Supplementary Material for: An Efficient Sequential Approach for k-parametric Dynamic Generalised Linear Models

Abstract

A novel sequential inferential method for Bayesian dynamic generalised linear models is presented, addressing both univariate and multivariate k-parametric exponential families. It efficiently handles diverse responses, including multinomial, gamma, normal, and Poisson distributed outcomes, by leveraging the conjugate and predictive structure of the exponential family. The approach integrates information geometry concepts, such as the projection theorem and Kullback-Leibler divergence, and aligns with recent advances in variational inference. Applications to both synthetic and real datasets highlight its computational efficiency and scalability, surpassing alternative methods. The approach supports the strategic integration of new information, facilitating monitoring, intervention, and the application of discount factors, which are typical in sequential analyses. The R package kDGLM is available for direct use by applied researchers, facilitating the implementation of the method for specific k-parametric dynamic generalised models.

Keywords: Bayesian analysis, Dynamic models, Information geometry, Monitoring and intervention, Sequential analysis.

S.1 Results for the simulated example of the dynamic multinomial model

In this section, we evaluate the effect of varying sample sizes T (T=2,5,20,50), as well as different specifications of the number of allocation categories, N, on the approximated inference produced by the sequential kDGLM proposal. We consider synthetic data simulated as described in Section 4.1 and compare the results of our proposed method to those obtained via non-sequential NUTS in Stan.

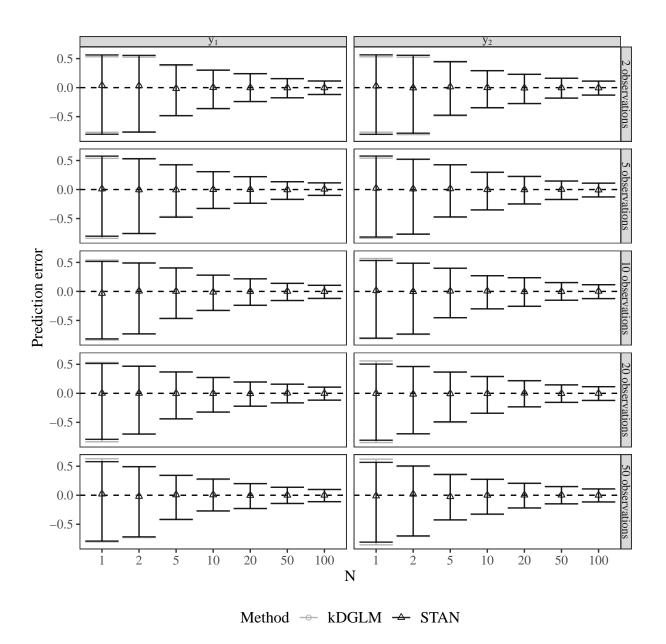


Figure S.1: The bias for the one-steap-ahead prediction for $Y_{1,T+1}$ and $Y_{2,T+1}$. Gray circles and black triangles represent the average bias among all samples, while the error bars represent the 0.975 and 0.025 quantiles. To facilitate visualization, the bias is divided by N. Gray: Sequential kDGLM. Black: Stan.

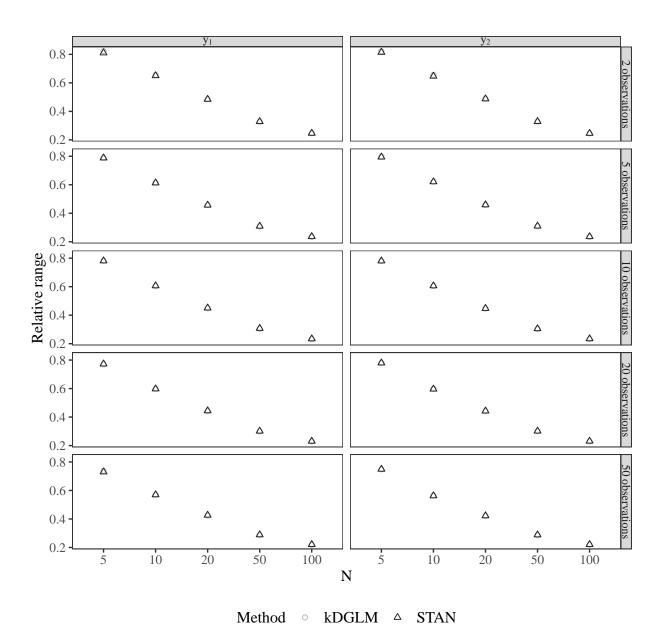


Figure S.2: The range of the 95% credibility interval for one-steap-ahead prediction for $Y_{1,T+1}$ and $Y_{2,T+1}$. Gray circles: Sequential kDGLM. Black triangles: Stan.

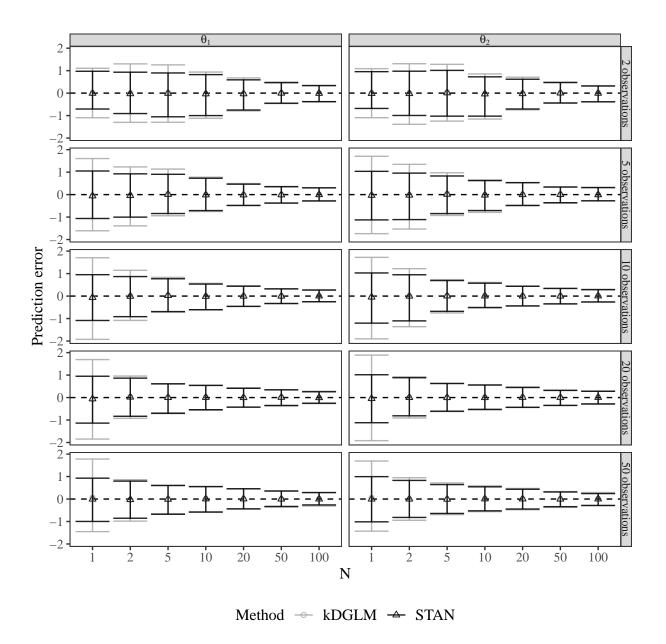


Figure S.3: The bias for the latent states at time T. Gray circles and black triangles represent the average bias among all samples, while the error bars represent the 0.975 and 0.025 quantiles. Gray: Sequential kDGLM. Black: Stan.

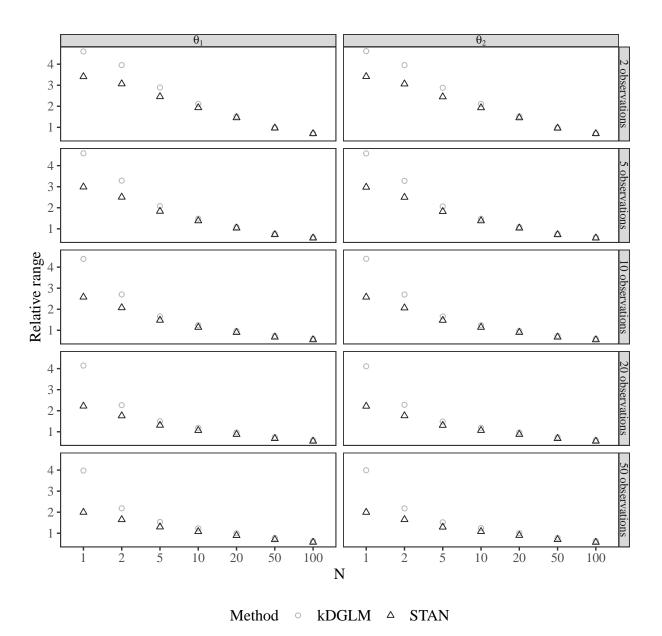


Figure S.4: The range of the 95% credibility interval for the latent states at at time T. Gray: Sequential kDGLM. Black: Stan.

S.2 Analytical development of the proposed method for normal and gamma members of the exponential family

This section provides a detailed description of the analytical developments involved in updating information for the normal and gamma distributions within the kDGLM sequential method. These are the sole instances presented in this work in which the application of the suggested methodology is not straightforward. The outlined procedures are executed at each time point t during the information updating process.

Normal distribution with unknown mean and precision: Consider the general formulation and parametrization defined in Section 4.2 for the normal model with dynamic predictive structure for the mean and precision. Since $\lambda \sim N_2(\mathbf{f}, \mathbf{Q})$, a system of equations in τ remains to be solved, that is, $E_q[\mathbf{H}_q(\mu,\phi)] = E_p[\mathbf{H}_q(\mu,\phi)]$, with q denoting a normal-gamma prior distribution and p, the density implied by the joint normality of the linear predictors for the mean ad log-precision. Using the facts that: i) $E_q[\mathbf{H}_q(\mu,\phi)] = \nabla b(\tau_1,\tau_2,\tau_3,\tau_0)$; ii) $E_p[\mathbf{H}_q(\mu,\phi)] = E_{p(\phi)}[E_{p(\mu|\phi)}\mathbf{H}_q(\mu,\phi)]$ which can be obtained by normal conditioning properties; iii) under p, $\phi \sim LN(f_2,q_2)$ and $\ln \phi \sim N(f_2,q_2)$, results in the system:

$$\begin{cases}
\frac{(2\tau_0+1)\tau_2^2}{2\tau_1\tau_2^2-8\tau_1^2\tau_3} - \frac{1}{2\tau_1} &= \exp(f_2 + q_2/2)[(f_1 + q_{12})^2 + q_1] \\
\frac{-(2\tau_0+1)\tau_2}{\tau_2^2-4\tau_1\tau_3} &= \exp(f_2 + q_2/2)(f_1 + q_{12}) \\
\frac{4\tau_1(\tau_0+1/2)}{\tau_2^2-4\tau_1\tau_3} &= \exp(f_2 + q_2/2) \\
\gamma(\tau_0+1/2) - \ln\left(\frac{\tau_2^2}{4\tau_1} - \tau_3\right) &= f_2
\end{cases}$$
(S.1)

with $\gamma(\cdot)$ denoting the digamma function. This system can be analytically solved, as follows: applying the parametrization:

$$c_0 = -2\tau_1, \quad \mu_0 = -\frac{\tau_2}{2\tau_1}, \quad \frac{d_0}{2} = \frac{\tau_2^2}{4\tau_1} - \tau_3, \quad \frac{n_0}{2} = \tau_0 + 1/2,$$
 (S.2)

to the third equation in (S.1), it follows that $n_0/d_0 = \exp(f_2 + q_2/2)$. Multiplying the third equation by $-\frac{2\tau_2}{4\tau_1} = \mu_0$ we obtain the second equation in (S.1) resulting in $\mu_0 = f_1 + q_{12}$. Doing some algebra in the first equation of (S.1), it follows that $c_0 = [q_1 \exp(f_2 + q_2/2)]^{-1}$. Finally, applying the approximation $\gamma(u) \approx \ln(u) - 1/(2u)$ to the fourth equation of the

system, results in $n_0 = 2/q_2$. Applying the reparametrization in (S.2) we obtain the following closed analytical expressions:

$$\begin{cases}
\tau_0 = 1/q_2 - 1/2 \\
\tau_1 = -[2q_1 \exp(f_2 + q_2/2)]^{-1} \\
\tau_2 = (f_1 + q_{12})[q_1 \exp(f_2 + q_2/2)]^{-1} \\
\tau_3 = -(f_1 + q_{12})^2[2q_1 \exp(f_2 + q_2/2)]^{-1} - [q_2 \exp(f_2 + q_2/2)]^{-1}.
\end{cases}$$

After observing \boldsymbol{y} , we obtain $\boldsymbol{\tau}^* = (\tau_1^*, \tau_2^*, \tau_3^*, \tau_0^*) = (\tau_1 - 1/2, \tau_2 + y, \tau_3 - y^2/2, \tau_0 + 1/2)$, the updated canonical parameters of a normal-gamma posterior density for (μ, ϕ) , and need to evaluate the parameters $(\boldsymbol{f}^*, \boldsymbol{Q}^*)$ of the posterior distribution of the linear predictors that are compatible with them, solving: $E_q[\boldsymbol{H}_q] = E_p[\boldsymbol{H}_q]$, now considering that p is a normal-gamma density for (μ, ϕ) and q is a bivariate normal for $(\mu, \ln \phi)$. $\boldsymbol{H}'_q = (\boldsymbol{\lambda}, \boldsymbol{\lambda} \boldsymbol{\lambda}')$, so the following system is trivially solved:

$$\begin{cases} f_1^* &= -\frac{\tau_2^*}{2\tau_1^*} \\ f_2^* &= \gamma(\tau_0^* + 1/2) - \ln\left(\frac{(\tau_2^*)^2}{4\tau_1^*} - \tau_3^*\right) \\ Q_{11}^* &= \frac{\tau_2^{*2} - 4\tau_1\tau_3}{4\tau_0\tau_1 - 2\tau_1} \\ Q_{12}^* &= 0 \\ Q_{22}^* &= \gamma'(\tau_0^* + 1/2). \end{cases}$$

Notice that, no matter the values of τ_0^* , τ_1^* , τ_2^* and τ_3^* , $Q_{12}^* = 0$, since the parameters that minimize the KL divergence from p to q must satisfy $Cov_p[\mu, \ln(\phi)] = Cov_q[\mu, \ln(\phi)]$ and, as a property of the normal-gamma distribution, we have that (μ, ϕ) are uncorrelated (although not independent), consequently, it can be shown that $(\mu, \ln \phi)$ are also uncorrelated.

Using the reparametrization $c_0^* = -2\tau_1^*$, $\mu_0^* = -\frac{\tau_2^*}{2\tau_1^*}$, $\frac{d_0^*}{2} = \frac{\tau_2^{*2}}{4\tau_1^*} - \tau_3^*$ and $\frac{n_0^*}{2} = \tau_0^* + 1/2$, we obtain:

$$\begin{cases} f_1^* &= \mu_0^* \\ f_2^* &= \gamma(n_0^*/2) - \ln(d_0^*/2) \\ Q_{11}^* &= \frac{d_0^*/2}{c_0^*(n_0^*/2 - 1)} \\ Q_{12}^* &= 0 \\ Q_{22}^* &= \gamma'(n_0^*/2). \end{cases}$$

An important point to discuss is the equation that defines Q_{11}^* . Specifically, notice that Q_{11}^* is not defined for $n_0^* \leq 2$. Indeed, by the Projection Theorem, the parameters

that minimize the KL divergence from p to q must satisfy $Var_p[\mu] = Var_q[\mu]$, but in the conjugated posterior distribution $\mu|\phi \sim N(\mu_0^*, (c_0^*\phi)^{-1})$ and $\phi \sim \mathcal{G}(n_0^*/2, d_0^*/2)$, which implies that $\mu \sim t\left(n_0^*, \mu_0^*, \frac{d_0^*/2}{c_0^*n_0^*/2}\right)$, so that when $n_0^* \leq 2$, $Var_p[\mu] = +\infty$, and the system $E_q[\mathbf{H}_q] = E_p[\mathbf{H}_q]$ has no valid solutions. Indeed, one can always reduce the KL divergence from p to q by increasing Q_{11}^* . This situation is quite undesirable since it compromises the use of our proposed methodology. Yet, we can restrict the parameter space for the normal posterior distribution to guarantee that, inside that restricted space, we always have a minimum for the divergence from p to q. A natural restriction for the parameter space is to set $Q_{11}^* = \frac{d_0^*/2}{c_0^*n_0^*/2}$, since:

- This restriction guarantees that the scale parameter for the marginal distribution of μ is identical in both the normal and the conjugated distribution.
- for a large value of n_0^* (which we expect to have for a reasonable sample size), $\frac{d_0^*/2}{c_0^*n_0^*/2} \approx \frac{d_0^*/2}{c_0^*(n_0^*/2-1)}$, i.e., for large values of n_0^* this restriction has no significant effect, in the sense that the optimum in the restricted space will be very close to the global optimum.
- $Q_{11}^* = \frac{d_0^*/2}{c_0^* n_0^*/2}$ becomes numerically well behaved for any possible value of n_0^* , since, after updating our knowledge of μ , it is guaranteed that $n_0^* \geq 1$, avoiding a division by values close to zero.

Gamma distribution with unknown shape α : Let $\eta = (\alpha, \mu)$ and $(y|\eta) \sim \mathcal{G}(\alpha, \alpha/\mu)$, $\alpha, \mu > 0$. Unlike other particular cases discussed in this work, the normalising constant of the joint prior for the pair (α, μ) is unknown and the conjugate prior is $\pi(\alpha, \mu) \propto \exp\left\{\tau_0\left(\alpha \ln\left(\frac{\alpha}{\mu}\right) - \ln(\Gamma(\alpha)\right) + \tau_1\alpha - \tau_2\frac{\alpha}{\mu}\right\}$. Consider a vector of linear predictors $\lambda' = (\lambda_1, \lambda_2) = (\ln(\mu), \ln(\phi)) = (F'_1\theta_1, F'_2\theta_2)$. Following Algorithm 1:

- Step 1: Given posterior moments for the states, obtain the prior moments of $\lambda \sim N(f, Q)$.
- Step 2.1: Obtain the conjugate prior sufficient statistics vector: $\boldsymbol{H}_q = \left(\alpha \ln(\mu) \alpha \ln(\alpha) + \ln(\Gamma(\alpha)), \alpha, \frac{\alpha}{\mu}\right)'.$
- Step 2.2: Solve the system $E_q[\mathbf{H}_q(\boldsymbol{\eta})] = E_p[\mathbf{H}_q(\boldsymbol{\eta})]$, where p is the prior density induced by the normal specification for the linear predictors. It is easily seen that $\mu|\alpha \sim \mathcal{IG}(\tau_0\alpha 1, \alpha\tau_2)$

where \mathcal{IG} denotes the inverse gamma density, if $\alpha > 1/\tau_0$. Using the fact that $E_q[\mathbf{H}_q] = E_q[E_q[\mathbf{H}_q|\alpha]]$ and that $E_q[\ln(\mu)|\alpha] = -\gamma(\tau_0\alpha - 1) + \ln(\alpha\tau_2)$, it follows that:

$$E_{q}\left[\frac{\alpha}{\mu}\right] = E_{q}\left[\alpha E_{q}\left[\frac{1}{\mu}|\alpha\right]\right] = E_{q}\left[\alpha \frac{\tau_{0}\alpha + 1}{\alpha\tau_{2}}\right] = \frac{\tau_{0}E_{q}\left[\alpha\right] - 1}{\tau_{2}};$$

$$E_{q}[\alpha\ln(\mu)] = E_{q}[\alpha E_{q}[\ln(\mu)|\alpha]] = E_{q}\left[\alpha(-\gamma(\tau_{0}\alpha + 1) + \ln(\alpha\tau_{2}))\right].$$

Assuming $E_p[H_q]$ known, and using the fact that $E_q[\alpha] = E_p[\alpha]$ (from the system to be solved), it follows that $\tau_2 = \frac{\tau_0 E_q[\alpha] - 1}{E_q \left[\frac{\alpha}{\mu}\right]}$. One can trivially find that $E_q[\alpha \ln(\mu) - \alpha \ln(\alpha) + \ln(\Gamma(\alpha)] = E_q[\ln(\Gamma(\alpha) - \alpha\gamma(\tau_0\alpha + 1)] + E_q[\alpha]\ln(\tau_2)$. Thus, $E_q[\mathbf{H}_q]$ can be expressed in terms of expected values which solely depend on the marginal prior for α : $\pi(\alpha) = \int_0^{+\infty} \pi(\alpha, \mu) d\mu \propto \frac{\alpha^{\tau_0\alpha}}{\alpha^{-\tau_0}\Gamma(\alpha+1)^{\tau_0}} \frac{\Gamma(\tau_0\alpha+1)}{\alpha^{\tau_0\alpha+1}\tau_2^{\tau_0\alpha+1}} \exp\{\tau_1\alpha\}$. Applying Stirling's formula $\Gamma(x+1) \approx \sqrt{2\pi x} \left(\frac{x}{e}\right)^x$ to this last expression results in: $\pi(\alpha) \propto \alpha^{\frac{\tau_0-1}{2}-1} \exp\{-\left[\tau_0 \ln\left(\frac{\tau_2}{\tau_0}\right) - \tau_1\right]\alpha\}$, that is, α approximately follows a $\mathcal{G}\left(\frac{\tau_0+1}{2}, \tau_0 \ln\left(\frac{\tau_2}{\tau_0}\right) - \tau_1\right)$ marginal conjugate prior, so that:

$$E_q[\alpha] = \frac{\tau_0 + 1}{2\left(\tau_0 \ln\left(\frac{\tau_2}{\tau_0}\right) - \tau_1\right)}.$$

Using a second order approximation for the digamma function, $\gamma(x) \approx \ln(x) - \frac{1}{2x} - \frac{1}{12x^2}$, and the fact that $\gamma(x+1) = \gamma(x) + \frac{1}{x}$ results in:

$$-\alpha\gamma(\tau_0\alpha + 1) = -\alpha\left(\gamma(\tau_0\alpha) + \frac{1}{\tau_0\alpha}\right)$$

$$\approx -\alpha\left(\ln(\tau_0\alpha) - \frac{1}{2\tau_0\alpha} - \frac{1}{12\tau_0^2\alpha^2} + \frac{1}{\tau_0\alpha}\right)$$

$$\approx -\alpha\ln(\alpha) - \alpha\ln(\tau_0) + \frac{1}{12\tau_0^2\alpha} - \frac{1}{2\tau_0}.$$

Once again using a second order approximation for the digamma function and the fundamental theorem of Calculus:

$$\ln(\Gamma(x)) = \int_1^x \gamma(t)dt \approx \int_1^x \ln(t) - \frac{1}{2t} - \frac{1}{12t^2}dt = x\ln(x) - x - \frac{1}{2}\ln(x) + \frac{1}{12x} + \frac{11}{12}.$$

The stated facts allow us to calculate the expected value of the remaining sufficient statistic, resulting in the following system to be solved:

$$\begin{cases}
\mathbb{E}_{p} \left[(\alpha \ln(\mu) - \alpha \ln(\alpha) + \ln(\Gamma(\alpha)) \right] &= \frac{\tau_{0} + 1}{2(\tau_{0} \ln(\frac{\tau_{2}}{\tau_{0}}) - \tau_{1})} (\ln(\tau_{2} / \tau_{0}) - 1) - \frac{1}{2} \psi\left(\frac{\tau_{0} + 1}{2}\right) \\
+ \frac{1}{2} \ln\left(\tau_{0} \ln\left(\frac{\tau_{2}}{\tau_{0}}\right) - \tau_{1}\right) + \frac{\tau_{0} + 1}{\tau_{0} - 1} \frac{\ln\left(\frac{\tau_{2}}{\tau_{0}}\right) - \frac{\tau_{1}}{\tau_{0}}}{6} \\
+ \frac{11}{12} - \frac{1}{2\tau_{0}} \\
\exp\{f_{1} + Q_{11} / 2\} &= \frac{\tau_{0} + 1}{2(\tau_{0} \ln(\frac{\tau_{2}}{\tau_{0}}) - \tau_{1})} \\
\exp\{f_{1} - f_{2} + Q_{11} / 2 + Q_{12} + Q_{22} / 2\} &= \frac{\tau_{0}(\tau_{0} + 1) - 1}{\tau_{2} 2(\tau_{0} \ln(\frac{\tau_{2}}{\tau_{0}}) - \tau_{1})}
\end{cases}$$

with $\mathbb{E}_p\left[\left(\alpha \ln(\mu) - \alpha \ln(\alpha) + \ln(\Gamma(\alpha))\right)\right]$ evaluated by numerical integration, with negligible cost.

- Step 3: Update the hyperparameters of the conjugate specification: $\boldsymbol{\tau}^* = (\tau_0^*, \tau_1^*, \tau_2^*) = (\tau_0 + 1, \tau_1 + \ln(y), \tau_2 + y).$
- Step 4.1: Let q denote a multivariate normal density for the vector of linear predictors, with sufficient statistics vector: $\mathbf{H}'_q = (\lambda, \ \lambda \lambda')$.
- Step 4.2: Solve the system $E_q[\mathbf{H}_q(\boldsymbol{\lambda})] = E_p[\mathbf{H}_q(\boldsymbol{\lambda})]$, where p is the updated distribution, obtaining:: $f_1^* = E_p[\ln(\mu)], \ f_2^* = E_p[\ln(\alpha)]; \ Q_{11}^* = V_p[\ln(\mu)], \ Q_{12}^* = Cov_p[\ln(\mu), \ln(\alpha)], \ Q_{22}^* = E_p[\ln(\alpha)]$. Using the fact that α approximately follows a $\mathcal{G}\left(\frac{\tau_0^*+1}{2}, \tau_0^* \ln\left(\frac{\tau_2^*}{\tau_0^*}\right) \tau_1^*\right)$ conjugate marginal posterior, it follows that:

$$E_p[\ln(\alpha)] \approx \gamma \left(\frac{\tau_0^* + 1}{2}\right) - \ln\left(\tau_0^* \ln\left(\frac{\tau_2^*}{\tau_0^*}\right) - \tau_1^*\right)$$

$$V_p[\ln(\alpha)] \approx \gamma' \left(\frac{\tau_0^* + 1}{2}\right).$$

Since, a posteriori, $\mu | \alpha \sim \mathcal{IG}(\tau_0^* \alpha + 1, \alpha \tau_2^*)$, it follows that:

$$E_{p}[\ln(\mu)] = E_{p}[E_{p}[\ln(\mu)|\alpha]] = -E_{p}[\gamma(\tau_{0}^{*}\alpha + 1) - \ln(\alpha\tau_{2}^{*})];$$

$$V_{p}[\ln(\mu)] = V_{p}[\ln(1/\mu)] = E_{p}[V_{p}[\ln(1/\mu)|\alpha]] + V_{p}[E_{p}[\ln(1/\mu)|\alpha]]$$

$$= E_{p}[\gamma(\tau_{0}^{*}\alpha + 1)] + V_{p}[\gamma(\tau_{0}^{*}\alpha + 1) - \ln(\alpha\tau_{2}^{*})];$$

$$E_{p}[\ln(\alpha)\ln(\mu)] = -E_{p}[\ln(\alpha)\ln(1/\mu)] = -E_{p}[\ln(\alpha)E_{p}[\ln(1/\mu)|\alpha]]$$

$$= -E_{p}[\ln(\alpha)(\gamma(\tau_{0}^{*}\alpha + 1) - \ln(\alpha\tau_{2}^{*}))].$$

All the integrals involved in the expected values above are evaluated by numerical integration, with negligible cost.

• Step 5: The updated moments of the states θ are trivially obtained, applying normal distribution properties.

Since the normalizing constant of the conjugate prior is unknown, we trivially obtain a sample from the predictive distribution, first sampling from the posterior λ and then sampling from the observational distribution of $y^*|\lambda$. Since we assume that λ follows a multivariate normal distribution, the computational cost of this process is negligible.

S.3 Comparison between kDGLM and NUTS via Stan for the stochastic volatility example

Section 4.2.1 presents a comparison between the results obtained using the sequential kDGLM method proposed in this work and those from a NUTS scheme implemented via Stan. The comparison relies on synthetic, artificially generated data. As discussed, the differences observed in the inference produced by the two methods are largely attributable to the linearization employed for sequential inference for the autoregressive parameter γ in the kDGLM method.

In Figure S.5, we display the inference results under both methods, kDGLM and NUTS, with the parameter γ fixed at its true value. As shown, the two methods yield nearly identical results. Therefore, for models where the evolution matrix \mathbf{G} is known, the kDGLM method is expected to produce inferences closely aligned with those generated by Markov Chain Monte Carlo schemes.

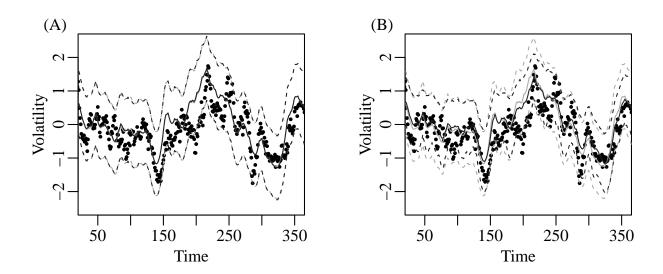


Figure S.5: Comparison of smoothed means (solid lines) and credible intervals (dashed lines) for volatility estimation from synthetic data, with fixed autoregressive parameter γ . The solid circles indicate theoretical volatilities. (A): results of the proposed sequential approach for normal (grey) and gamma (black) formulations. (B): comparison of normal formulation using the sequential approach (grey) and NUTS via Stan (black).