

Maximum likelihood estimation for integrated diffusion processes

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Abstract

We propose a method for obtaining maximum likelihood estimates of parameters in diffusion models when the data is a discrete time sample of the integral of the process, while no direct observations of the process itself are available. The data are, moreover, assumed to be contaminated by measurement errors. Integrated volatility is an example of this type of observations. Another example is ice-core data on oxygen isotopes used to investigate paleo-temperatures.

The data can be viewed as incomplete observations of a model with a tractable likelihood function. Therefore we propose a simulated EM-algorithm to obtain maximum likelihood estimates of the parameters in the diffusion model. As part of the algorithm, we use a recent simple method for approximate simulation of diffusion bridges. In a simulation study for the Ornstein-Uhlenbeck process the methods works well. The method is applied to a set of integrated paleo-temperature data obtained from an ice-core.

1 Introduction.

We consider estimation for a general one-dimensional diffusion process $X = \{X_t\}_{t \geq 0}$. Likelihood based estimation (including Bayesian) for discretely observed diffusion processes has been investigated by Ozaki (1985), Pedersen (1995), Poulsen (1999), Elerian, Chib & Shephard (2001), Eraker (2001), Roberts & Stramer (2001), Aït-Sahalia (2002), Durham & Galant (2002), Aït-Sahalia & Mykland (2003), Beskos et al. (2006) Aït-Sahalia (2008). Martingale estimating functions for discretely observed diffusions are reviewed in Sørensen (1997) and Sørensen (2010).

In this paper we consider maximum likelihood estimation in the situation where we do not observe the process X itself directly, but instead observe integrals of the process over disjoint time-intervals. These observations are, moreover, assumed to be contaminated by measurement errors. Integrated diffusion processes play an important role in finance as models for realized volatility, see e.g. Andersen et al. (2001b), Andersen et al. (2001a), Bollerslev & Zhou (2002), and Barndorff-Nielsen & Shephard (2002). These processes are also used for modelling purposes in fields of engineering and the sciences. An example is provided by the records of the concentration of oxygen isotopes in ice-core data from Greenland and Antarctica, see e.g. Ditlevsen, Ditlevsen & Andersen (2002). Such data are used to investigate the paleo-climate.

The likelihood function for a discretely sampled integrated diffusion with observation error is in almost all cases not explicitly available. Moreover, the integrated process is not a Markov process, so there is no easily calculated martingales. Therefore martingale estimating functions are not a feasible alternative, but prediction-based estimating function can be applied, see Sørensen (2000). In the present paper, we note instead that the data can be viewed as incomplete observations from a model with a tractable likelihood function. The full data set is a continuous time record of the diffusion process and the observation errors. We can therefore find maximum likelihood estimates by applying the Expectation-Maximization (EM) algorithm, see Dempster, Laird & Rubin (1977) and McLachlan & Krishnan (1997). To do so we need to calculate the conditional expectation of the log-likelihood function for the full model given the observations. We do this by simulating sample paths of the diffusion process given the data using ideas from Chib, Pitt & Shephard (2006). An essential step in doing this is to simulate a part of a sample path given the rest, which corresponds to simulation a diffusion bridge. This is done by applying the method for approximate diffusion bridge simulation recently proposed by Bladt & Sørensen (2009).

Parametric inference for integrated diffusion process has been considered by Gloter (2000), Bollerslev & Zhou (2002), Ditlevsen & Sørensen (2004), Gloter (2006), and Forman & Sørensen (2008). Nonparametric inference has been considered in Comte, Genon-Catalot & Rozenholc (2009).

In Section 2 the model of an integrated diffusion process with measurement error and its assumptions are presented. Section 3 contains the calculation of the likelihood function with full diffusion observation and the EM-algorithm. In Section 4 we present a method for simulation of a diffusion process conditional on integrals observed with measurement error. In Section 5 we consider the Ornstein-Uhlenbeck process in detail and a simulation study for this model is reported. In Section 6 the parameters of an Ornstein-Uhlenbeck model are estimated from a set of integrated paleo-temperature data obtained from an ice-core from Greenland. Some concluding remarks are given in Section 7.

2 Model and data

We consider likelihood estimation the the general one-dimensional diffusion process $X = \{X_t\}_{t \geq 0}$ given by the stochastic differential equation

$$dX_t = b(X_t; \psi)dt + \sigma(X_t; \psi)dW_t \quad (2.1)$$

where $W = \{W_t\}$ is a standard Wiener process, and where the drift and diffusion coefficients depend on an unknown p -dimensional parameter ψ belonging to the parameter set $\Psi \subseteq \mathbb{R}^p$. We assume that the solution X is an ergodic, stationary diffusion with invariant measure with density function $\nu_\psi(x)$ ($X_0 \sim \nu_\psi$ is independent of W). We also assume that the stochastic differential equation has a unique weak solution, i.e. a solution exists and all solutions have identical finite-dimensional distributions; see e.g. Kutoyants (2004). It is well-known that sufficient conditions for these assumptions can be expressed in terms of the so-called scale function and speed measure; see e.g. Karlin & Taylor (1981).

In this paper we consider the situation where the process X has not been observed directly. Instead the data are integrals of X_t over intervals $[t_{i-1}, t_i]$ observed with measurement error, i.e.

$$Y_i = \int_{t_{i-1}}^{t_i} X_s ds + Z_i, \quad i = 1, \dots, n, \quad (2.2)$$

where $Z_i \sim N(0, \tau^2)$, $i = 1, \dots, n$ are mutually independent and independent of X . We assume that $t_0 = 0$, so the total interval of observation is $[0, t_n]$. Note that the variance of the measurement error, τ^2 , is an extra unknown parameter. Thus we need to estimate the $p + 1$ -dimensional parameter $\theta = (\psi, \tau^2)$.

Conditionally on the sample path of X , the observations Y_i , $i = 1, \dots, n$ are independent and normal distributed:

$$Y_i | X_t : t \in [0, t_n] \sim N \left(\int_{t_{i-1}}^{t_i} X_s ds, \tau^2 \right), \quad (2.3)$$

We assume that the coefficients of the stochastic differential equation (2.1) satisfy the following conditions which we need in the following sections.

Condition 2.1 *The drift and diffusion coefficients of (2.1), $b(x; \psi)$ and $\sigma(x; \psi)$ satisfy that for all $\psi \in \Psi$*

- $b(x; \psi)$ is continuously differentiable w.r.t. x
- $\sigma(x; \psi)$ is twice continuously differentiable w.r.t. x
- $\sigma(x; \psi) > 0$ for all x in the state space of X

3 The likelihood function and the EM-Algorithm

We can think of the data set $Y = (Y_1, \dots, Y_n)$ as an incomplete observation of a full data set given by the sample path X_t , $t \in [0, t_n]$ and the measurement errors Z_1, \dots, Z_n , or equivalently X_t , $t \in [0, t_n]$ and $Y = (Y_1, \dots, Y_n)$. Therefore likelihood based estimation can be done by means of the EM-algorithm or MCMC-methods. In this paper we concentrate

on the EM-algorithm. We need to find the likelihood function for the full data set and the conditional expectation of this full log-likelihood function given the observations $Y = (Y_1, \dots, Y_n)$.

3.1 Likelihood with full diffusion observation

The full observation of a diffusion sample path in the time interval $[0, t_n]$ is an element in the space C of continuous functions from $[0, t_n]$ to \mathbb{R} . We equip this space with the usual σ -algebra, \mathcal{C} , generated by the cylinder sets, and consider the probability measures induced on (C, \mathcal{C}) by the solutions to (2.1). These measures are in general singular because the diffusion coefficient depends on the parameter ψ . In order to obtain a likelihood function, we use the standard 1-1 transformation

$$h(x; \psi) = \int_{x^*}^x \frac{1}{\sigma(u; \psi)} du, \quad (3.1)$$

where x^* is some arbitrary element of the state space of X . By this parameter dependent transformation, we obtain a diffusion process with unit diffusion coefficient. Specifically, we obtain (by Ito's formula) that

$$U_t = h(X_t; \psi)$$

satisfies the stochastic differential equation

$$dU_t = \mu(U_t; \psi)dt + dW_t, \quad (3.2)$$

with

$$\mu(u; \psi) = \frac{b(h^{-1}(u; \psi); \psi)}{\sigma(h^{-1}(u; \psi); \psi)} - \frac{\sigma'(h^{-1}(u; \psi); \psi)}{2},$$

where σ' denotes the derivative of σ w.r.t. x . In (3.2) the diffusion coefficient does not depend on the parameters, so the probability measures induced on (C, \mathcal{C}) by the solution to (3.2) are equivalent and the likelihood function can be found.

We can express the observations Y_i in terms of the process U . By inserting $X_s = h^{-1}(U_s; \psi)$ in (2.2), we find that

$$Y_i = \int_{t_{i-1}}^{t_i} h^{-1}(U_s; \psi) ds + Z_i, \quad i = 1, \dots, n.$$

Therefore we will think of the full dataset as U_t , $t \in [0, t_n]$ and $Y = (Y_1, \dots, Y_n)$. Since conditionally on the sample path of U the observations Y_i , $i = 1, \dots, n$ are independent, we have that the likelihood of Y conditional on the sample path of U in $[0, t_n]$ is

$$L(Y_1, \dots, Y_n | U_t, t \in [0, t_n]) = \prod_{i=1}^n \phi(Y_i; \int_{t_{i-1}}^{t_i} h^{-1}(U_s; \psi) ds, \tau^2) \quad (3.3)$$

where $\phi(u; a_1, a_2)$ denotes the density of the normal distribution with mean a_1 and variance a_2 evaluated at u .

Let P_ψ be the probability measure induced by $U = \{U_t\}_{t \in [0, t_n]}$ on (C, \mathcal{C}) , i.e. the probability measure with respect to which the coordinate process has the same distribution as U , and let Q be the Wiener measure on (C, \mathcal{C}) . We assume that the coefficient μ satisfies

conditions ensuring that the Girsanov theorem holds so that we have the Radon-Nykodym derivative

$$\frac{dP_\psi}{dQ}(B) = \exp \left\{ \int_0^{t_n} \mu(B_t; \psi) dB_t - \frac{1}{2} \int_0^{t_n} \mu^2(B_t; \psi) dt \right\}; \quad (3.4)$$

see e.g. Liptser & Shiryaev (1977), Jacod & Shiryaev (1987) or Øksendal (1998).

The evaluation of $\frac{dP_\psi}{dQ}$ is impossible because of the Ito integral term. To simplify the likelihood function, we apply the transformation

$$a(x; \psi) = \int^x \mu(u; \psi) du$$

(any antiderivative of μ), which under Condition 2.1 is twice continuously differentiable. By Ito's formula

$$\int_0^{t_n} \mu(B_t) dB_t = a(B_{t_n}; \psi) - a(B_0; \psi) - \frac{1}{2} \int_0^{t_n} \mu'(B_t; \psi) dt,$$

where μ' denotes the derivative of $\mu(u; \psi)$ w.r.t. u . We can now write the likelihood function (3.4) as

$$\frac{dP_\psi}{dQ}(B) = \exp \left\{ a(B_{t_n}; \psi) - a(B_0; \psi) - \frac{1}{2} \int_0^{t_n} [\mu(B_t; \psi)^2 + \mu'(B_t; \psi)] dt \right\}.$$

By combining this expression and (3.3), we see that the log-likelihood function for θ based on the full data set U_t , $t \in [0, t_n]$ and $Y = (Y_1, \dots, Y_n)$ is given by

$$\begin{aligned} \log L(\theta; Y_1, \dots, Y_n, U_t, t \in [0, t_n]) &= \sum_{i=1}^n \log \phi(Y_i; \int_{t_{i-1}}^{t_i} h^{-1}(U_s; \psi) ds, \tau^2) \\ &+ a(U_{t_n}; \psi) - a(U_0; \psi) - \frac{1}{2} \int_0^{t_n} (\mu(U_t; \psi)^2 + \mu'(U_t; \psi)) dt. \end{aligned} \quad (3.5)$$

3.2 EM Algorithm.

We can now apply the EM-algorithm to the full log-likelihood function (3.5) to obtain the maximum likelihood estimate of the parameter θ .

As the initial value for the algorithm, let $\hat{\theta}$ be any value of the parameter vector $\theta = (\psi, \tau^2) \in \Psi \times (0, \infty)$. Then the EM-algorithm works as follow.

1. E-STEP.

Generate M sample paths of the diffusion process X , $X^{(k)}$, $k = 1, \dots, M$, conditional on the observations Y_1, \dots, Y_n using the parameter value $\hat{\theta} = (\hat{\psi}, \hat{\tau}^2)$, and calculate

$$g(\theta) = \frac{1}{M - M_0} \sum_{k=M_0+1}^M \log L(\theta; Y_1, \dots, Y_n, h(X_t^{(k)}; \hat{\psi}), t \in [0, t_n]),$$

for a suitable burn-in period M_0 and M sufficiently large.

2. M-STEP.

$$\hat{\theta} = \operatorname{argmax} g(\theta).$$

3. Go to 1.

To implement this algorithm, the main issue is how to generate sample paths of X conditionally on Y_1, \dots, Y_n , where the relation between the Y_i s and X is given by (2.2). The algorithm must produce a sequence $X^{(k)}$, $k = 1, \dots, M$, that is sufficiently mixing to ensure that $g(\theta)$ approximates the conditional expectation of the full log-likelihood function (3.5) given the data. This problem is discussed at the next section.

4 Conditional diffusion process simulation.

In this section we present a method for generating a sample from

$$\{X_t; t \in [0, t_n]\} | (Y_1, \dots, Y_n)$$

for a given value of the parameter vector θ , i.e. for simulating the diffusion X conditional on the observations $Y = (Y_1, \dots, Y_n)$ of integrals of X over subintervals $[t_{j-1}, t_j]$, $j = 1, \dots, n$ perturbed by measurement errors. This can be done by means of a Metropolis-Hastings algorithm, see e.g. Chib & Greenberg (1995) or Gilks, Richardson & Spiegelhalter (1996). However, if the sample path in the entire time interval $[0, t_n]$ is updated in one step, the rejection probability is typically very large. Therefore it is more efficient to randomly divide the time interval into subintervals and update the sample path in each of the subintervals conditional on the rest of the sample path. This corresponds to simulating a (conditional) diffusion bridge in each subinterval (except the end-intervals). The method outlined in this section is a modification of the method in Chib, Pitt & Shephard (2006), where we use the algorithm for approximate diffusion bridge simulation proposed by Bladt & Sørensen (2009).

In the following the parameter value $\theta = (\psi, \tau^2)$ is fixed.

Algorithm 1

1. Generate an initial unrestricted stationary sample path, $\{X_t^{(0)} : t \in [0, t_n]\}$, of the diffusion given by (2.1) using for instance the Milstein scheme or one of the other methods in Kloeden & Platen (1999).
2. Set $l = 1$.
3. Generate a sample path $\{X_t^{(l)} : t \in [0, t_n]\}$ conditional on Y by updating the subsets of the sample path:

- (a) Randomly split the time interval from 0 to t_n in K blocks, and write these sub-sampling times as

$$0 = \tau_0 \leq \tau_1 \leq \dots \leq \tau_K = t_n,$$

where each τ_i is one of the end-points of the integration intervals, t_j , $j = 0, \dots, n$. Let $Y_{\{k\}}$ denote the collection of all observations Y_j for which $\tau_{k-1} < t_j \leq \tau_k$.

- (b) Draw $X_0^{(l)}$ from the stationary distribution, ν_ψ , and simulate the conditional sub-path

$$\{X_t^{(l)} : t \in [\tau_{k-1}, \tau_k]\} | Y_{\{k\}}, X_{\tau_{k-1}}^{(l)}, X_{\tau_k}^{(l-1)} \quad (4.1)$$

for $k = 1, \dots, K - 1$. Finally, simulate a sample path from

$$\{X_t^{(l)} : t \in [\tau_{K-1}, \tau_K]\} | Y_{\{K\}}, X_{\tau_{K-1}}^{(l)}.$$

4. $l=l+1$.

5. Go to 3.

To implement of this algorithm, the main issue is how to sample variables of the type (4.1), which is a non-linear diffusion bridge. We use the method for approximate diffusion bridge simulation proposed by Bladt & Sørensen (2009). The idea (in the case of a diffusion bridge in the time interval $[0, 1]$) is to let one diffusion process move forward from time zero out of one given point, a , until it meets another diffusion process that independently moves backwards from time one out of another given point, b . Conditional on the event that the two diffusions intersect, the process constructed in this way is an approximation to a realization of a diffusion bridge between a and b . The diffusions can be simulated by means of simple procedures like the Euler scheme or the Milstein scheme, see Kloeden & Platen (1999). The method is therefore very easy to implement. The resulting sample path is an approximation to a diffusion bridge in the sense that it has the distribution of a diffusion bridge from a to b conditional on the event that the bridge is hit by an independent diffusion with stochastic differential equation (2.1) and initial distribution with density $p_1(b, \cdot)$. Simulation studies in Bladt & Sørensen (2009) indicate that the approximation is very good for bridges between points that are likely to appear on a sample path of the diffusion, which is the type of bridges that are relevant to this paper.

Alternative methods that provide exact diffusion bridges have been proposed by Beskos, Papaspiliopoulos & Roberts (2006) and Beskos, Papaspiliopoulos & Roberts (2007). When the drift and diffusion coefficients satisfy certain boundedness conditions, this algorithm is relatively simple, but under weaker condition it is more complex. A simulation study in Bladt & Sørensen (2009) indicates that for the method which we use here, the CPU-time is linear in the length of the interval where the diffusion bridge is defined, whereas for the method in Beskos, Papaspiliopoulos & Roberts (2006), the CPU time increases exponentially with the interval length. This is an advantage of the method in Bladt & Sørensen (2009) in the present context. MCMC algorithms for simulation of diffusion bridges were proposed by Roberts & Stramer (2001), Durham & Gallant (2002), and Chib, Pitt & Shephard (2006).

To generate the random subintervals in step 3 (a) of Algorithm 1, we use the following algorithm, where the number of integration subintervals $[t_{j-1}, t_j]$ included in one of the random subintervals is a Poisson distributed random number plus 1. The draws in the algorithm are independent. First choose the expectation of the Poisson distribution, $\lambda \geq 1$.

Algorithm 2

1. Draw $k_1 \sim \text{Poisson}(\lambda) + 1$: if $k_1 \geq n$ set $k_1 = n$, $K = 1$ and stop, otherwise set $i = 2$.
2. Draw $k_i \sim \text{Poisson}(\lambda) + 1$, if $\sum_{j=1}^i k_j \geq n$ set $k_i = n$, $K = i$ and stop, else set $i = i + 1$ and repeat 2.

Finally define $\tau_i = t_{k_i}$, $i = 1, \dots, K$.

We have discussed how to simulate diffusion bridges, but we need diffusion bridges conditional on the data Y . Sample paths of the conditional bridges (4.1) can be obtained by the following Metropolis-Hastings algorithm. By a (t, a, s, b) -bridge, we mean a diffusion bridge in the time interval $[t, s]$ with $X_t = a$ and $X_s = b$. After a burn-in period the following algorithm will output samples from a $(\tau_{k-1}, a, \tau_k, b)$ -bridge conditional on $Y_{\{k\}}$, the data in $(\tau_{k-1}, \tau_k]$. To formulate the algorithm we need to specify that the end-point τ_{k-1} is equal to t_j , and that there are n_k observations in the interval $(\tau_{k-1}, \tau_k]$, namely, $Y_{j+1}, \dots, Y_{j+n_k}$.

Algorithm 3

1. Simulate a $(\tau_{k-1}, a, \tau_k, b)$ -bridge, $X^{(0)}$, and set $l = 1$.
2. Propose a new sample paths by simulating a $(\tau_{k-1}, a, \tau_k, b)$ -bridge, $X^{(l)}$.
3. Accept the proposed diffusion bridge with probability

$$\min \left(1, \prod_{i=1}^{n_k} \frac{\phi(Y_{j+i}; \int_{t_{j+i-1}}^{t_{j+i}} X_s^{(l)} ds, \tau^2)}{\phi(Y_{j+i}; \int_{t_{j+i-1}}^{t_{j+i}} X_s^{(l-1)} ds, \tau^2)} \right).$$

Otherwise set $X^{(l)} = X^{(l-1)}$.

4. Set $l = l + 1$ and go to 2.

As previously, $\phi(x; \mu, \tau^2)$ denotes the density function of the normal distribution with mean μ and variance τ^2 .

5 The Ornstein-Uhlenbeck process: a simulation study.

In this section we apply the method developed above to the Ornstein-Uhlenbeck process, which is a solution of the stochastic differential equation

$$dX_t = -\alpha X_t dt + \sigma dW_t, \tag{5.1}$$

where $\alpha > 0$ and $\sigma > 0$ are unknown parameters to be estimated, and W is a standard Wiener process. We investigate the bias of the estimators in a simulation study.

5.1 The likelihood and the EM-algorithm

The transformation (3.1) is here given by

$$h(x; \sigma) = \frac{x}{\sigma},$$

so $h^{-1}(x; \sigma) = \sigma x$. Hence $U_t = h(X_t; \sigma) = X_t/\sigma$, solves the stochastic differential equation

$$dU_t = -\alpha U_t dt + dW_t.$$

We have $\mu(u; \alpha, \sigma) = -\alpha u$, so

$$a(u; \alpha, \sigma) = -\frac{1}{2}\alpha u^2.$$

Thus the full log-likelihood function (3.5) is given by

$$\begin{aligned} & \log L(\theta; Y_1, \dots, Y_n, U_t, t \in [0, t_n]) \\ &= \sum_{i=1}^n \log \phi \left(Y_i; \sigma \int_{t_{i-1}}^{t_i} U_s ds, \tau^2 \right) + \frac{\alpha}{2}(U_0^2 - U_{t_n}^2 + t_n) - \frac{\alpha^2}{2} \int_0^{t_n} U_t^2 dt, \end{aligned} \quad (5.2)$$

where $\theta = (\alpha, \sigma, \tau^2)$.

Now, the EM algorithm works as follow.

E-STEP

The objective function $g(\theta)$ is for the Ornstein-Uhlenbeck process given by

$$\begin{aligned} g(\theta) &= \frac{1}{M - M_0} \sum_{k=M_0+1}^M \sum_{i=1}^n -\frac{(Y_i - \sigma \int_{t_{i-1}}^{t_i} U_t^{(m)} dt)^2}{2\tau^2} - \frac{n}{2} \log(2\pi\tau^2) + \frac{\alpha}{2} t_n \\ &+ \frac{\alpha}{2(M - M_0)} \sum_{k=M_0+1}^M ((U_0^{(m)})^2 - (U_{t_n}^{(m)})^2) - \frac{\alpha^2}{2(M - M_0)} \sum_{k=M_0+1}^M \sum_{i=1}^n \int_{t_{i-1}}^{t_i} (U_t^{(m)})^2 dt. \end{aligned}$$

Here $U_t^{(m)} = X_t^{(m)}/\hat{\sigma}$, where $X_t^{(m)}$ is the m -th sample path of the process X simulated conditionally on the data Y using the Algorithms 1 – 3 with the parameter value obtained in the previous step $(\hat{\alpha}, \hat{\sigma}, \hat{\tau}^2)$.

M-STEP

The maximum $\hat{\theta}$ is obtained as the solution to the following system of equations

$$\frac{\partial g(\theta)}{\partial \alpha} = \frac{1}{2} t_n + \frac{\sum_{k=M_0+1}^M [(U_0^{(m)})^2 - (U_{t_n}^{(m)})^2]}{2(M - M_0)} - \frac{\alpha \sum_{k=M_0+1}^M \sum_{i=1}^n \int_{t_{i-1}}^{t_i} (U_t^{(m)})^2 dt}{M - M_0} = 0, \quad (5.3)$$

$$\frac{\partial g(\theta)}{\partial \sigma} = \frac{\sum_{k=M_0+1}^M \sum_{i=1}^n (Y_i - \sigma \int_{t_{i-1}}^{t_i} U_t^{(m)} dt) (\int_{t_{i-1}}^{t_i} U_t^{(m)} dt)}{\tau^2 (M - M_0)} = 0 \quad (5.4)$$

and

$$\frac{\partial g(\theta)}{\partial \tau^2} = \frac{\sum_{k=M_0+1}^M \sum_{i=1}^n (Y_i - \sigma \int_{t_{i-1}}^{t_i} U_t^{(m)} dt)^2}{2\tau^4 (M - M_0)} - \frac{n}{2\tau^2} = 0. \quad (5.5)$$

From (5.3) we have

$$\hat{\alpha} = \frac{t_n (M - M_0) + \sum_{k=M_0+1}^M [(U_0^{(m)})^2 - (U_{t_n}^{(m)})^2]}{2 \sum_{k=M_0+1}^M \sum_{i=1}^n \int_{t_{i-1}}^{t_i} (U_t^{(m)})^2 dt},$$

and from (5.4)

$$\hat{\sigma} = \frac{\sum_{k=M_0+1}^M \sum_{i=1}^n Y_i \int_{t_{i-1}}^{t_i} U_t^{(m)} dt}{\sum_{k=M_0+1}^M \sum_{i=1}^n (\int_{t_{i-1}}^{t_i} U_t^{(m)} dt)^2}. \quad (5.6)$$

Now inserting $\hat{\sigma}$ given by (5.6) in (5.5) we obtain

$$\hat{\tau}^2 = \frac{(M - M_0)(\sum_{i=1}^n Y_i^2) \left[\sum_{k=M_0+1}^M \sum_{i=1}^n (\int_{t_{i-1}}^{t_i} U_t^{(m)} dt)^2 \right] - \left[\sum_{k=M_0+1}^M \sum_{i=1}^n Y_i \int_{t_{i-1}}^{t_i} U_t^{(m)} dt \right]^2}{n(M - M_0) \sum_{k=M_0+1}^M \sum_{i=1}^n (\int_{t_{i-1}}^{t_i} U_t^{(m)} dt)^2}.$$

The Hessian matrix of $g(\theta)$ evaluated at $\hat{\theta}$ is negative definite, so $\hat{\theta}$ is maximum.

5.2 A simulation study

In this section we present the result of a small simulation study, in which we simulated 1000 datasets and for each of them obtained estimates by means of the EM-algorithm proposed in the present paper. Each data set was obtained by simulating a sample path of length 1500 with initial distribution $X_0 \sim N(0, \sigma^2/(2\alpha))$, and then calculating data Y_i , $i = 1, \dots, 1500$ by (2.2) with $t_i = i$, $i = 0, \dots, n$. The parameter values were $\alpha = 0.1$, $\sigma = 0.5$ and $\tau^2 = 1.25$.

The EM-algorithm was run with $M = 10000$ and $M_0 = 1000$ and for three different values of λ , namely $\lambda = 10, 20, 30$. The average of the estimates obtained for the 1000 dataset are given in Table 5.1. The bias is small, and is overall most satisfactory for $\lambda = 20$.

λ	α	σ	τ^2
10	0.106	0.523	1.229
20	0.101	0.507	1.235
30	0.084	0.458	1.252

Table 5.1: Average of parameter estimates obtained from 1000 simulated datasets with parameter values $\alpha = 0.1$, $\sigma = 0.5$ and $\tau^2 = 1.25$.

6 Ice-core data

In this section we consider as an example a preliminary analysis of ice-cores data. Ice-cores from Greenland and Antarctica are cut in pieces each of which represents a time interval in the past, with time increasing as a function of the depth. The isotope ratio $^{18}O/^{16}O$ in the ice, measured as an average in each piece of ice, is a proxy for paleo-temperatures averaged over the time interval in which the snow that formed the ice fell. The difference, $\delta^{18}O$, between the ^{18}O content in the ice and the ^{18}O content in present day's ocean water has been shown empirically to have an approximately constant linear relationship with the air temperature where the precipitation falls. The variation of the paleo-temperature can be modelled by a stochastic differential equation, and the ice-core data can thus be modelled as an integrated diffusion process, see Ditlevsen, Ditlevsen & Andersen (2002), who modelled the temperature with an Ornstein-Uhlenbeck process in suitable subintervals, where the process could be assumed to be reasonably stationary.

We consider a data set obtained from an ice core obtained from the Greenland ice-sheet, see Svensson et al. (2008). The record of integrated $\delta^{18}O$ goes back 60,000 years, and each data point is the average over a time interval of length 20 years. Thus there are 3000 data points. The time series is plotted in Figure 6.1. It is obviously not stationary, so in a preliminary analysis we divide it into three subintervals, $[-60000, -30000]$, $[-30000, -10000]$,

and $[-10000, 0]$. In each of these intervals we model the observations by (2.2) where X is an Ornstein-Uhlenbeck process, and the integrals are over intervals of length 20. It is quite obvious that the Ornstein-Uhlenbeck process does not fit the data well at this time scale, so in a more full analysis either much shorter subintervals must be considered, or preferably, a more complex model must be developed.

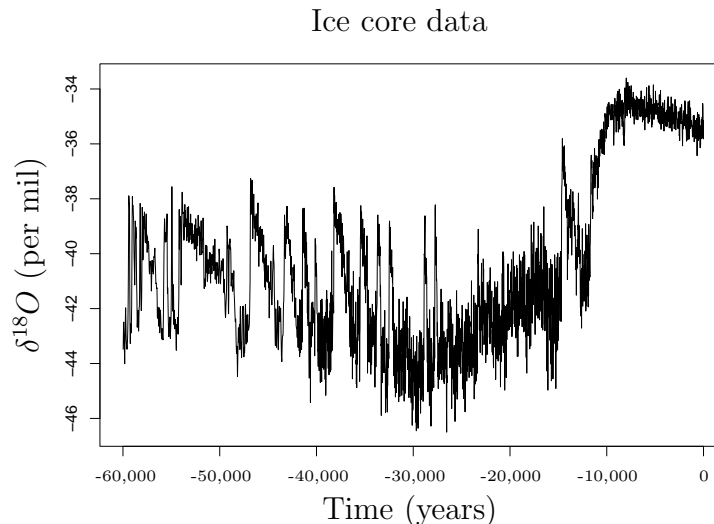


Figure 6.1: $\delta^{18}O$ -values integrated over 20 years intervals obtained from ice core data from the Greenland ice-sheet.

The EM algorithm is run with $M = 10000$, $M_0 = 1000$, and $\lambda = 20$ for each of the three blocks of data. The resulting estimates are reported in Table 6.1. The parameter α^{-1} can be interpreted as a correlation time, and $\sigma/\sqrt{2\alpha}$ is the standard deviation of the stationary distribution. When compared to the plot, these parameter values seem reasonable.

Parameter	$[-10000,0]$	$[-30000,-10000]$	$[-60000,-30000]$
α^{-1}	205.3	669.8	321.1
σ	0.0303	0.1167	0.1395
$\sigma/\sqrt{2\alpha}$	0.307	2.136	1.767
τ^2	0.1063	0.6967	0.2260

Table 6.1: Estimates obtained by the EM-algorithm for the ice core data using an Ornstein-Uhlenbeck model for each of the three subintervals.

7 Concluding remarks

We have presented an EM-algorithm for obtaining maximum likelihood estimates of parameters in diffusion models when the data is a discrete time sample of the integral of the diffusion process contaminated by measurement errors, while no direct observations of the

process itself are available. This was done by viewing the data as an incomplete observation, where the full data set includes a continuous time record of the diffusion process.

It is not difficult to generalize the method presented in this paper to the situation, where the diffusion process is integrated w.r.t. a more general measure than the Lebesgue measure considered in this paper. This would allow analysis of e.g. weighted averages of diffusion processes, and discrete time observation would be a particular case. Note also that a Gibbs sampler could easily be set up in close analogy to the EM-algorithm used in the present paper. This would be much closer to the approach in Chib, Pitt & Shephard (2006).

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