
Unified Inference for Variational Bayesian Linear Gaussian State-Space Models

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Abstract

Linear Gaussian State-Space Models are widely used and a Bayesian treatment of parameters is therefore of considerable interest. The approximate Variational Bayesian method applied to these models is an attractive approach, used successfully in applications ranging from acoustics to bioinformatics. The most challenging aspect of implementing the method is in performing inference on the hidden state sequence of the model. We show how to convert the inference problem so that standard and stable Kalman Filtering/Smoothing recursions from the literature may be applied. This is in contrast to previously published approaches based on Belief Propagation. Our framework both simplifies and unifies the inference problem, so that future applications may be easily developed. We demonstrate the elegance of the approach on Bayesian temporal ICA, with an application to finding independent components in noisy EEG signals.

1 Linear Gaussian State-Space Models

Linear Gaussian State-Space Models (LGSSMs)¹ are fundamental in time-series analysis [1, 2, 3]. In these models the observations $v_{1:T}$ ² are generated from an underlying dynamical system on $h_{1:T}$ according to

$$v_t = Bh_t + \eta_t^v, \quad \eta_t^v \sim \mathcal{N}(\mathbf{0}_V, \Sigma_V); \quad h_t = Ah_{t-1} + \eta_t^h, \quad \eta_t^h \sim \mathcal{N}(\mathbf{0}_H, \Sigma_H),$$

where $\mathcal{N}(\mu, \Sigma)$ denotes a Gaussian with mean μ and covariance Σ , and $\mathbf{0}_X$ denotes an X -dimensional zero vector. The observation v_t has dimension V and the hidden state h_t dimension H . Probabilistically, the LGSSM is defined by:

$$p(v_{1:T}, h_{1:T} | \Theta) = p(v_1 | h_1) p(h_1) \prod_{t=2}^T p(v_t | h_t) p(h_t | h_{t-1}),$$

with $p(v_t | h_t) = \mathcal{N}(Bh_t, \Sigma_V)$, $p(h_t | h_{t-1}) = \mathcal{N}(Ah_{t-1}, \Sigma_H)$, $p(h_1) = \mathcal{N}(\mu, \Sigma)$ and where $\Theta = \{A, B, \Sigma_H, \Sigma_V, \mu, \Sigma\}$ denotes the model parameters. Because of the widespread use of these models, a Bayesian treatment of parameters is of considerable interest [4, 5, 6, 7, 8].

An exact implementation of the Bayesian LGSSM is formally intractable [8], and recently a Variational Bayesian (VB) approximation has been studied [4, 5, 6, 7, 9]. The most challenging part of implementing the VB method is performing inference over $h_{1:T}$, and previous authors have developed their own specialized routines, based on Belief Propagation, since standard LGSSM inference routines appear, at first sight, not to be applicable.

¹Also called Kalman Filters/Smoothers, Linear Dynamical Systems.

² $v_{1:T}$ denotes v_1, \dots, v_T .

A key contribution of this paper is to show how the Variational Bayesian treatment of the LGSSM *can* be implemented using standard inference routines. Based on the insight we provide, any standard inference method may be applied, including those specifically addressed to improve numerical stability [10, 11, 2]. In this article, we decided to describe the standard predictor-corrector and Rauch-Tung-Striebel recursions [2], and also suggest a small modification that reduces computational cost.

The Bayesian LGSSM is particularly of interest when strong prior constraints are needed to find adequate solutions. One such case is in EEG signal analysis, whereby we wish to extract sources that evolve independently through time. Since EEG is particularly noisy [12], a prior that encourages sources to have preferential spectral properties is advantageous in recovering meaningful sources. This application is discussed in Section 4, and demonstrates the ease of applying our VB framework.

2 Bayesian Linear Gaussian State-Space Models

In the Bayesian treatment of the LGSSM, instead of considering the model parameters Θ as fixed, we define a prior distribution $p(\Theta|\hat{\Theta})$, where $\hat{\Theta}$ is a set of hyperparameters. Then:

$$p(v_{1:T}|\hat{\Theta}) = \int_{\Theta} p(v_{1:T}|\hat{\Theta}, \Theta)p(\Theta|\hat{\Theta}). \quad (1)$$

In a full Bayesian treatment we would define additional prior distributions over the hyperparameters $\hat{\Theta}$. Here we take instead the ML-II ('evidence') framework, in which the optimal set of hyperparameters is found by maximizing $p(v_{1:T}|\hat{\Theta})$ with respect to $\hat{\Theta}$ [6, 7, 9].

For the parameter priors, we define Gaussians on the columns of A and B :

$$p(A|\alpha, \Sigma_H) \propto \prod_{j=1}^H e^{-\frac{\alpha_j}{2}(A_j - \hat{A}_j)^\top \Sigma_H^{-1}(A_j - \hat{A}_j)}, \quad p(B|\beta, \Sigma_V) \propto \prod_{j=1}^H e^{-\frac{\beta_j}{2}(B_j - \hat{B}_j)^\top \Sigma_V^{-1}(B_j - \hat{B}_j)},$$

which has the effect of biasing the transition and emission matrices to desired forms \hat{A} and \hat{B} . The conjugate priors for the covariances Σ_H and Σ_V are Inverse Wishart distributions [7]³. In the simpler and more common case of assuming diagonal covariances these become Inverse Gamma distributions [7, 5]. The hyperparameters are then $\hat{\Theta} = \{\alpha, \beta\}$ ⁴.

Variational Bayes

Optimizing Eq. (1) with respect to $\hat{\Theta}$ is difficult due to the intractability of the integrals. Instead, in VB, one considers the lower bound [6, 7, 9]⁵:

$$\mathcal{L} = \log p(v_{1:T}|\hat{\Theta}) \geq H_q(\Theta, h_{1:T}) + \left\langle \log p(\Theta|\hat{\Theta}) \right\rangle_{q(\Theta)} + \left\langle E(h_{1:T}, \Theta|\hat{\Theta}) \right\rangle_{q(\Theta, h_{1:T})} \equiv \mathcal{F},$$

where

$$E(h_{1:T}, \Theta|\hat{\Theta}) \equiv \log p(v_{1:T}, h_{1:T}|\Theta, \hat{\Theta}).$$

The notation $H_d(x)$ signifies the entropy of the distribution $d(x)$, and $\langle \cdot \rangle_{d(x)}$ denotes the expectation operator.

The key approximation in VB is $q(\Theta, h_{1:T}) \equiv q(\Theta)q(h_{1:T})$, from which one may show that, for optimality of \mathcal{F} ,

$$q(h_{1:T}) \propto e^{\langle E(h_{1:T}, \Theta|\hat{\Theta}) \rangle_{q(\Theta)}}, \quad q(\Theta) \propto p(\Theta) e^{\langle E(h_{1:T}, \Theta|\hat{\Theta}) \rangle_{q(h_{1:T})}}.$$

These coupled equations need to be iterated to convergence. The updates for the parameters $q(\Theta)$ are straightforward and are given in Appendices A and B. Once converged, the hyperparameters are updated by maximizing \mathcal{F} with respect to $\hat{\Theta}$, which lead to simple update formulae [7].

Our main concern is with the update for $q(h_{1:T})$, for which this paper makes a departure from treatments previously presented.

³For expositional simplicity, we do not put priors on μ and Σ .

⁴For simplicity, we keep the parameters of the Inverse Wishart priors fixed.

⁵Strictly we should write throughout $q(\cdot|v_{1:T})$. We omit the dependence on $v_{1:T}$ for notational convenience.

3 Unified Inference on $q(h_{1:T})$

Optimally $q(h_{1:T})$ is Gaussian since $\langle E(h_{1:T}, \Theta | \hat{\Theta}) \rangle_{q(\Theta)}$ is quadratic in $h_{1:T}$, being namely⁶

$$-\frac{1}{2} \sum_{t=1}^T \left[\langle (v_t - Bh_t)^\top \Sigma_V^{-1} (v_t - Bh_t) \rangle_{q(B, \Sigma_V)} + \langle (h_t - Ah_{t-1})^\top \Sigma_H^{-1} (h_t - Ah_{t-1}) \rangle_{q(A, \Sigma_H)} \right]. \quad (2)$$

Optimally, $q(A|\Sigma_H)$ and $q(B|\Sigma_V)$ are Gaussians (see Appendix A), so we can easily carry out the averages. The further averages over $q(\Sigma_H)$ and $q(\Sigma_V)$ are also easy due to conjugacy. Whilst this defines the distribution $q(h_{1:T})$, quantities such as $q(h_t)$, which are required for the parameter updates (see the Appendices), need to be inferred from this distribution. Clearly, in the non-Bayesian case, the averages over the parameters are not present, and the above simply represents an LGSSM whose visible variables have been clamped into their evidential states. In that case, inference can be performed using any standard method. Our aim, therefore, is to try to represent the averaged Eq. (2) directly as an LGSSM $\tilde{q}(h_{1:T}|\tilde{v}_{1:T})$, for some suitable parameter settings.

Mean + Fluctuation Decomposition

A useful decomposition is to write

$$\langle (v_t - Bh_t)^\top \Sigma_V^{-1} (v_t - Bh_t) \rangle_{q(B, \Sigma_V)} = \underbrace{(v_t - \langle B \rangle h_t)^\top \langle \Sigma_V^{-1} \rangle (v_t - \langle B \rangle h_t)}_{\text{mean}} + \underbrace{h_t^\top S_B h_t}_{\text{fluctuation}},$$

and similarly

$$\langle (h_t - Ah_{t-1})^\top \Sigma_H^{-1} (h_t - Ah_{t-1}) \rangle_{q(A, \Sigma_H)} = \underbrace{(h_t - \langle A \rangle h_{t-1})^\top \langle \Sigma_H^{-1} \rangle (h_t - \langle A \rangle h_{t-1})}_{\text{mean}} + \underbrace{h_{t-1}^\top S_A h_{t-1}}_{\text{fluctuation}},$$

where the parameter covariances are $S_B = V H_B^{-1}$ and $S_A = H H_A^{-1}$ (see Appendix A). The mean terms simply represent a clamped LGSSM with averaged parameters. However, the extra contributions from the fluctuations mean that Eq. (2) cannot be written as a clamped LGSSM with averaged parameters. In order to deal with these extra terms, our idea is to treat the fluctuations as arising from an augmented visible variable, for which Eq. (2) can then be considered as a clamped LGSSM.

Inference Using an Augmented LGSSM

To represent Eq. (2) as a LGSSM $\tilde{q}(h_{1:T}|\tilde{v}_{1:T})$, we augment v_t and B as⁷:

$$\tilde{v}_t = \text{vert}(v_t, \mathbf{0}_H, \mathbf{0}_H), \quad \tilde{B} = \text{vert}(\langle B \rangle, U_A, U_B),$$

where U_A is the Cholesky decomposition of S_A , so that $U_A^\top U_A = S_A$. Similarly, U_B is the Cholesky decomposition of S_B . The equivalent LGSSM $\tilde{q}(h_{1:T}|\tilde{v}_{1:T})$ is then completed by specifying⁸

$$\tilde{A} \equiv \langle A \rangle, \quad \tilde{\Sigma}_H \equiv \langle \Sigma_H^{-1} \rangle^{-1}, \quad \tilde{\Sigma}_V \equiv \text{diag}(\langle \Sigma_V^{-1} \rangle^{-1}, I_H, I_H), \quad \tilde{\mu} \equiv \mu, \quad \tilde{\Sigma} \equiv \Sigma.$$

The validity of this parameter assignment can be checked by showing that, up to negligible constants, the exponent of this augmented LGSSM has the same form as Eq. (2). Now that this has been written as an LGSSM $\tilde{q}(h_{1:T}|\tilde{v}_{1:T})$, standard inference routines in the literature may be applied to compute $q(h_t) = \tilde{q}(h_t|\tilde{v}_{1:T})$ [1, 11, 2]⁹.

For completeness, we decided to describe the standard predictor-corrector form of a Kalman filter, together with the Rauch-Tung-Striebel recursions [2] for performing inference in an LGSSM. These

⁶For simplicity of exposition, we ignore the first time-point here.

⁷The notation $\text{vert}(x_1, \dots, x_n)$ stands for vertically concatenating the arguments x_1, \dots, x_n .

⁸Strictly, we need a time-dependent emission $\tilde{B}_t = \tilde{B}$, for $t = 1, \dots, T-1$. For time T , \tilde{B}_T has the Cholesky factor U_A replaced by $\mathbf{0}_{H,H}$.

⁹Note that, since the augmented LGSSM $\tilde{q}(h_{1:T}|\tilde{v}_{1:T})$ is designed to match the *fully* clamped distribution $q(h_{1:T})$, filtering $\tilde{q}(h_{1:T}|\tilde{v}_{1:T})$ does not correspond to filtering $q(h_{1:T})$.

Algorithm 1 LGSSM: Forward and backward recursive updates. The smoothed posterior $p(h_t|v_{1:T})$ is returned in the mean \hat{h}_t^T and covariance P_t^T .

procedure FORWARD

1a: $P \leftarrow \Sigma$

1b: $P \leftarrow D\Sigma$, where $D \equiv I - \Sigma U_{AB} (I + U_{AB}^T \Sigma U_{AB})^{-1} U_{AB}^T$

2a: $\hat{h}_1^0 \leftarrow \mu$

2b: $\hat{h}_1^0 \leftarrow D\mu$

3: $K \leftarrow PB^T(BPB^T + \Sigma_V)^{-1}$, $P_1^1 \leftarrow (I - KB)P$, $\hat{h}_1^1 \leftarrow \hat{h}_1^0 + K(v_1 - B\hat{h}_1^0)$

for $t \leftarrow 2, T$ **do**

4: $P_t^{t-1} \leftarrow AP_{t-1}^{t-1}A^T + \Sigma_H$

5a: $P \leftarrow P_t^{t-1}$

5b: $P \leftarrow D_t P_t^{t-1}$, where $D_t \equiv I - P_t^{t-1} U_{AB} (I + U_{AB}^T P_t^{t-1} U_{AB})^{-1} U_{AB}^T$

6a: $\hat{h}_t^{t-1} \leftarrow A\hat{h}_{t-1}^{t-1}$

6b: $\hat{h}_t^{t-1} \leftarrow D_t A\hat{h}_{t-1}^{t-1}$

7: $K \leftarrow PB^T(BPB^T + \Sigma_V)^{-1}$, $P_t^t \leftarrow (I - KB)P$, $\hat{h}_t^t \leftarrow \hat{h}_t^{t-1} + K(v_t - B\hat{h}_t^{t-1})$

end for

end procedure

procedure BACKWARD

for $t \leftarrow T - 1, 1$ **do**

$\overleftarrow{A}_t \leftarrow P_t^t A^T (P_{t+1}^t)^{-1}$

$P_t^T \leftarrow P_t^t + \overleftarrow{A}_t (P_{t+1}^T - P_{t+1}^t) \overleftarrow{A}_t^T$

$\hat{h}_t^T \leftarrow \hat{h}_t^t + \overleftarrow{A}_t (\hat{h}_{t+1}^T - A\hat{h}_t^t)$

end for

end procedure

are given in Algorithm 1. To compute $\tilde{q}(h_t|\tilde{v}_{1:T})$, we then call the FORWARD and BACKWARD procedures.

We present two variants of the FORWARD pass. Either we may call procedure FORWARD in Algorithm 1 with parameters $\tilde{A}, \tilde{B}, \tilde{\Sigma}_H, \tilde{\Sigma}_V, \tilde{\mu}, \tilde{\Sigma}$ and the augmented visible variables \tilde{v}_t in which we use steps 1a, 2a, 5a and 6a. This is exactly the predictor-corrector form of a Kalman filter [2]. Otherwise, in order to reduce the computational cost, we may call procedure FORWARD with the parameters $\langle A \rangle, \langle B \rangle, \langle \Sigma_H^{-1} \rangle^{-1}, \langle \Sigma_V^{-1} \rangle^{-1}, \mu, \Sigma$ and the original visible variable v_t in which we use steps 1b (where $U_{AB}^T U_{AB} \equiv S_A + S_B$), 2b, 5b and 6b. The two algorithms are mathematically equivalent. Computing $q(h_t) = \tilde{q}(h_t|\tilde{v}_{1:T})$ is then completed by calling the common BACKWARD pass¹⁰.

The important point here is that the reader may supply any standard Kalman Filtering/Smoothing routine, and simply call it with the appropriate parameters. In some parameter regimes, or in very long time series, numerical stability may be a serious concern, for which several stabilized algorithms have been developed over the years, for example the square-root forms [10, 11, 2]. By converting the problem to a standard form, we have therefore unified and simplified inference, so that future applications may be more readily developed.

¹⁰The cross-moment required for learning in Section 4 can be easily computed using:

$$\left\langle h_{t-1} h_t^T \right\rangle_{p(h_{t-1:t}|v_{1:T})} = \overleftarrow{A}_{t-1} P_t^T + \hat{h}_{t-1}^T (\hat{h}_t^T)^T.$$

This is much simpler than formulae surprisingly continued in the literature [3, 13].

3.1 Relation to Previous Approaches

An alternative approach to the one above, and taken in [7, 5], is to recognize that the posterior is

$$\log q(h_{1:T}) = \sum_{t=2}^T \phi_t(h_{t-1}, h_t) + \text{const.}$$

for suitably defined quadratic forms $\phi_t(h_{t-1}, h_t)$. Here the potentials $\phi_t(h_{t-1}, h_t)$ encode the averaging over the parameters A, B, Σ_H, Σ_V . The approach taken in [7] is to recognize this as a pairwise Markov chain, for which the Belief Propagation recursions may be applied. The backward pass from Belief Propagation makes use of the observations $v_{1:T}$, so that any approximate online treatment would be difficult. The approach in [5] is based on a Kullback-Leibler minimization of the posterior with a chain structure, which is algorithmically equivalent to Belief Propagation. Whilst mathematically valid procedures, the resulting algorithms do not correspond to any of the standard forms in the Kalman Filtering/Smoothing literature, whose properties have been well studied [14].

A stated aim in [7] is to find a sequential form for smoothing, since this has potential advantages in online situations, whereby high-dimensional observations can be discarded once they have been filtered. Our algorithm provides exactly this sequential form for smoothing.

4 An Application to Bayesian ICA

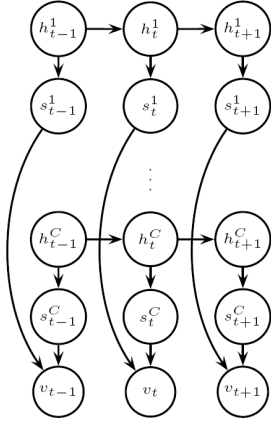


Figure 1: The structure of the LGSSM for ICA.

Σ_H and Σ_V which are instead considered as hyperparameters. More details of the model are given in [15]. The constraint $B = WP$ requires a minor modification from Section 3, as we discuss below.

Inference on $q(h_{1:T})$

A small modification of the mean + fluctuation decomposition for B occurs, namely:

$$\langle (v_t - Bh_t)^\top \Sigma_V^{-1} (v_t - Bh_t) \rangle_{q(W)} = (v_t - \langle B \rangle h_t)^\top \Sigma_V^{-1} (v_t - \langle B \rangle h_t) + h_t^\top P^\top S_W P h_t,$$

where $\langle B \rangle \equiv \langle W \rangle P$ and $S_W = V H_W^{-1}$. The quantities $\langle W \rangle$ and H_W are obtained as in Appendix A.1 with the replacement $h_t \leftarrow P h_t$. To represent the above as a LGSSM, we augment v_t and B as

$$\tilde{v}_t = \text{vert}(v_t, \mathbf{0}_H, \mathbf{0}_C), \quad \tilde{B} = \text{vert}(\langle B \rangle, U_A, U_W P),$$

where U_W is the Cholesky decomposition of S_W . The equivalent LGSSM is then completed by specifying $\tilde{A} \equiv \langle A \rangle$, $\tilde{\Sigma}_H \equiv \Sigma_H$, $\tilde{\Sigma}_V \equiv \text{diag}(\Sigma_V, I_H, I_C)$, $\tilde{\mu} \equiv \mu$, $\tilde{\Sigma} \equiv \Sigma$, and inference for $q(h_{1:T})$ performed using Algorithm 1. This demonstrates the elegance and unity of the approach in Section 3, since no new algorithm needs to be developed to perform inference, even in this special constrained parameter case.

A particular case for which the Bayesian LGSSM is of interest is in extracting independent source signals underlying a multivariate time-series [15, 5]. This will demonstrate how the approach developed in Section 3 makes VB easily to apply. The sources s^i are modeled as independent in the following sense:

$$p(s_{1:T}^i, s_{1:T}^j) = p(s_{1:T}^i) p(s_{1:T}^j), \quad \text{for } i \neq j, \quad i, j = 1, \dots, C.$$

Independence implies block diagonal transition and state noise matrices A , Σ_H and Σ , where each block c has dimension H_c . A one dimensional source s_t^c for each independent dynamical subsystem is then formed from $s_t^c = \mathbf{1}_c^\top h_t^c$, where $\mathbf{1}_c$ is a unit vector and h_t^c is the state of dynamical system c . Combining the sources, we can write $s_t = P h_t$, where $P = \text{diag}(\mathbf{1}_1^\top, \dots, \mathbf{1}_C^\top)$, $h_t = \text{vert}(h_t^1, \dots, h_t^C)$. The resulting emission matrix is constrained to be of the form $B = WP$, where W is the $V \times C$ mixing matrix. This means that the observations are formed from linearly mixing the sources, $v_t = W s_t + \eta_t^v$. The graphical structure of this model is presented in Fig 1. To encourage redundant components to be removed, we place a zero mean Gaussian prior on W . In this case, we do not define a prior for the parameters

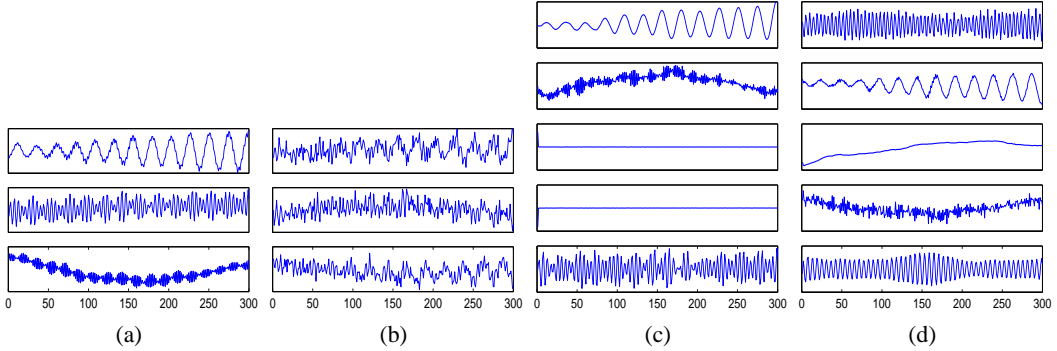


Figure 2: (a) Original sources s_t . (b) Observations resulting from mixing the original sources, $v_t = Ws_t + \eta_t^v$, $\eta_t^v \sim \mathcal{N}(0, I)$. (c) Recovered sources using the Bayesian LGSSM. (d) Sources found with MAP LGSSM.

4.1 Demonstration

As a simple demonstration, we used a LGSSM to generate 3 sources s_t^c with random 5×5 transition matrices A^c , $\mu = \mathbf{0}_H$ and $\Sigma \equiv \Sigma_H \equiv I_H$. The sources were mixed into three observations $v_t = Ws_t + \eta_t^v$, for W chosen with elements from a zero mean unit variance Gaussian distribution, and $\Sigma_V = I_V$. We then trained a Bayesian LGSSM with 5 sources and 7×7 transition matrices A^c . To bias the model to find the simplest sources, we used $\hat{A}^c \equiv \mathbf{0}_{H_c, H_c}$ for all sources. In Fig2a and Fig 2b we see the original sources and the noisy observations respectively. In Fig2c we see the estimated sources from our method after convergence of the hyperparameter updates. Two of the 5 sources have been removed, and the remaining three are a reasonable estimation of the original sources. Another possible approach for introducing prior knowledge is to use a Maximum a Posteriori (MAP) procedure by adding a prior term to the original log-likelihood $\log p(v_{1:T}|A, W, \Theta) + \log p(A|\alpha) + \log p(W|\beta)$. However, it is not clear how to reliably find the hyperparameters α and β in this case. One solution is to estimate them by optimizing the new objective function jointly with respect to the parameters and hyperparameters (this is the so-called joint map estimation – see for example [16]). A typical result of using this joint MAP approach on the artificial data is presented in Fig2d. The joint MAP does not estimate the hyperparameters well, and the incorrect number of sources is identified.

4.2 Application to EEG Analysis

In Fig 3a we plot three seconds of EEG data recorded from 4 channels (located in the right hemisphere) while a subject is performing imagined movement of the right hand. As is typical in EEG, each channel shows drift terms below 1 Hz which correspond to artifacts of the instrumentation, together with the presence of 50 Hz mains contamination and masks the rhythmical activity related to the mental task, mainly centered at 10 and 20 Hz [17]. We would therefore like a method which enables us to extract components in these information-rich 10 and 20 Hz frequency bands. Standard ICA methods such as FastICA do not find satisfactory sources based on raw ‘noisy’ data, and preprocessing with band-pass filters is usually required. Additionally, in EEG research, flexibility in the number of recovered sources is important since there may be many independent oscillators of interest underlying the observations and we would like some way to automatically determine their effective number. To preferentially find sources at particular frequencies, we specified a block diagonal matrix \hat{A}^c for each source c , where each block is a 2×2 rotation matrix at the desired frequency. We defined the following 16 groups of frequencies: [0.5], [0.5], [0.5], [0.5]; [10,11], [10,11], [10,11], [10,11]; [20,21], [20,21], [20,21], [20,21]; [50], [50], [50], [50]. The temporal evolution of the sources obtained after training the Bayesian LGSSM is given in Fig 3(b,c,d,e) (grouped by frequency range). The Bayes LGSSM removed 4 unnecessary sources from the mixing matrix W , that is one [10,11] Hz and three [20,21] Hz sources. The first 4 sources contain dominant low frequency drift, source 5, 6 and 8 contain [10,11] Hz, while source 10 contains [20,21] Hz centered activity. Of the 4 sources initialized to 50 Hz, only 2 retained 50 Hz activity, while the A^c of the other two have changed to model other frequencies present in the EEG. This method demonstrates the usefulness and applicability of the VB method in a real-world situation.

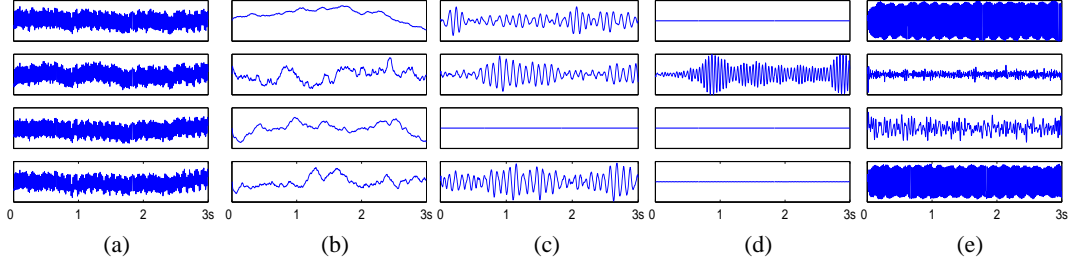


Figure 3: (a) Original raw EEG recordings from 4 channels. (b-e) 16 sources s_t estimated by the Bayesian LGSSM.

5 Conclusion

We considered the application of Variational Bayesian learning to Linear Gaussian State-Space Models. This is an important class of models with widespread application, and finding a simple way to implement this approximate Bayesian procedure is of considerable interest. The most demanding part of the procedure is inference of the hidden states of the model. Previously, this has been achieved using Belief Propagation, which differs from inference in the Kalman Filtering/Smoothing literature, for which highly efficient and stabilized procedures exist. A central contribution of this paper is to show how inference *can* be written using the standard Kalman Filtering/Smoothing recursions by augmenting the original model. Additionally, a minor modification to the standard Kalman Filtering routine may be applied for computational efficiency. We demonstrated the elegance and unity of our approach by showing how to easily apply a Variational Bayes analysis of temporal ICA. Specifically, our Bayes ICA approach successfully extracts independent processes underlying EEG signals, biased towards preferred frequency ranges. We hope that this simple and unifying interpretation of Variational Bayesian LGSSMs may therefore facilitate the further application to related models.

A Parameter Updates for A and B

A.1 Determining $q(B|\Sigma_V)$

By examining \mathcal{F} , the contribution of $q(B|\Sigma_V)$ can be interpreted as the negative KL divergence between $q(B|\Sigma_V)$ and a Gaussian. Hence, optimally, $q(B|\Sigma_V)$ is a Gaussian. The covariance $[\Sigma_B]_{ij,kl} \equiv \langle (B_{ij} - \langle B_{ij} \rangle)(B_{kl} - \langle B_{kl} \rangle) \rangle$ (averages wrt $q(B|\Sigma_V)$) is given by:

$$[\Sigma_B]_{ij,kl} = [H_B^{-1}]_{jl} [\Sigma_V]_{ik}$$

where

$$[H_B]_{jl} \equiv \sum_{t=1}^T \langle h_t^j h_t^l \rangle_{q(h_t)} + \beta_j \delta_{jl}.$$

The mean is given by $\langle B \rangle = N_B H_B^{-T}$, where $[N_B]_{ij} \equiv \sum_t \langle h_t^j \rangle v_t^i + \beta_j \hat{B}_{ij}$.

Determining $q(A|\Sigma_H)$

Optimally, $q(A|\Sigma_H)$ is a Gaussian with covariance

$$[\Sigma_A]_{ij,kl} = [H_A^{-T}]_{jl} [\Sigma_H]_{ik}$$

where

$$[H_A]_{jl} \equiv \sum_{t=1}^{T-1} \langle h_t^j h_t^l \rangle_{q(h_t)} + \alpha_j \delta_{jl}.$$

The mean is given by $\langle A \rangle = N_A H_A^{-1}$, where $[N_A]_{ij} \equiv \sum_{t=2}^T \langle h_{t-1}^j h_t^i \rangle + \alpha_j \hat{A}_{ij}$.

B Covariance Updates

By specifying an Inverse Wishart prior for the covariances, conjugate update formulae are possible. In practice, it is more common to specify diagonal covariances, for which the corresponding priors are simply Inverse Gamma distributions [7, 5]. For this simple diagonal case, the explicit updates are given below.

Determining $q(\Sigma_V)$

For the constraint, $(\Sigma_V)^{-1} = \text{diag}(\rho)$ where each diagonal element follows a Gamma prior $Ga(b_1, b_2)$ [7], $q(\rho)$ factorizes and the optimal updates are

$$q(\rho_i) = Ga\left(b_1 + \frac{T}{2}, b_2 + \frac{1}{2} \left(\sum_t (v_t^i)^2 - [G_B]_{i,i} + \sum_j \beta_j \hat{B}_{ij}^2 \right)\right),$$

where $G_B \equiv N_B H_B^{-1} N_B^T$.

Determining $q(\Sigma_H)$

Analogously, for $(\Sigma_H)^{-1} = \text{diag}(\tau)$ with prior $Ga(a_1, a_2)$ [5], the updates are

$$q(\tau_i) = Ga\left(a_1 + \frac{T-1}{2}, a_2 + \frac{1}{2} \left(\sum_{t=2}^T \langle (h_t^i)^2 \rangle - [G_A]_{i,i} + \sum_j \alpha_j \hat{A}_{ij}^2 \right)\right),$$

where $G_A \equiv N_A H_A^{-1} N_A^T$.

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