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# A Novel Gaussian Sum Smoother for Approximate Inference in Switching Linear Dynamical Systems\*

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## Abstract

We introduce a method for approximate smoothed inference in a class of switching linear dynamical systems, based on a novel form of Gaussian Sum smoother. This class includes the switching Kalman Filter and the more general case of switch transitions dependent on the continuous latent state. The method improves on the standard Kim smoothing approach by dispensing with one of the key approximations, thus making fuller use of the available future information. Whilst the only central assumption required is projection to a mixture of Gaussians, we show that an additional conditional independence assumption results in a simpler but accurate alternative. Unlike the alternative Expectation Propagation procedure, our method consists only of a single forward and backward pass and is reminiscent of the standard smoothing ‘correction’ recursions in the simpler linear dynamical system. The algorithm performs well on both toy experiments and in a large scale application to noise robust speech recognition.

## 1 Switching Linear Dynamical System

The Linear Dynamical System (LDS) [1] is a key temporal model in which a latent linear process generates the observed series. For complex time-series which are not well described globally by a single LDS, we may break the time-series into segments, each modeled by a potentially different LDS. This is the basis for the Switching LDS (SLDS) [2, 3, 4, 5] where, for each time  $t$ , a switch variable  $s_t \in 1, \dots, S$  describes which of the LDSs is to be used. The observation (or ‘visible’)  $v_t \in \mathcal{R}^V$  is linearly related to the hidden state  $h_t \in \mathcal{R}^H$  with additive noise  $\eta$  by

$$v_t = B(s_t)h_t + \eta^v(s_t) \equiv p(v_t|h_t, s_t) = \mathcal{N}(B(s_t)h_t, \Sigma^v(s_t)) \quad (1)$$

where  $\mathcal{N}(\mu, \Sigma)$  denotes a Gaussian distribution with mean  $\mu$  and covariance  $\Sigma$ . The transition dynamics of the continuous hidden state  $h_t$  is linear,

$$h_t = A(s_t)h_{t-1} + \eta^h(s_t), \equiv p(h_t|h_{t-1}, s_t) = \mathcal{N}(A(s_t)h_{t-1}, \Sigma^h(s_t)) \quad (2)$$

The switch  $s_t$  may depend on both the previous  $s_{t-1}$  and  $h_{t-1}$ . This is an augmented SLDS (aSLDS), and defines the model

$$p(v_{1:T}, h_{1:T}, s_{1:T}) = \prod_{t=1}^T p(v_t|h_t, s_t)p(h_t|h_{t-1}, s_t)p(s_t|h_{t-1}, s_{t-1})$$

The standard SLDS[4] considers only switch transitions  $p(s_t|s_{t-1})$ . At time  $t = 1$ ,  $p(s_1|h_0, s_0)$  simply denotes the prior  $p(s_1)$ , and  $p(h_1|h_0, s_1)$  denotes  $p(h_1|s_1)$ .

The aim of this article is to address how to perform inference in the aSLDS. In particular we desire the *filtered* estimate  $p(h_t, s_t|v_{1:t})$  and the *smoothed* estimate  $p(h_t, s_t|v_{1:T})$ , for any  $1 \leq t \leq T$ . Both filtered and smoothed inference in the SLDS is intractable, scaling exponentially with time [4].

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\*Version to appear in NIPS 2006

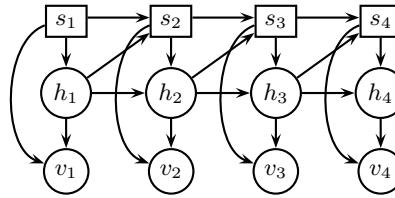


Figure 1: The independence structure of the aSLDS. Square nodes denote discrete variables, round nodes continuous variables. In the SLDS links from  $h$  to  $s$  are not normally considered.

## 2 Expectation Correction

Our approach to approximate  $p(h_t, s_t | v_{1:T})$  mirrors the Rauch-Tung-Striebel ‘correction’ smoother for the simpler LDS [1]. The method consists of a single forward pass to recursively find the filtered posterior  $p(h_t, s_t | v_{1:t})$ , followed by a single backward pass to correct this into a smoothed posterior  $p(h_t, s_t | v_{1:T})$ . The forward pass we use is equivalent to standard Assumed Density Filtering (ADF) [6]. The main contribution of this paper is a novel form of backward pass, based only on collapsing the smoothed posterior to a mixture of Gaussians. Together with the ADF forward pass, we call the method Expectation Correction, since it corrects the moments found from the forward pass. A more detailed description of the method, including pseudocode, is given in [7].

### 2.1 Forward Pass (Filtering)

Readers familiar with ADF may wish to continue directly to Section (2.2). Our aim is to form a recursion for  $p(s_t, h_t | v_{1:t})$ , based on a Gaussian mixture approximation of  $p(h_t | s_t, v_{1:t})$ . Without loss of generality, we may decompose the filtered posterior as

$$p(h_t, s_t | v_{1:t}) = p(h_t | s_t, v_{1:t})p(s_t | v_{1:t}) \quad (3)$$

The exact representation of  $p(h_t | s_t, v_{1:t})$  is a mixture with  $O(S^t)$  components. We therefore approximate this with a smaller  $I$ -component mixture

$$p(h_t | s_t, v_{1:t}) \approx \sum_{i_t=1}^I p(h_t | i_t, s_t, v_{1:t})p(i_t | s_t, v_{1:t})$$

where  $p(h_t | i_t, s_t, v_{1:t})$  is a Gaussian parameterised with mean  $f(i_t, s_t)$  and covariance  $F(i_t, s_t)$ . To find a recursion for these parameters, consider

$$p(h_{t+1} | s_{t+1}, v_{1:t+1}) = \sum_{s_t, i_t} p(h_{t+1} | s_t, i_t, s_{t+1}, v_{1:t+1})p(s_t, i_t | s_{t+1}, v_{1:t+1}) \quad (4)$$

#### Evaluating $p(h_{t+1} | s_t, i_t, s_{t+1}, v_{1:t+1})$

We find  $p(h_{t+1} | s_t, i_t, s_{t+1}, v_{1:t+1})$  from the joint distribution  $p(h_{t+1}, v_{t+1} | s_t, i_t, s_{t+1}, v_{1:t})$ , which is a Gaussian with covariance and mean elements:

$$\begin{aligned} \Sigma_{hh} &= A(s_{t+1})F(i_t, s_t)A^\top(s_{t+1}) + \Sigma^h(s_{t+1}), & \Sigma_{vv} &= B(s_{t+1})\Sigma_{hh}B^\top(s_{t+1}) + \Sigma^v(s_{t+1}) \\ \Sigma_{vh} &= B(s_{t+1})F(i_t, s_t), & \mu_v &= B(s_{t+1})A(s_{t+1})f(i_t, s_t), & \mu_h &= A(s_{t+1})f(i_t, s_t) \end{aligned} \quad (5)$$

To find  $p(h_{t+1} | s_t, i_t, s_{t+1}, v_{1:t+1})$  we may then condition  $p(h_{t+1}, v_{t+1} | s_t, i_t, s_{t+1}, v_{1:t})$  on  $v_{t+1}$  using conditioning<sup>1</sup>. For the case  $S = 1$ , this forms the usual Kalman Filter recursions[1].

#### Evaluating $p(s_t, i_t | s_{t+1}, v_{1:t+1})$

The mixture weight in (4) can be found from the decomposition

$$p(s_t, i_t | s_{t+1}, v_{1:t+1}) \propto p(v_{t+1} | i_t, s_t, s_{t+1}, v_{1:t})p(s_{t+1} | i_t, s_t, v_{1:t})p(i_t | s_t, v_{1:t})p(s_t | v_{1:t}) \quad (6)$$

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<sup>1</sup> $p(x|y)$  is a Gaussian with mean  $\mu_x + \Sigma_{xy}\Sigma_{yy}^{-1}(y - \mu_y)$  and covariance  $\Sigma_{xx} - \Sigma_{xy}\Sigma_{yy}^{-1}\Sigma_{yx}$ .

The first factor in (6),  $p(v_{t+1}|i_t, s_t, s_{t+1}, v_{1:t})$  is given as a Gaussian with mean  $\mu_v$  and covariance  $\Sigma_{vv}$ , as given in (5). The last two factors  $p(i_t|s_t, v_{1:t})$  and  $p(s_t|v_{1:t})$  are given from the previous iteration. Finally,  $p(s_{t+1}|i_t, s_t, v_{1:t})$  is found from

$$p(s_{t+1}|i_t, s_t, v_{1:t}) = \langle p(s_{t+1}|h_t, s_t) \rangle_{p(h_t|i_t, s_t, v_{1:t})} \quad (7)$$

where  $\langle \cdot \rangle_p$  denotes expectation with respect to  $p$ . In the SLDS, (7) is replaced by the Markov transition  $p(s_{t+1}|s_t)$ . In the aSLDS, however, (7) will generally need to be computed numerically.

### Closing the recursion

We are now in a position to calculate (4). For each setting of the variable  $s_{t+1}$ , we have a mixture of  $I \times S$  Gaussians which we numerically collapse back to  $I$  Gaussians to form

$$p(h_{t+1}|s_{t+1}, v_{1:t+1}) \approx \sum_{i_{t+1}=1}^I p(h_{t+1}|i_{t+1}, s_{t+1}, v_{1:t+1}) p(i_{t+1}|s_{t+1}, v_{1:t+1})$$

Any method of choice may be supplied to collapse a mixture to a smaller mixture; our code simply repeatedly merges low-weight components. In this way the new mixture coefficients  $p(i_{t+1}|s_{t+1}, v_{1:t+1})$ ,  $i_{t+1} \in 1, \dots, I$  are defined, completing the description of how to form a recursion for  $p(h_{t+1}|s_{t+1}, v_{1:t+1})$  in (3). A recursion for the switch variable is given by

$$p(s_{t+1}|v_{1:t+1}) \propto \sum_{s_t, i_t} p(v_{t+1}|s_{t+1}, i_t, s_t, v_{1:t}) p(s_{t+1}|i_t, s_t, v_{1:t}) p(i_t|s_t, v_{1:t}) p(s_t|v_{1:t})$$

where all terms have been computed during the recursion for  $p(h_{t+1}|s_{t+1}, v_{1:t+1})$ .

The likelihood  $p(v_{1:T})$  may be found by recursing  $p(v_{1:t+1}) = p(v_{t+1}|v_{1:t})p(v_{1:t})$ , where

$$p(v_{t+1}|v_t) = \sum_{i_t, s_t, s_{t+1}} p(v_{t+1}|i_t, s_t, s_{t+1}, v_{1:t}) p(s_{t+1}|i_t, s_t, v_{1:t}) p(i_t|s_t, v_{1:t}) p(s_t|v_{1:t})$$

## 2.2 Backward Pass (Smoothing)

The main contribution of this paper is to find a suitable way to ‘correct’ the filtered posterior  $p(s_t, h_t|v_{1:t})$  obtained from the forward pass into a smoothed posterior  $p(s_t, h_t|v_{1:T})$ . We derive this for the case of a single Gaussian representation. The extension to the mixture case is straightforward and is presented in [7]. We approximate the smoothed posterior  $p(h_t|s_t, v_{1:T})$  by a Gaussian with mean  $g(s_t)$  and covariance  $G(s_t)$ , and our aim is to find a recursion for these parameters. A useful starting point for a recursion is:

$$p(h_t, s_t|v_{1:T}) = \sum_{s_{t+1}} p(s_{t+1}|v_{1:T}) p(h_t|s_t, s_{t+1}, v_{1:T}) p(s_t|s_{t+1}, v_{1:T})$$

The term  $p(h_t|s_t, s_{t+1}, v_{1:T})$  may be computed as

$$p(h_t|s_t, s_{t+1}, v_{1:T}) = \int_{h_{t+1}} p(h_t|h_{t+1}, s_t, s_{t+1}, v_{1:t}) p(h_{t+1}|s_t, s_{t+1}, v_{1:T}) \quad (8)$$

The recursion therefore requires  $p(h_{t+1}|s_t, s_{t+1}, v_{1:T})$ , which we can write as

$$p(h_{t+1}|s_t, s_{t+1}, v_{1:T}) \propto p(h_{t+1}|s_{t+1}, v_{1:T}) p(s_t|s_{t+1}, h_{t+1}, v_{1:t}) \quad (9)$$

The difficulty here is that the functional form of  $p(s_t|s_{t+1}, h_{t+1}, v_{1:t})$  is not squared exponential in  $h_{t+1}$ , so that  $p(h_{t+1}|s_t, s_{t+1}, v_{1:T})$  will not be Gaussian. One possibility would be to approximate the non-Gaussian  $p(h_{t+1}|s_t, s_{t+1}, v_{1:T})$  by a Gaussian (or mixture thereof) by minimising the Kullback-Leibler divergence between the two, or performing moment matching in the case of a single Gaussian. A simpler alternative (which forms ‘standard’ EC) is to make the assumption  $p(h_{t+1}|s_t, s_{t+1}, v_{1:T}) \approx p(h_{t+1}|s_{t+1}, v_{1:T})$ , where  $p(h_{t+1}|s_{t+1}, v_{1:T})$  is already known from the previous backward recursion. Under this assumption, the recursion becomes

$$p(h_t, s_t|v_{1:T}) \approx \sum_{s_{t+1}} p(s_{t+1}|v_{1:T}) p(s_t|s_{t+1}, v_{1:T}) \langle p(h_t|h_{t+1}, s_t, s_{t+1}, v_{1:t}) \rangle_{p(h_{t+1}|s_{t+1}, v_{1:T})} \quad (10)$$

**Evaluating**  $\langle p(h_t|h_{t+1}, s_t, s_{t+1}, v_{1:t}) \rangle_{p(h_{t+1}|s_{t+1}, v_{1:T})}$

$\langle p(h_t|h_{t+1}, s_t, s_{t+1}, v_{1:t}) \rangle_{p(h_{t+1}|s_{t+1}, v_{1:T})}$  is a Gaussian in  $h_t$ , whose statistics we will now compute. First we find  $p(h_t|h_{t+1}, s_t, s_{t+1}, v_{1:t})$  which may be obtained from the joint distribution

$$p(h_t, h_{t+1}|s_t, s_{t+1}, v_{1:t}) = p(h_{t+1}|h_t, s_{t+1})p(h_t|s_t, v_{1:t}) \quad (11)$$

which itself can be found from a forward dynamics from the filtered estimate  $p(h_t|s_t, v_{1:t})$ . The statistics for the marginal  $p(h_t|s_t, s_{t+1}, v_{1:t})$  are simply those of  $p(h_t|s_t, v_{1:t})$ , since  $s_{t+1}$  carries no extra information about  $h_t$ . The remaining statistics are the mean of  $h_{t+1}$ , the covariance of  $h_{t+1}$  and cross-variance between  $h_t$  and  $h_{t+1}$ , which are given by

$$\langle h_{t+1} \rangle = A(s_{t+1})f_t(s_t), \quad \Sigma_{t+1,t+1} = A(s_{t+1})F_t(s_t)A^T(s_{t+1}) + \Sigma^h(s_{t+1}), \quad \Sigma_{t+1,t} = A(s_{t+1})F_t(s_t)$$

Given the statistics of (11), we may now condition on  $h_{t+1}$  to find  $p(h_t|h_{t+1}, s_t, s_{t+1}, v_{1:t})$ . Doing so effectively constitutes a reversal of the dynamics,

$$h_t = \overleftarrow{A}(s_t, s_{t+1})h_{t+1} + \overleftarrow{\eta}(s_t, s_{t+1})$$

where  $\overleftarrow{A}$  and  $\overleftarrow{\eta}(s_t, s_{t+1}) \sim \mathcal{N}(\overleftarrow{m}(s_t, s_{t+1}), \overleftarrow{\Sigma}(s_t, s_{t+1}))$  are easily found using conditioning. Averaging the above reversed dynamics over  $p(h_{t+1}|s_{t+1}, v_{1:T})$ , we find that  $\langle p(h_t|h_{t+1}, s_t, s_{t+1}, v_{1:t}) \rangle_{p(h_{t+1}|s_{t+1}, v_{1:T})}$  is a Gaussian with statistics

$$\mu_t = \overleftarrow{A}(s_t, s_{t+1})g(s_{t+1}) + \overleftarrow{m}(s_t, s_{t+1}), \quad \Sigma_{t,t} = \overleftarrow{A}(s_t, s_{t+1})G(s_{t+1})\overleftarrow{A}^T(s_t, s_{t+1}) + \overleftarrow{\Sigma}(s_t, s_{t+1})$$

These equations directly mirror the standard RTS backward pass[1].

**Evaluating**  $p(s_t|s_{t+1}, v_{1:T})$

The main departure of EC from previous methods is in treating the term

$$p(s_t|s_{t+1}, v_{1:T}) = \langle p(s_t|h_{t+1}, s_{t+1}, v_{1:t}) \rangle_{p(h_{t+1}|s_{t+1}, v_{1:T})} \quad (12)$$

The term  $p(s_t|h_{t+1}, s_{t+1}, v_{1:t})$  is given by

$$p(s_t|h_{t+1}, s_{t+1}, v_{1:t}) = \frac{p(h_{t+1}|s_{t+1}, s_t, v_{1:t})p(s_t, s_{t+1}|v_{1:t})}{\sum_{s'_t} p(h_{t+1}|s_{t+1}, s'_t, v_{1:t})p(s'_t, s_{t+1}|v_{1:t})} \quad (13)$$

Here  $p(s_t, s_{t+1}|v_{1:t}) = p(s_{t+1}|s_t, v_{1:t})p(s_t|v_{1:t})$ , where  $p(s_{t+1}|s_t, v_{1:t})$  occurs in the forward pass, (7). In (13),  $p(h_{t+1}|s_{t+1}, s_t, v_{1:t})$  is found by marginalising (11).

Computing the average of (13) with respect to  $p(h_{t+1}|s_{t+1}, v_{1:T})$  may be achieved by any numerical integration method desired. The simplest approximation is to evaluate the integrand at the mean value of the averaging distribution  $p(h_{t+1}|s_{t+1}, v_{1:T})$ . Otherwise, sampling from the Gaussian  $p(h_{t+1}|s_{t+1}, v_{1:T})$  has the advantage that covariance information is used<sup>2</sup>.

### Closing the Recursion

We have now computed both the continuous and discrete factors in (8), which we wish to use to write the smoothed estimate in the form  $p(h_t, s_t|v_{1:T}) = p(s_t|v_{1:T})p(h_t|s_t, v_{1:T})$ . The distribution  $p(h_t|s_t, v_{1:T})$  is readily obtained from the joint (8) by conditioning on  $s_t$  to form the mixture

$$p(h_t|s_t, v_{1:T}) = \sum_{s_{t+1}} p(s_{t+1}|s_t, v_{1:T})p(h_t|s_t, s_{t+1}, v_{1:T})$$

which may be collapsed to a  $J$ -component mixture of Gaussians. The smoothed posterior  $p(s_t|v_{1:T})$  is given by

$$p(s_t|v_{1:T}) = \sum_{s_{t+1}} p(s_{t+1}|v_{1:T}) \langle p(s_t|h_{t+1}, s_{t+1}, v_{1:t}) \rangle_{p(h_{t+1}|s_{t+1}, v_{1:T})}. \quad (14)$$

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<sup>2</sup>This is a form of exact sampling since drawing samples from a Gaussian is easy. This should not be confused with meaning that this use of sampling renders EC a sequential Monte-Carlo sampling scheme.

### 2.3 Relation to other methods

The standard-EC Backpass procedure is closely related to Kim’s method [8]. In both standard-EC and Kim’s method, the approximation  $p(h_{t+1}|s_t, s_{t+1}, v_{1:T}) \approx p(h_{t+1}|s_{t+1}, v_{1:T})$ , is used to form a numerically simple backward pass. The other ‘approximation’ in EC is to numerically compute the average in (14). In Kim’s method, however, an update for the discrete variables is formed by replacing the required term in (14) by

$$\langle p(s_t|h_{t+1}, s_{t+1}, v_{1:t}) \rangle_{p(h_{t+1}|s_{t+1}, v_{1:T})} \approx p(s_t|s_{t+1}, v_{1:t}) \quad (15)$$

This approximation decouples the discrete backward pass in Kim’s method from the continuous dynamics, since  $p(s_t|s_{t+1}, v_{1:t}) \propto p(s_{t+1}|s_t)p(s_t|v_{1:t})/p(s_{t+1}|v_{1:t})$  can be computed simply from the filtered results alone. The fundamental difference therefore between EC and Kim’s method is that the approximation, (15), is not required by EC. The EC backward pass therefore makes fuller use of the future information, resulting in a recursion which intimately couples the continuous and discrete variables. Unlike [8] and [4], where  $g_t, G_t \equiv f_t, F_t$  and only the backward pass mixture weights are updated from the forward pass, EC actually changes the Gaussian parameters  $g_t, G_t$  in a non-trivial way. The resulting effect on the quality of the approximation can be profound, as we will see in the experiments.

The Expectation Propagation algorithm makes the central assumption, as in EC, of collapsing the posteriors to a Gaussian family [5]. However, in EP, collapsing to a mixture of Gaussians is difficult – indeed, even working with a single Gaussian may be numerically unstable. In contrast, EC works largely with moment parameterisations of Gaussians, for which relatively few numerical difficulties arise. As explained in the derivation of (10), the conditional independence assumption  $p(h_{t+1}|s_t, s_{t+1}, v_{1:T}) \approx p(h_{t+1}|s_{t+1}, v_{1:T})$  is not strictly necessary in EC. We motivate it by computational simplicity, since finding an appropriate moment matching approximation of  $p(h_{t+1}|s_t, s_{t+1}, v_{1:T})$  in (9) requires a relatively expensive non-Gaussian integration. The important point here is that, if we did treat  $p(h_{t+1}|s_t, s_{t+1}, v_{1:T})$  more correctly, the only assumption in EC would be a collapse to a mixture of Gaussians, as in EP. As a point of interest, as in EC, the exact computation requires only a single forward and backward pass, whilst EP is an ‘open’ procedure requiring iteration to convergence.

In [9] a related dynamics reversed is proposed. However, the singularities resulting from incorrectly treating  $p(v_{t+1:T}|h_t, s_t)$  as a density are heuristically finessed.

In [10] a variational method approximates the joint distribution  $p(h_{1:T}, s_{1:T}|v_{1:T})$  rather than the marginal inference  $p(h_t, s_t|v_{1:T})$ . This is a disadvantage when compared to other methods that directly approximate the marginal.

Sequential Monte Carlo methods (Particle Filters)[11], are essentially mixture of delta-function approximations. Whilst potentially powerful, these typically suffer in high-dimensional hidden spaces, unless techniques such as Rao-Blackwellisation are performed. ADF is generally preferential to Particle Filtering since in ADF the approximation is a mixture of non-trivial distributions, and is therefore better at capturing the variability of the posterior.

## 3 Demonstration

Testing EC in a problem with a reasonably long temporal sequence,  $T$ , is important since numerical instabilities may not be apparent in timeseries of just a few points. To do this, we sequentially generate hidden and visible states from a given model, here with  $H = 3, S = 2, V = 1$  – see Figure(2) for full details of the experimental setup. Then, given only the parameters of the model and the visible observations (but not any of the hidden states  $h_{1:T}, s_{1:T}$ ), the task is to infer  $p(h_t|s_t, v_{1:T})$  and  $p(s_t|v_{1:T})$ . Since the exact computation is exponential in  $T$ , a simple alternative is to assume that the original sample states  $s_{1:T}$  are the ‘correct’ inferences, and compare how our most probable posterior smoothed estimates  $\arg \max_{s_t} p(s_t|v_{1:T})$  compare with the assumed correct sample  $s_t$ . We chose conditions that, from the viewpoint of classical signal processing, are difficult, with changes in the switches occurring at a much higher rate than the typical frequencies in the signal  $v_t$ .

For EC we use the mean approximation for the numerical integration of (12). We included the Particle Filter merely for a point of comparison with ADF, since they are not designed to approximate

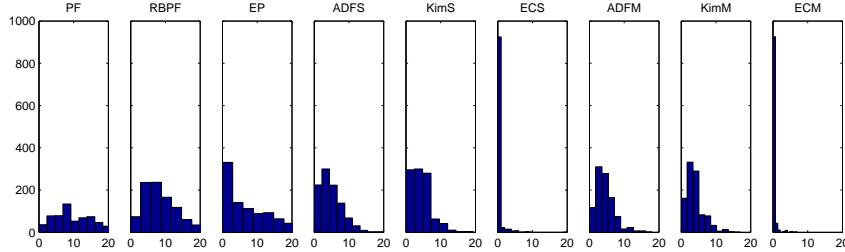


Figure 2: The number of errors in estimating  $p(s_t|v_{1:T})$  for a binary switch ( $S = 2$ ) over a time series of length  $T = 100$ . Hence 50 errors corresponds to random guessing. Plotted are histograms of the errors are over 1000 experiments. The  $x$ -axes are cut off at 20 errors to improve visualisation of the results. (PF) Particle Filter. (RBPF) Rao-Blackwellised PF. (EP) Expectation Propagation. (ADFS) Assumed Density Filtering using a Single Gaussian. (KimS) Kim’s smoother using the results from ADFS. (ECS) Expectation Correction using a Single Gaussian ( $I = J = 1$ ). (ADFM) ADF using a multiple of  $I = 4$  Gaussians. (KimM) Kim’s smoother using the results from ADFM. (ECM) Expectation Correction using a mixture with  $I = J = 4$  components.  $S = 2, V = 1$  (scalar observations),  $T = 100$ , with zero output bias.  $A(s) = 0.9999 * \text{orth}(\text{randn}(H, H))$ ,  $B(s) = \text{randn}(V, H)$ .  $H = 3$ ,  $\Sigma^h(s) = I_H$ ,  $\Sigma^v(s) = 0.1I_V$ ,  $p(s_{t+1}|s_t) \propto 1_{S \times S} + I_S$ . At time  $t = 1$ , the priors are  $p_1 = \text{uniform}$ , with  $h_1$  drawn from  $\mathcal{N}(10 * \text{randn}(H, 1), I_H)$ .

the smoothed estimate, for which 1000 particles were used, with Kitagawa resampling. For the Rao-Blackwellised Particle Filter [11], 500 particles were used, with Kitagawa resampling. We found that EP<sup>3</sup> was numerically unstable and often struggled to converge. To encourage convergence, we used the damping method in [12], performing 20 iterations with a damping factor of 0.5. Nevertheless, the disappointing performance of EP is most likely due to conflicts resulting from numerical instabilities introduced by the frequent conversions between moment and canonical representations.

The best filtered results are given using ADF, since this is better able to represent the variance in the filtered posterior than the sampling methods. Unlike Kim’s method, EC makes good use of the future information to clean up the filtered results considerably. One should bear in mind that both EC and Kim’s method use the same ADF filtered results. These demonstrates that EC may dramatically improve on Kim’s method, so that the small amount of extra work in making a numerical approximation of  $p(s_t|s_{t+1}, v_{1:T})$ , (12), may bring significant benefits. We found similar conclusions for experiments with an aSLDS[7].

## 4 Application to Noise Robust ASR

Here we briefly present an application of the SLDS to robust Automatic Speech Recognition (ASR), for which the intractable inference is performed by EC, and serves to demonstrate how EC scales well to a large-scale application. Fuller details are given in [13]. The standard approach to noise robust ASR is to provide a set of noise-robust features to a standard Hidden Markov Model (HMM) classifier, which is based on modeling the acoustic feature vector. For example, the method of Unsupervised Spectral Subtraction (USS) [14] provides state-of-the-art performance in this respect. Incorporating noise models directly into such feature-based HMM systems is difficult, mainly because the explicit influence of the noise on the features is poorly understood. An alternative is to model the raw speech signal directly, such as the SAR-HMM model [15] for which, under *clean* conditions, isolated spoken digit recognition performs well. However, the SAR-HMM performs poorly under noisy conditions, since no explicit noise processes are taken into account by the model.

The approach we take here is to extend the SAR-HMM to include an explicit noise process, so that the observed signal  $v_t$  is modeled as a noise corrupted version of a clean *hidden* signal  $v_t^h$ :

$$v_t = v_t^h + \tilde{\eta}_t \quad \text{with} \quad \tilde{\eta}_t \sim \mathcal{N}(0, \tilde{\sigma}^2)$$

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<sup>3</sup>Generalised EP [5], which groups variables together improves on the results, but is still far inferior to the EC results presented here – Onno Zoeter personal communication.

Noise Variance	SNR (dB)	HMM	SAR-HMM	AR-SLDS
0	26.5	100.0%	97.0%	96.8%
$10^{-7}$	26.3	100.0%	79.8%	96.8%
$10^{-6}$	25.1	90.9%	56.7%	96.4%
$10^{-5}$	19.7	86.4%	22.2%	94.8%
$10^{-4}$	10.6	59.1%	9.7%	84.0%
$10^{-3}$	0.7	9.1%	9.1%	61.2%

Table 1: Comparison of the recognition accuracy of three models when the test utterances are corrupted by various levels of Gaussian noise.

The dynamics of the clean signal is modeled by a switching AR process

$$v_t^h = \sum_{r=1}^R c_r(s_t) v_{t-r}^h + \eta_t^h(s_t), \quad \eta_t^h(s_t) \sim \mathcal{N}(0, \sigma^2(s_t))$$

where  $s_t \in \{1, \dots, S\}$  denotes which of a set of AR coefficients  $c_r(s_t)$  are to be used at time  $t$ , and  $\eta_t^h(s_t)$  is the so-called *innovation* noise. When  $\sigma^2(s_t) \equiv 0$ , this model reproduces the SAR-HMM of [15], a specially constrained HMM. Hence inference and learning for the SAR-HMM are tractable and straightforward.

For the case  $\sigma^2(s_t) \geq 0$  the model can be recast as an SLDS. To do this we define  $h_t$  as a vector which contains the  $R$  most recent clean hidden samples

$$h_t = [v_t^h \quad \dots \quad v_{t-r+1}^h]^\top \quad (16)$$

and we set  $A(s_t)$  to be a  $R \times R$  matrix where the first row contains the AR coefficients  $-c_r(s_t)$  and the rest is a shifted down identity matrix. For example, for a third order ( $R = 3$ ) AR process,

$$A(s_t) = \begin{bmatrix} -c_1(s_t) & -c_2(s_t) & -c_3(s_t) \\ 1 & 0 & 0 \\ 0 & 1 & 0 \end{bmatrix}. \quad (17)$$

The hidden covariance matrix  $\Sigma_h(s)$  has all elements are zero, except the top-left most which is set to the innovation variance. To extract the first component of  $h_t$  we use the (switch independent)  $1 \times R$  projection matrix  $B = [1 \ 0 \ \dots \ 0]$ . The (switch independent) visible scalar noise variance is given by  $\Sigma_v \equiv \sigma_v^2$ .

A well-known issue with raw speech signal models is that the energy of a signal may vary from one speaker to another or because of a change in recording conditions. For this reason the innovation  $\Sigma_h$  is adjusted by maximising the likelihood of an observed sequence with respect to the innovation covariance, a process called *Gain Adaptation* [15].

#### 4.1 Training & Evaluation

Following [15], we trained a separate SAR-HMM for each of the eleven digits (0–9 and ‘oh’) from the TI-DIGITS database [16]. The training set for each digit was composed of 110 single digit utterances down-sampled to 8 kHz, each one pronounced by a male speaker. Each SAR-HMM was composed of ten states with a left-right transition matrix. Each state was associated with a 10th-order AR process and the model was constrained to stay an integer multiple of  $K = 140$  time steps (0.0175 seconds) in the same state. We refer the reader to [15] for a detailed explanation of the training procedure used with the SAR-HMM.

An AR-SLDS was built for each of the eleven digits by copying the parameters of the corresponding trained SAR-HMM, i.e., the AR coefficients  $c_r(s)$  are copied into the first row of the hidden transition matrix  $A(s)$  and the same discrete transition distribution  $p(s_t | s_{t-1})$  is used. The models were then evaluated on a test set composed of 112 corrupted utterances of each of the eleven digits, each pronounced by different male speakers than those used in the training set.

The recognition accuracy obtained by the models on the corrupted test sets is presented in Table 1. As expected, the performance of the SAR-HMM rapidly decreases with noise. Thanks to USS,

the feature-based HMM has above 90% accuracy as long as the SNR is bigger than 20 dB. In contrast, the AR-SLDS achieves a recognition accuracy of 61.2% at a SNR close to 0 dB, while the performance of the two other methods is equivalent to random guessing (9.1%).

Whilst other inference methods may also perform well in this case, we found that EC performs admirably, without numerical instabilities, even for time-series with several thousand time-steps.

## 5 Discussion

We presented a method for approximate smoothed inference in an augmented class of switching linear dynamical systems. Our approximation is based on the idea that due to the forgetting which commonly occurs in Markovian models, a finite number of mixture components may provide a reasonable approximation. Clearly, in systems with very long correlation times our method may require too many mixture components to produce a satisfactory result, although we are unaware of other techniques that would be able to cope well in that case. The main benefit of EC over Kim smoothing is that future information is more accurately dealt with. Whilst EC is not as general as EP, EC carefully exploits the properties of singly-connected distributions, such as the aSLDS to provide a relatively numerically stable procedure. The relaxed version of EC makes the same basic assumptions as EP (in singly-connected distributions), but results only in a single forward and backward pass, each being based on a stable update procedure. We have successfully applied EC to a problem in automatic speech recognition where we model a one dimensional speech signal using a SLDS [13]. The signal consists of many thousands of timepoints, and numerical stability is an important concern. We hope that the ideas presented here may therefore help facilitate the practical application of dynamic hybrid networks. Code for EC is available from \*.

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