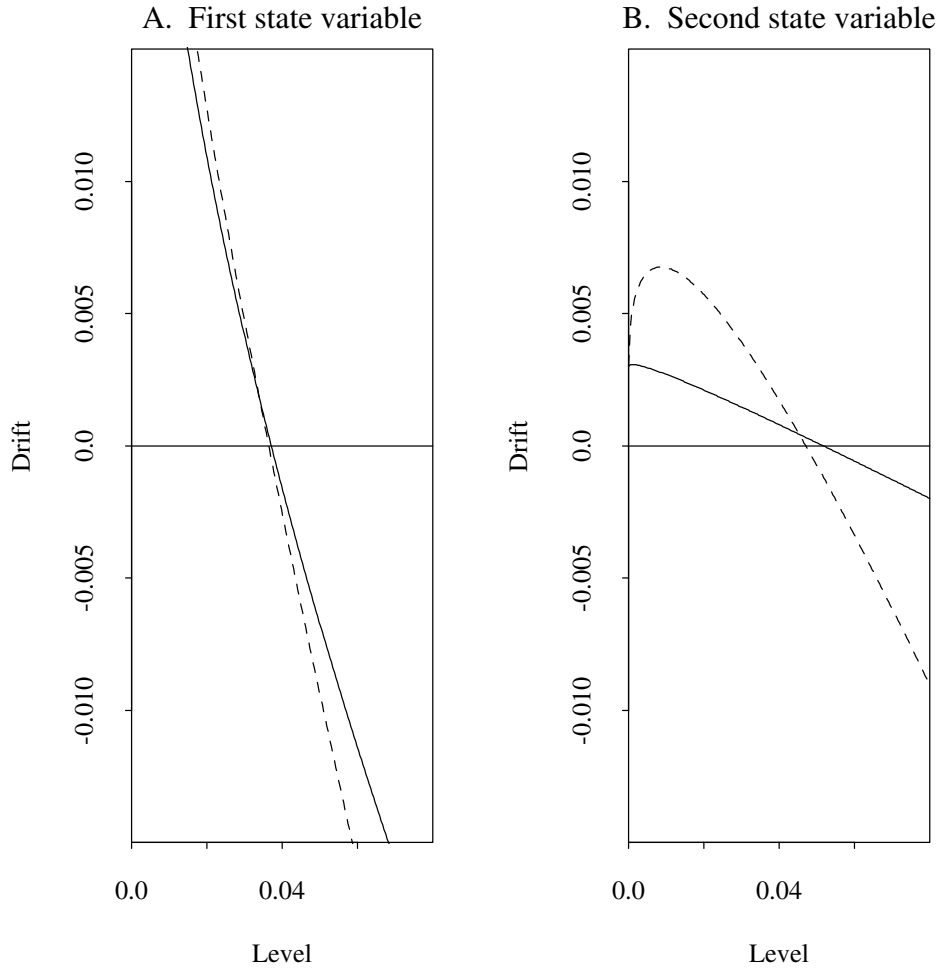


Figure 3: True and estimated drift functions for a two-factor nonlinear model

Simulated data of 1000 weeks of yields on three bonds are generated by a two-factor square-root diffusion model with nonlinear drifts. The maturities of the bonds range from three months to ten years. Two of the bonds' yields are observed without error and the other bond's yield is observed with normally distributed, iid measurement error of 20 basis points. The true model's drift functions for the state variables are indicated by the solid lines. The dashed lines are drift functions implied by mean parameter estimates from 500 Monte Carlo simulations of simulated maximum likelihood.