

Particle filters in the dynamichazard package

BENJAMIN CHRISTOFFERSEN

JANUARY 8, 2019

This vignette covers the particle filters and smoothers implemented in the `dynamichazard` package in R. Some prior knowledge of particle filters is assumed. Doucet and Johansen (2009) provide a tutorial on particle filters and Kantas et al. (2015) cover parameter estimation with particle filters. See also Cappé et al. (2005) for a general introduction to Hidden Markov models. This vignette relies heavily on Fearnhead et al. (2010) and there is a big overlap between what is presented here and the paper.

1 Method

The models implemented in the package is survival analysis models for terminal events. These can be in discrete times where we have binary indicators $Y_{ik} = 1_{\{T_i \in (t_{k-1}, t_k]\}}$ which is one if the random event time of individual i denoted by $T_i \in (0, \infty)$ is in the interval $(t_{k-1}, t_k]$ and zero otherwise. It can also be in continuous time where we model the distribution of the event time of individual i , T_i , with a piecewise constant exponential distribution conditional on observable covariates and the path a discrete latent variable. To be more concrete, the model is

$$\begin{aligned}
 y_{it} &\sim g(y_{it}|\eta_{it}) \\
 \eta_t &= \mathbf{X}_t \mathbf{R}^+ \boldsymbol{\alpha}_t + \mathbf{o}_t + \mathbf{Z}_t \boldsymbol{\omega} & i = 1, \dots, n_t \\
 \boldsymbol{\alpha}_t &= \mathbf{F} \boldsymbol{\alpha}_{t-1} + \mathbf{R} \boldsymbol{\epsilon}_t & \boldsymbol{\epsilon}_t \sim \mathcal{N}(\mathbf{0}, \mathbf{Q}), \quad t = 1, \dots, d \\
 & & \boldsymbol{\alpha}_0 \sim \mathcal{N}(\boldsymbol{\mu}_0, \mathbf{Q}_0)
 \end{aligned} \tag{1}$$

where I denote the conditional densities as $g_t(\mathbf{y}_t|\boldsymbol{\alpha}_t) = g(\mathbf{y}_t|\mathbf{X}_t \mathbf{R}^+ \boldsymbol{\alpha}_t + \mathbf{o}_t + \mathbf{Z}_t \boldsymbol{\omega})$ and $f(\boldsymbol{\alpha}_t|\boldsymbol{\alpha}_{t-1})$. For each $t = 1, \dots, d$, we have risk set given by $R_t \subseteq \{1, 2, \dots, n\}$. Further, we let $n_t = |R_t|$ denote the number of observation at risk at time t and $n_{\max} = \max_{t \in \{1, \dots, d\}} n_t$. The observed outcomes are denoted by $\mathbf{y}_t = \{y_{it}\}_{i \in R_t}$. \mathbf{X}_t is the design matrix of the covariates and $\boldsymbol{\alpha}_t$ is the state vector containing the time-varying coefficients. The \mathbf{Z}_t is the design matrix for the covariates with time-invariant coefficients and $\boldsymbol{\omega}$ are the corresponding coefficients.

The i 'th row of \mathbf{X}_t is \mathbf{x}_{it} , $\mathbf{x}_{it}, \boldsymbol{\epsilon}_t \in \mathbb{R}^r$, $\boldsymbol{\mu}_0, \boldsymbol{\alpha}_t \in \mathbb{R}^p$, $\mathbf{F} \in \mathbb{R}^{p \times p}$, $\mathbf{Q} \in \mathbb{R}^{r \times r}$ is a positive definite matrix, $\mathbf{Q}_0 \in \mathbb{R}^{p \times p}$ is a positive definite matrix, \mathbf{o}_t s are known offsets, and \mathbf{R} is a $p \times r$ matrix with $p \geq r$ which contains a subset of r columns of \mathbf{I}_p . We will order the entires of \mathbf{R} such that the first r columns are the first r columns of \mathbf{I}_p . I.e.,

$$\mathbf{R} = \begin{pmatrix} \mathbf{I}_r \\ \mathbf{0} \end{pmatrix}$$

Superscript $+$ denotes the Moore-Penrose inverse, $\mathbf{R}^+ = \mathbf{R}^\top$ and \mathbf{R} is left inverse (i.e., $\mathbf{R}^\top \mathbf{R} = \mathbf{I}_r$). $\mathbf{R}\mathbf{R}^\top$ is a $p \times p$ diagonal matrix where the r first diagonal entries has value

1 and the rest of the diagonal entries are zero. The data sets we are working with have $n_{\max} \gg p \geq r$ (e.g. $n_{\max} = 100000$ and $r = 5$). We let

$$\boldsymbol{\xi}_t = \mathbf{R}^+ \boldsymbol{\alpha}_t$$

The above allows us to have an o th order vector autoregressions, VAR(o), by settings

$$\boldsymbol{\alpha}_t = (\boldsymbol{\xi}_t, \boldsymbol{\xi}_{t-1}, \dots, \boldsymbol{\xi}_{t-o+1})$$

$$\mathbf{F} = \begin{pmatrix} \mathbf{F}_1 & \cdots & \cdots & \mathbf{F}_{o-1} & \mathbf{F}_o \\ \mathbf{I}_r & \mathbf{0} & \cdots & \mathbf{0} & \mathbf{0} \\ \mathbf{0} & \mathbf{I}_r & \ddots & \vdots & \vdots \\ \vdots & \ddots & \ddots & \mathbf{0} & \mathbf{0} \\ \mathbf{0} & \cdots & \mathbf{0} & \mathbf{I}_r & \mathbf{0} \end{pmatrix}, \quad \mathbf{F}_i \in \mathbb{R}^{r \times r}$$

I will use a particle filter and smoother to get smoothed estimates of $\boldsymbol{\alpha}_1, \dots, \boldsymbol{\alpha}_d$ given the outcomes $\mathbf{y}_{1:d} = \{\mathbf{y}_1, \mathbf{y}_2, \dots, \mathbf{y}_d\}$ and use an EM-algorithm to estimate \mathbf{Q} , $\boldsymbol{\omega}$, and $\boldsymbol{\mu}_0$. One choice of smoother is shown in Fearnhead et al. (2010) and another is the the generalized two-filter smoother shown by Briers et al. (2009). The rest of vignette is structured as follows: first I give a brief introduction to the implemented particle filters and smoothers. Then I cover what the effect is of some the arguments to particle functions in R in the packages. The implemented particle filter and smoother from Fearnhead et al. (2010) is presented next, followed by the used EM-algorithm, the smoother in Briers et al. (2009), and other miscellaneous topics.

1.1 Overview

As a quick recap which will ease reading the next sections, we will start by recalling an application of importance sampling, use this to motivate particle filtering, and give a brief idea of the implemented particle smoothers. Suppose we want to approximate a density $c(x) = \zeta \tilde{c}(x)$ where we only know $\tilde{c}(x)$ and not the normalization constant ζ . One way to approximate this density is to

- sample x_1, x_2, \dots, x_N from a distribution with density $b(x)$.
- Compute the unnormalized weights $\bar{w}_i = \tilde{c}(x_i)/b(x_i)$.
- Normalize the weights $w_i = \bar{w}_i / \sum_{i=1}^N \bar{w}_i$.

This gives us the following discrete approximation of the density

$$c(x) \approx \sum_{i=1}^N w_i \delta_{x_i}(x)$$

where δ_x is the Dirac delta function which has unit point mass at x . This is directly applicable to the model in Equation (1) as at time 1 we want to approximate

$$p(\boldsymbol{\alpha}_1 | \mathbf{y}_1) = \frac{g_1(\mathbf{y}_1 | \boldsymbol{\alpha}_1) \int f(\boldsymbol{\alpha}_1 | \boldsymbol{\alpha}_0) \phi(\boldsymbol{\alpha}_0 | \boldsymbol{\mu}_0, \mathbf{Q}_0) d\boldsymbol{\alpha}_0}{p(\mathbf{y}_1)}$$

where $\phi(\cdot|\mathbf{m}, \mathbf{M})$ is the density function of a multivariate normal distribution with mean \mathbf{m} and covariance matrix \mathbf{M} . We can easily evaluate the numerator for each α_1 but not the normalization constant, $p(\mathbf{y}_1)$.

The extension to a particle filter (which I will call a forward particle filter) is that at time 2 we want to approximate

$$p(\alpha_{1:2}|\mathbf{y}_{1:2}) = p(\alpha_1|\mathbf{y}_1) \frac{g_2(\mathbf{y}_2|\alpha_2) f(\alpha_2|\alpha_1)}{p(\mathbf{y}_2|\mathbf{y}_1)}$$

Now, we can use the discrete approximation at time 1 of $p(\alpha_1|\mathbf{y}_1)$, sample α_2 given each sampled α_1 , and apply importance sampling again. We can repeat this with similar arguments at time 3, 4, \dots , d giving us an approximation of $p(\alpha_{1:d}|\mathbf{y}_{1:d})$. We will call the last element of a sampled path at time t a particle. Further, we will denote the j th particle at time t and its' associated weight by $\alpha_t^{(j)}$ and $w_t^{(j)}$ respectively. One issue that may arise is that our samples (particles) may degenerate so essentially only one sampled path of $\alpha_{1:d}$ has any weight in the end. To avoid this, we may introduce a re-sampling step. One way to re-sample is using the weights and letting the re-sampling weights be $\beta_{t+1}^{(j)} = w_t^{(j)}$ where $\beta_{t+1}^{(j)}$ is the re-sampling weight of particle j at time t . We then sample with replacement using $\beta_{t+1}^{(j)}$. Another option when we re-sample the particles from time t is to use the information of the outcomes at time $t+1$, \mathbf{y}_{t+1} . This is called an auxiliary particle filter and is introduced in Pitt and Shephard (1999).

However, we may end up with few or only one unique value at the early time points (say α_1) when we re-sample. Thus, it will be useful to use a smoother to get a better approximation of the marginal density $p(\alpha_t|\mathbf{y}_{1:d})$. To do so, one idea is to use the two-filter formula from Kitagawa (1994). Though, this requires that we can evaluate $p(\mathbf{y}_{t:d}|\alpha_t)$. It turns out that we can approximate this up to a constant which just what need (see Section 7).

The approximation uses a particle filter which is run backwards in time and which approximates an artificial distribution. The arguments for the backward particle filter is very similar to the forward particle filter presented above. The k th particle in the backward particle filter at time t , its' re-sampling weight, and the associated weight will be denoted by respectively $\tilde{\alpha}_t^{(k)}$, $\tilde{\beta}_{t-1}^{(k)}$ and $\tilde{w}_t^{(k)}$. The final i th smoothed particle and weight at time t will be denoted by $\hat{\alpha}_t^{(i)}$ and $\hat{w}_t^{(i)}$. The latter gives us the following approximation of the marginal density of $\alpha_t | \mathbf{y}_{1:d}$

$$p(\alpha_t|\mathbf{y}_{1:d}) \approx \sum_{i=1}^{N_S} \hat{w}_t^{(i)} \delta_{\hat{\alpha}_t^{(i)}}(\alpha_t)$$

if we sampled N_S smoothed particles at time t . The smoothing algorithm from Fearnhead et al. (2010) is shown in Algorithm 1, the forward particle filter is shown Algorithm 2, and the backward particle filter is shown in Algorithm 3.

1.2 Methods in the package

The PF_EM method in the `dynamichazard` package contains an implementation of the described methods. You specify the number of particles by the `N_first`, `N_fw_n_bw` and `N_smooth` argument for respectively the N_f , N and N_s in the Algorithm 1-3. We may want more particles in the smoothing step, $N_s > N$, as pointed out in the discussion in Fearnhead

et al. (2010, page 460 and 461). Further, selecting $N_f > N$ may be preferable to ensure coverage of the state space at time 0 and $d + 1$.

We do not need to sample the time 0 and $d + 1$ particles. Instead we can make a special proposal distribution for time 1 and time d . This is not implemented though...

The `method` argument specify how the filters are set up. The argument can take the following values

- `"bootstrap_filter"` for a bootstrap filter. This is where we sample using Equation (5), (11) and (14). This is fast but the proposal distribution may be a poor approximation of the distribution we want to target.
- `"PF_normal_approx_w_cloud_mean"` and `"AUX_normal_approx_w_cloud_mean"` for the Taylor approximation of the conditional density of \mathbf{y}_t made using the mean of the parent particles and/or mean of the child particles. See Section 2. The PF and AUX prefix specifies whether or not the auxiliary version should be used.
- `"PF_normal_approx_w_particles"` and `"AUX_normal_approx_w_particles"` for the Taylor approximation of the conditional density of \mathbf{y}_t made using the parent and/or child particle. See Section 2. The PF and AUX prefix specifies whether or not the auxiliary version should be used.

The smoother is selected with the `smoother` argument. `"Fearnhead_0_N"` gives the smoother in Algorithm 1 and `"Brier_0_N_square"` gives the smoother in Algorithm 4. The *Systematic Re-sampling* (Kitagawa, 1996) is used in all re-sampling steps. See Douc and Cappé (2005) for a comparison of re-sampling methods. The rest of the arguments to `PF_EM` are similar to those of the `ddhazard` function.

It is not clear what will give the best performance for a given data set. An advice is to use the `trace` argument and check the effective sample at each point in time doing the estimation. `"bootstrap_filter"` may not be that much cheaper in terms of computation time as we still have to evaluate g_t in Equation (16), (17), and (19) which is $\mathcal{O}(n_{\max}Nr)$ or $\mathcal{O}(n_{\max}N_Sr)$ run times. On the other hand, the `"...w_particles"` methods are $\mathcal{O}(n_{\max}Nr^2)$ or $\mathcal{O}(n_{\max}N_Sr^2)$ run times with a potentially much larger constant. Thus, the `"...w_cloud_mean"` may be preferred.

The rest of the vignette covers the implemented methods. It is mainly included to show exactly what is computed and why. Further, I cover some currently not implemented extensions that may be implemented in the future.

1.3 Proposal distributions and re-sampling weights

Algorithm 1 shows one of the particle smoothers shown by Fearnhead et al. (2010) in the first order state space model. In this situation $\mathbf{R} = \mathbf{I}_r$, $r = p$ and $\boldsymbol{\alpha}_t = \boldsymbol{\xi}_t$. We need to specify a series of proposal distributions and re-sampling weights. To show what is implemented and why, we first consider the model where

$$\mathbf{y}_t \mid \boldsymbol{\alpha}_t \sim \mathcal{N}(\mathbf{X}_t\boldsymbol{\alpha}_t + \mathbf{o}_t + \mathbf{Z}_t\boldsymbol{\omega}, \mathbf{H}_t)$$

for some known positive definite matrix \mathbf{H}_t . This is not implemented in this package but deriving optimal re-sampling weights and proposal distributions is possible in this case. In fact, it makes little sense to use a particle filter and particle smoother in this case since the

Kalman filter and an exact smoother can be applied. However, the results here will turn out to be useful to motivate the approximations we use later. The state space model is

$$\begin{aligned} \mathbf{y}_t &\sim \mathcal{N}(\boldsymbol{\eta}_t, \mathbf{H}_t) \\ \boldsymbol{\eta}_t &= \mathbf{X}_t \boldsymbol{\alpha}_t + \mathbf{o}_t + \mathbf{Z}_t \boldsymbol{\omega} & i = 1, \dots, n_t \\ \boldsymbol{\alpha}_t &= \mathbf{F} \boldsymbol{\alpha}_{t-1} + \mathbf{R} \boldsymbol{\epsilon}_t & \boldsymbol{\epsilon}_t \sim \mathcal{N}(\mathbf{0}, \mathbf{Q}), \quad t = 1, \dots, d \\ & & \boldsymbol{\alpha}_0 \sim \mathcal{N}(\boldsymbol{\mu}_0, \mathbf{Q}_0) \end{aligned}$$

We let $\mathbf{h}_t = \mathbf{o}_t + \mathbf{Z}_t \boldsymbol{\omega}$ such that $\boldsymbol{\eta}_t = \mathbf{X}_t \boldsymbol{\alpha}_t + \mathbf{h}_t$ to ease the notation. We first turn to the forward particle filter in Algorithm 2. Ideally, we want the re-sampling weights to be

$$\begin{aligned} \beta_t^{(j)} &\propto p(\mathbf{y}_t | \boldsymbol{\alpha}_{t-1}^{(j)}) w_{t-1}^{(j)} & (2) \\ &= \int g_t(\mathbf{y}_t | \boldsymbol{\alpha}_t) f(\boldsymbol{\alpha}_t | \boldsymbol{\alpha}_{t-1}^{(j)}) d\boldsymbol{\alpha}_t w_{t-1}^{(j)} \\ &= \phi(\mathbf{y}_t | \mathbf{X}_t \mathbf{F} \boldsymbol{\alpha}_{t-1}^{(j)} + \mathbf{h}_t, \mathbf{X}_t \mathbf{Q} \mathbf{X}_t^\top + \mathbf{H}_t) w_{t-1}^{(j)} \end{aligned}$$

We can notice that setting $\beta_t^{(j)} = w_{t-1}^{(j)}$ yields the so-called sequential importance re-sampling algorithm. For the proposal distribution, the optimal proposal density is

$$q(\boldsymbol{\alpha}_t | \boldsymbol{\alpha}_{t-1}^{(j)}, \mathbf{y}_t) = p(\boldsymbol{\alpha}_t | \boldsymbol{\alpha}_{t-1}^{(j)}, \mathbf{y}_t)$$

where we find that

$$\begin{aligned} \log p(\boldsymbol{\alpha}_t | \boldsymbol{\alpha}_{t-1}^{(j)}, \mathbf{y}_t) &= \log p(\boldsymbol{\alpha}_t, \mathbf{y}_t | \boldsymbol{\alpha}_{t-1}^{(j)}) + \dots \\ &= \log g_t(\mathbf{y}_t | \boldsymbol{\alpha}_t) + \log f(\boldsymbol{\alpha}_t | \boldsymbol{\alpha}_{t-1}^{(j)}) + \dots \\ &= -\frac{1}{2}(\mathbf{y}_t - \mathbf{X}_t \boldsymbol{\alpha}_t - \mathbf{h}_t)^\top \mathbf{H}_t^{-1} (\mathbf{y}_t - \mathbf{X}_t \boldsymbol{\alpha}_t - \mathbf{h}_t) \\ &\quad - \frac{1}{2}(\boldsymbol{\alpha}_t - \mathbf{F} \boldsymbol{\alpha}_{t-1}^{(j)})^\top \mathbf{Q}^{-1} (\boldsymbol{\alpha}_t - \mathbf{F} \boldsymbol{\alpha}_{t-1}^{(j)}) + \dots \\ &= -\frac{1}{2} \boldsymbol{\alpha}_t^\top \boldsymbol{\Sigma}_t^{-1} \boldsymbol{\alpha}_t + \boldsymbol{\alpha}_t^\top \boldsymbol{\Sigma}_t^{-1} \boldsymbol{\mu}(\boldsymbol{\alpha}_{t-1}^{(j)}) + \dots \\ \boldsymbol{\Sigma}_t &= (\mathbf{Q}^{-1} + \mathbf{X}_t^\top \mathbf{H}_t^{-1} \mathbf{X}_t)^{-1} & (3) \\ \boldsymbol{\mu}(\mathbf{x}) &= \boldsymbol{\Sigma}_t (\mathbf{Q}^{-1} \mathbf{F} \mathbf{x} + \mathbf{X}_t^\top \mathbf{H}_t^{-1} (\mathbf{y}_t - \mathbf{h}_t)) & (4) \end{aligned}$$

The ... are terms of the normalization constant. We recognize the multivariate normal distribution density and thus the optimal proposal density is

$$q(\boldsymbol{\alpha}_t | \boldsymbol{\alpha}_{t-1}^{(j)}, \mathbf{y}_t) = \phi(\boldsymbol{\alpha}_t | \boldsymbol{\mu}(\boldsymbol{\alpha}_{t-1}^{(j)}), \boldsymbol{\Sigma}_t)$$

Alternatively, we can use the so-called *bootstrap filter* and let

$$q(\boldsymbol{\alpha}_t | \boldsymbol{\alpha}_{t-1}^{(j)}, \mathbf{y}_t) = \phi(\boldsymbol{\alpha}_t | \mathbf{F} \boldsymbol{\alpha}_{t-1}^{(j)}, \mathbf{Q}) \quad (5)$$

which we can sample from in $\mathcal{O}(Np^2)$ time if we have a pre-computed Cholesky decomposition of \mathbf{Q} . This is computationally cheap compared to optimal solution which is $\mathcal{O}(Np^2 + p^3 + n_{\max} p^2)$ but it is not optimal.

Backward filter (Algorithm 3)

We need to specify the artificial prior $\gamma_t(\boldsymbol{\alpha}_t)$ for our artificial backward distribution. Briers et al. (2009, page 69 and 70) provides recommendation on the selection. One suggestion is the artificial density function

$$\begin{aligned}\gamma_t(\boldsymbol{\alpha}_t) &= \phi\left(\boldsymbol{\alpha}_t \mid \overleftarrow{\mathbf{m}}_t, \overleftarrow{\mathbf{P}}_t\right) \\ \overleftarrow{\mathbf{m}}_t &= \mathbf{F}^t \boldsymbol{\mu}_0 \\ \overleftarrow{\mathbf{P}}_t &= \begin{cases} \mathbf{Q}_0 & t = 0 \\ \mathbf{F} \overleftarrow{\mathbf{P}}_{t-1} \mathbf{F}^\top + \mathbf{Q} & t > 0 \end{cases}\end{aligned}\quad (6)$$

The backward arrows are added to stress that these are means and covariance matrices which we use in the artificial distribution we target in the backward particle filter. The artificial distribution we target in backward particle filters has the following conditional density functions

$$\begin{aligned}\tilde{p}(\boldsymbol{\alpha}_{t:d} \mid \mathbf{y}_{t:d}) &\propto \gamma_t(\boldsymbol{\alpha}_t) \prod_{i=t}^d g_i(\mathbf{y}_i \mid \boldsymbol{\alpha}_i) \prod_{i=t}^{d-1} f(\boldsymbol{\alpha}_{i+1} \mid \boldsymbol{\alpha}_i) \\ \tilde{p}(\boldsymbol{\alpha}_t \mid \mathbf{y}_{(t+1):d}) &\propto \gamma_t(\boldsymbol{\alpha}_t) \int \tilde{p}(\boldsymbol{\alpha}_{t+1} \mid \mathbf{y}_{(t+1):d}) \frac{f(\boldsymbol{\alpha}_{t+1} \mid \boldsymbol{\alpha}_t)}{\gamma_{t+1}(\boldsymbol{\alpha}_{t+1})} d\boldsymbol{\alpha}_{t+1} \\ \tilde{p}(\boldsymbol{\alpha}_t \mid \mathbf{y}_{t:d}) &\propto g_t(\mathbf{y}_t \mid \boldsymbol{\alpha}_t) \tilde{p}(\boldsymbol{\alpha}_t \mid \mathbf{y}_{(t+1):d}) \\ \tilde{p}(\boldsymbol{\alpha}_t \mid \boldsymbol{\alpha}_{t+1}) &= \frac{f(\boldsymbol{\alpha}_{t+1} \mid \boldsymbol{\alpha}_t) \gamma_t(\boldsymbol{\alpha}_t)}{\gamma_{t+1}(\boldsymbol{\alpha}_{t+1})}\end{aligned}\quad (7)$$

where we have left out some of the normalization constants. Sampling from this artificial distribution turns out to be useful as it gives us an approximation of a conditional density we need up to a constant (see Section 7). To derive the re-sampling weight, we first find an expression for the density $\tilde{p}(\boldsymbol{\alpha}_t \mid \boldsymbol{\alpha}_{t+1})$. We can observe that

$$\begin{aligned}\log \tilde{p}(\boldsymbol{\alpha}_t \mid \boldsymbol{\alpha}_{t+1}) &= \log f(\boldsymbol{\alpha}_t \mid \boldsymbol{\alpha}_{t+1}) + \log \gamma_t(\boldsymbol{\alpha}_t) + \dots \\ &= -\frac{1}{2} \boldsymbol{\alpha}_t^\top \overleftarrow{\mathbf{S}}_t^{-1} \boldsymbol{\alpha}_t - \boldsymbol{\alpha}_t^\top \overleftarrow{\mathbf{S}}_t^{-1} \overleftarrow{\mathbf{a}}_t(\boldsymbol{\alpha}_{t+1}) + \dots \\ \overleftarrow{\mathbf{S}}_t &= (\mathbf{P}_t^{-1} + \mathbf{F}^\top \mathbf{Q}^{-1} \mathbf{F})^{-1} \\ \overleftarrow{\mathbf{a}}_t(\mathbf{x}) &= \overleftarrow{\mathbf{S}}_t (\mathbf{Q}^{-1} \mathbf{m}_t + \mathbf{F}^\top \mathbf{Q}^{-1} \mathbf{x})\end{aligned}$$

so

$$\tilde{p}(\boldsymbol{\alpha}_t \mid \boldsymbol{\alpha}_{t+1}) = \phi\left(\boldsymbol{\alpha}_t \mid \overleftarrow{\mathbf{a}}_t(\boldsymbol{\alpha}_{t+1}), \overleftarrow{\mathbf{S}}_t\right)\quad (8)$$

As shown by Fearnhead et al. (2010), we can show that

$$\begin{aligned}\overleftarrow{\mathbf{S}}_t &= \overleftarrow{\mathbf{P}}_t \mathbf{F}^\top \overleftarrow{\mathbf{P}}_{t+1}^{-1} \mathbf{Q} \mathbf{F}^{-\top} \\ \overleftarrow{\mathbf{a}}_t(\mathbf{x}) &= \overleftarrow{\mathbf{P}}_t \mathbf{F}^\top \overleftarrow{\mathbf{P}}_{t+1}^{-1} \mathbf{x} + \overleftarrow{\mathbf{S}}_t \overleftarrow{\mathbf{P}}_t^{-1} \overleftarrow{\mathbf{m}}_t\end{aligned}\quad (9)$$

The first equality can be shown by

$$\begin{aligned}
& \left(\overleftarrow{\mathbf{P}}_t \mathbf{F}^\top \overleftarrow{\mathbf{P}}_{t+1}^{-1} \mathbf{Q} \mathbf{F}^{-\top} \right)^{-1} \left(\mathbf{P}_t^{-1} + \mathbf{F}^\top \mathbf{Q}^{-1} \mathbf{F} \right)^{-1} \\
&= \mathbf{F}^\top \mathbf{Q}^{-1} \overleftarrow{\mathbf{P}}_{t+1} \mathbf{F}^{-\top} \overleftarrow{\mathbf{P}}_t^{-1} \left(\mathbf{P}_t^{-1} + \mathbf{F}^\top \mathbf{Q} \mathbf{F} \right)^{-1} \\
&= \mathbf{F}^\top \mathbf{Q}^{-1} \overleftarrow{\mathbf{P}}_{t+1} \mathbf{F}^{-\top} \overleftarrow{\mathbf{P}}_t^{-1} \left(\overleftarrow{\mathbf{P}}_t - \overleftarrow{\mathbf{P}}_t \mathbf{F}^\top \left(\mathbf{Q} + \mathbf{F} \overleftarrow{\mathbf{P}}_t \mathbf{F}^\top \right)^{-1} \mathbf{F} \overleftarrow{\mathbf{P}}_t \right) \\
&= \mathbf{F}^\top \mathbf{Q}^{-1} \overleftarrow{\mathbf{P}}_{t+1} \mathbf{F}^{-\top} \left(\mathbf{I} - \mathbf{F}^\top \overleftarrow{\mathbf{P}}_{t+1}^{-1} \mathbf{F} \overleftarrow{\mathbf{P}}_t \right) \\
&= \mathbf{F}^\top \mathbf{Q}^{-1} \overleftarrow{\mathbf{P}}_{t+1} \mathbf{F}^{-\top} - \mathbf{F}^\top \mathbf{Q}^{-1} \mathbf{F} \overleftarrow{\mathbf{P}}_t \\
&= \mathbf{F}^\top \mathbf{Q}^{-1} \left(\mathbf{F} \overleftarrow{\mathbf{P}}_t \mathbf{F}^\top + \mathbf{Q} \right) \mathbf{F}^{-\top} - \mathbf{F}^\top \mathbf{Q}^{-1} \mathbf{F} \overleftarrow{\mathbf{P}}_t \\
&= \mathbf{I}
\end{aligned}$$

where we assume that all matrices are non-singular and we use the Woodbury matrix identity. Similar arguments can be used for $\overleftarrow{\mathbf{a}}_t(\mathbf{x})$. Using the above, we can find that the optimal re-sampling weights are

$$\begin{aligned}
\tilde{\beta}_t^{(k)} &\propto \tilde{p} \left(\mathbf{y}_t \mid \tilde{\alpha}_{t+1}^{(k)} \right) \tilde{w}_{t+1}^{(k)} \\
&\propto \int g_t(\mathbf{y}_t \mid \boldsymbol{\alpha}_t) \tilde{p} \left(\boldsymbol{\alpha}_t \mid \tilde{\alpha}_{t+1}^{(k)} \right) d\boldsymbol{\alpha}_t \tilde{w}_{t+1}^{(k)} \\
&= \phi \left(\mathbf{y}_t \mid \mathbf{X}_t \overleftarrow{\mathbf{a}}_t(\tilde{\alpha}_{t+1}^{(k)}) + \mathbf{h}_t, \mathbf{X}_t \overleftarrow{\mathbf{S}}_t \mathbf{X}_t^\top + \mathbf{H}_t \right) \tilde{w}_{t+1}^{(k)}
\end{aligned} \tag{10}$$

or we can set the re-sampling weights proportional to $\tilde{\beta}_t^{(k)} \propto \tilde{w}_{t+1}^{(k)}$ and get a sequential importance re-sampling like algorithm. As for the proposal distribution, the optimal density is

$$\begin{aligned}
\log \tilde{q} \left(\boldsymbol{\alpha}_t \mid \mathbf{y}_t, \tilde{\alpha}_{t+1}^{(k)} \right) &= \log \tilde{p} \left(\boldsymbol{\alpha}_t, \tilde{\alpha}_{t+1}^{(k)}, \mathbf{y}_t \right) + \dots \\
&= \log \gamma_t(\boldsymbol{\alpha}_t) + \log g_t(\mathbf{y}_t \mid \boldsymbol{\alpha}_t) + \log f \left(\boldsymbol{\alpha}_{t+1}^{(k)} \mid \boldsymbol{\alpha}_t \right) \\
&= -\frac{1}{2} \boldsymbol{\alpha}_t^\top \overleftarrow{\mathbf{\Sigma}}_t^{-1} \boldsymbol{\alpha}_t + \boldsymbol{\alpha}_t^\top \overleftarrow{\mathbf{\Sigma}}_t^{-1} \overleftarrow{\boldsymbol{\mu}}_t(\tilde{\alpha}_{t+1}^{(k)}) + \dots \\
\overleftarrow{\mathbf{\Sigma}}_t &= \left(\mathbf{P}_t^{-1} + \mathbf{F}^\top \mathbf{Q}^{-1} \mathbf{F} + \mathbf{X}_t^\top \mathbf{H}_t^{-1} \mathbf{X}_t \right)^{-1} \\
\overleftarrow{\boldsymbol{\mu}}_t(\mathbf{x}) &= \mathbf{\Sigma}_t \left(\mathbf{P}_t^{-1} \mathbf{m}_t + \mathbf{F}^\top \mathbf{Q}^{-1} \mathbf{x} + \mathbf{X}_t^\top \mathbf{H}_t^{-1} (\mathbf{y}_t - \mathbf{h}_t) \right)
\end{aligned}$$

Thus, we set

$$\tilde{q} \left(\boldsymbol{\alpha}_t \mid \mathbf{y}_t, \tilde{\alpha}_{t+1}^{(k)} \right) = \phi \left(\boldsymbol{\alpha}_t \mid \overleftarrow{\boldsymbol{\mu}}_t(\tilde{\alpha}_{t+1}^{(k)}), \overleftarrow{\mathbf{\Sigma}}_t \right)$$

A computationally simpler but non-optimal option is to use a bootstrap filter like method and set

$$\tilde{q} \left(\boldsymbol{\alpha}_t \mid \mathbf{y}_t, \tilde{\alpha}_{t+1}^{(k)} \right) = \tilde{p} \left(\boldsymbol{\alpha}_t \mid \tilde{\alpha}_{t+1}^{(k)} \right) \tag{11}$$

Combining / smoothing (Algorithm 1)

We end this example with the conditional Gaussian observable outcome model with the proposal distribution needed for Algorithm 1. We want to select

$$\begin{aligned} q\left(\boldsymbol{\alpha}_t \mid \boldsymbol{\alpha}_{t-1}^{(j)}, \mathbf{y}_t, \tilde{\boldsymbol{\alpha}}_{t+1}^{(k)}\right) &= p\left(\boldsymbol{\alpha}_t \mid \boldsymbol{\alpha}_{t-1}^{(j)}, \mathbf{y}_t, \tilde{\boldsymbol{\alpha}}_{t+1}^{(k)}\right) \\ &\propto g\left(\mathbf{y}_t \mid \boldsymbol{\alpha}_t\right) f\left(\boldsymbol{\alpha}_t \mid \boldsymbol{\alpha}_{t-1}^{(j)}\right) f\left(\tilde{\boldsymbol{\alpha}}_{t+1}^{(k)} \mid \boldsymbol{\alpha}_t\right) \end{aligned}$$

Looking at the log density as we did before, we find that

$$\begin{aligned} \log q\left(\boldsymbol{\alpha}_t \mid \boldsymbol{\alpha}_{t-1}^{(j)}, \mathbf{y}_t, \tilde{\boldsymbol{\alpha}}_{t+1}^{(k)}\right) &= \log g\left(\mathbf{y}_t \mid \boldsymbol{\alpha}_t\right) + \log f\left(\boldsymbol{\alpha}_t \mid \boldsymbol{\alpha}_{t-1}^{(j)}\right) + \log f\left(\tilde{\boldsymbol{\alpha}}_{t+1}^{(k)} \mid \boldsymbol{\alpha}_t\right) + \dots \\ &= -\frac{1}{2} \boldsymbol{\alpha}_t^\top \overleftarrow{\boldsymbol{\Sigma}}_t^{-1} \boldsymbol{\alpha}_t + \boldsymbol{\alpha}_t^\top \overleftarrow{\boldsymbol{\Sigma}}_t^{-1} \overleftarrow{\boldsymbol{\mu}}_t\left(\boldsymbol{\alpha}_{t-1}^{(j)}, \tilde{\boldsymbol{\alpha}}_{t+1}^{(k)}\right) + \dots \\ \overleftarrow{\boldsymbol{\Sigma}}_t &= \left(\mathbf{Q}^{-1} + \mathbf{F}^\top \mathbf{Q}^{-1} \mathbf{F} + \mathbf{X}_t^\top \mathbf{H}^{-1} \mathbf{X}_t\right)^{-1} \end{aligned} \tag{12}$$

$$\overleftarrow{\boldsymbol{\mu}}_t(\mathbf{x}, \overleftarrow{\mathbf{x}}) = \boldsymbol{\Sigma}_t \left(\mathbf{Q}^{-1} \mathbf{F} \mathbf{x} + \mathbf{F}^\top \mathbf{Q}^{-1} \overleftarrow{\mathbf{x}} + \mathbf{X}_t^\top \mathbf{H}_t^{-1} (\mathbf{y}_t - \mathbf{h}_t)\right) \tag{13}$$

so that

$$q\left(\boldsymbol{\alpha}_t \mid \boldsymbol{\alpha}_{t-1}^{(j)}, \mathbf{y}_t, \tilde{\boldsymbol{\alpha}}_{t+1}^{(k)}\right) = \phi\left(\boldsymbol{\alpha}_t \mid \overleftarrow{\boldsymbol{\mu}}_t\left(\boldsymbol{\alpha}_{t-1}^{(j)}, \tilde{\boldsymbol{\alpha}}_{t+1}^{(k)}\right), \boldsymbol{\Sigma}_t\right)$$

Alternatively, we can use a bootstrap filter like proposal distribution with

$$\begin{aligned} \overleftarrow{\boldsymbol{\Sigma}}_t &= \left(\mathbf{Q}^{-1} + \mathbf{F}^\top \mathbf{Q}^{-1} \mathbf{F}\right)^{-1} \\ \overleftarrow{\boldsymbol{\mu}}_t(\mathbf{x}, \overleftarrow{\mathbf{x}}) &= \boldsymbol{\Sigma}_t \left(\mathbf{Q}^{-1} \mathbf{F} \mathbf{x} + \mathbf{F}^\top \mathbf{Q}^{-1} \overleftarrow{\mathbf{x}}\right) \end{aligned} \tag{14}$$

This is not optimal but faster.

Algorithm 1 $\mathcal{O}(N)$ particle smoother using the method in Fearnhead et al. (2010).

Input:

$\mathbf{Q}, \mathbf{Q}_0, \mathbf{a}_0, \mathbf{X}_1, \dots, \mathbf{X}_d, \mathbf{Z}_1, \dots, \mathbf{Z}_d, \mathbf{o}_1, \dots, \mathbf{o}_d, \mathbf{y}_1, \dots, \mathbf{y}_d, R_1, \dots, R_d, \boldsymbol{\omega}$
 Proposal distribution with density

$$q\left(\boldsymbol{\alpha}_t \mid \boldsymbol{\alpha}_{t-1}^{(j)}, \mathbf{y}_t, \tilde{\boldsymbol{\alpha}}_{t+1}^{(k)}\right) \quad (15)$$

- 1: **procedure** FILTER FORWARD
- 2: Run a forward particle filter to get particle clouds $\left\{\boldsymbol{\alpha}_t^{(j)}, w_t^{(j)}, \beta_{t+1}^{(j)}\right\}_{j=1, \dots, N}$ approximating the density $p(\boldsymbol{\alpha}_t \mid \mathbf{y}_{1:t})$ for $t = 0, 1, \dots, d$. See Algorithm 2.
- 3: **procedure** FILTER BACKWARDS
- 4: Run a similar backward filter to get $\left\{\tilde{\boldsymbol{\alpha}}_t^{(k)}, \tilde{w}_t^{(k)}, \tilde{\beta}_{t-1}^{(k)}\right\}_{k=1, \dots, d}$ approximating the artificial density $\tilde{p}(\boldsymbol{\alpha}_t \mid \mathbf{y}_{t:d})$ for $t = d+1, d, d-1, \dots, 1$. See Algorithm 3.
- 5: **procedure** SMOOTH (COMBINE)
- 6: **for** $t = 1, \dots, d$ **do**
Re-sample
- 7: Sample $i = 1, 2, \dots, N_s$ pairs of $(j_i, k_i) \in N^2$ where each component is independently sampled using re-sampling weights $\beta_t^{(j)}$ and $\tilde{\beta}_t^{(k)}$.
Propagate
- 8: Sample particles $\hat{\boldsymbol{\alpha}}_t^{(i)}$ from the proposal distribution $\tilde{q}\left(\cdot \mid \boldsymbol{\alpha}_{t-1}^{(j_i)}, \mathbf{y}_t, \tilde{\boldsymbol{\alpha}}_{t+1}^{(k_i)}\right)$.
Re-weight
- 9: Assign each particle a weight

$$\hat{w}_t^{(i)} \propto \frac{f\left(\hat{\boldsymbol{\alpha}}_t^{(i)} \mid \boldsymbol{\alpha}_{t-1}^{(j_i)}\right) g_t\left(\mathbf{y}_t \mid \hat{\boldsymbol{\alpha}}_t^{(i)}\right) f\left(\tilde{\boldsymbol{\alpha}}_{t+1}^{(k_i)} \mid \hat{\boldsymbol{\alpha}}_t^{(i)}\right) w_{t-1}^{(j_i)} \tilde{w}_{t+1}^{(k_i)}}{\tilde{q}\left(\hat{\boldsymbol{\alpha}}_t^{(i)} \mid \boldsymbol{\alpha}_{t-1}^{(j_i)}, \mathbf{y}_t, \tilde{\boldsymbol{\alpha}}_{t+1}^{(k_i)}\right) \beta_t^{(j_i)} \tilde{\beta}_t^{(k_i)} \gamma_{t+1}\left(\tilde{\boldsymbol{\alpha}}_{t+1}^{(k_i)}\right)} \quad (16)$$

Algorithm 2 Forward filter as in Pitt and Shephard (1999). It is equivalent with Doucet and Johansen (2009, page 20 and 25). The version and notation below is from Fearnhead et al. (2010, page 449).

Input:

Proposal distribution with density

$$q\left(\boldsymbol{\alpha}_t \mid \boldsymbol{\alpha}_{t-1}^{(j)}, \mathbf{y}_t\right)$$

Function h to compute re-sampling weights

$$\beta_t^{(j)} \propto h(\mathbf{y}_t, \boldsymbol{\alpha}_{t-1}^{(j)}) w_{t-1}^{(j)}$$

- 1: Sample $\boldsymbol{\alpha}_0^{(1)}, \dots, \boldsymbol{\alpha}_0^{(N_f)}$ particles from $\mathcal{N}(\boldsymbol{\mu}_0, \mathbf{Q}_0)$ and set the weights $w_0^{(1)}, \dots, w_0^{(N_f)}$ to $1/N_f$.
- 2: **for** $t = 1, \dots, d$ **do**
- 3: **procedure** RE-SAMPLE
- 4: Compute re-sampling weights $\beta_t^{(j)}$ using h and re-sample according to $\beta_t^{(j)}$ to get indices j_1, \dots, j_N . If we do not re-sample then set $\beta_t^{(j)} = 1/N$ or $1/N_f$ at time $t = 1$.
- 5: **procedure** PROPAGATE
- 6: Sample new particles $\boldsymbol{\alpha}_t^{(i)}$ using the proposal distribution $q\left(\boldsymbol{\alpha}_t \mid \boldsymbol{\alpha}_{t-1}^{(j_i)}, \mathbf{y}_t\right)$.
- 7: **procedure** RE-WEIGHT
- 8: Re-weight particles using

$$w_t^{(i)} \propto \frac{g_t\left(\mathbf{y}_t \mid \boldsymbol{\alpha}_t^{(i)}\right) f\left(\boldsymbol{\alpha}_t^{(i)} \mid \boldsymbol{\alpha}_{t-1}^{(j_i)}\right) w_{t-1}^{(j_i)}}{q\left(\boldsymbol{\alpha}_t^{(i)} \mid \boldsymbol{\alpha}_{t-1}^{(j_i)}, \mathbf{y}_t\right) \beta_t^{(j_i)}} \quad (17)$$

Algorithm 3 Backwards filter. See Briers et al. (2009) and Fearnhead et al. (2010).

Input:

An artificial distribution

$$\tilde{p}(\boldsymbol{\alpha}_t | \mathbf{y}_{t:d}) \propto \gamma_t(\boldsymbol{\alpha}_t) p(\mathbf{y}_{t:d} | \boldsymbol{\alpha}_t) \quad (18)$$

with an artificial prior distribution $\gamma_t(\boldsymbol{\alpha}_t)$.

Proposal distribution

$$\tilde{q}(\boldsymbol{\alpha}_t | \mathbf{y}_t, \tilde{\boldsymbol{\alpha}}_{t+1}^{(k)})$$

Function h to compute re-sampling weights

$$\tilde{\beta}_t^{(k)} \propto h(\mathbf{y}_t, \tilde{\boldsymbol{\alpha}}_{t+1}^{(k)}) \tilde{w}_{t+1}^{(k)}$$

- 1: Sample $\tilde{\boldsymbol{\alpha}}_{d+1}^{(1)}, \dots, \tilde{\boldsymbol{\alpha}}_{d+1}^{(N_f)}$ particles from $\gamma_{d+1}(\cdot)$ and set the weights $\tilde{w}_{d+1}^{(1)}, \dots, \tilde{w}_{d+1}^{(N_f)}$ to $1/N_f$.
- 2: **for** $t = d, \dots, 1$ **do**
- 3: **procedure** RE-SAMPLE
- 4: Compute re-sampling weights $\tilde{\beta}_t^{(k)}$ using h and re-sample according to $\tilde{\beta}_t^{(k)}$ to get indices k_1, \dots, k_N . If we do not re-sample then set $\tilde{\beta}_t^{(k)} = 1/N$ or $1/N_f$ at time $t = d$.
- 5: **procedure** PROPAGATE
- 6: Sample new particles $\tilde{\boldsymbol{\alpha}}_t^{(i)}$ using the proposal distribution $\tilde{q}(\boldsymbol{\alpha}_t | \tilde{\boldsymbol{\alpha}}_{t+1}^{(k_i)}, \mathbf{y}_t)$.
- 7: **procedure** RE-WEIGHT
- 8: Re-weight particles using

$$\tilde{w}_t^{(i)} \propto \frac{g_t(\mathbf{y}_t | \tilde{\boldsymbol{\alpha}}_t^{(i)}) f(\tilde{\boldsymbol{\alpha}}_{t+1}^{(k_i)} | \tilde{\boldsymbol{\alpha}}_t^{(i)}) \gamma_t(\tilde{\boldsymbol{\alpha}}_t^{(i)}) \tilde{w}_{t+1}^{(k_i)}}{q(\tilde{\boldsymbol{\alpha}}_t^{(i)} | \tilde{\boldsymbol{\alpha}}_{t+1}^{(k_i)}, \mathbf{y}_t) \gamma_{t+1}(\tilde{\boldsymbol{\alpha}}_{t+1}^{(k_i)}) \beta_t^{(k_i)}} \quad (19)$$

2 Non-linear conditional observation model

If we go back to the model in Equation (1) then $\mathbf{y}_t \mid \boldsymbol{\alpha}_t$ is not a multivariate normal distribution for the implemented models. In this case, we have no closed form solutions for the optimal re-sampling weights, and we do not know the following conditional distributions: $\boldsymbol{\alpha}_t \mid \mathbf{y}_t, \boldsymbol{\alpha}_{t-1}$, $\boldsymbol{\alpha}_t \mid \mathbf{y}_t, \boldsymbol{\alpha}_{t+1}$ (in the artificial distribution $\tilde{\mathbb{P}}$), and $\boldsymbol{\alpha}_t \mid \mathbf{y}_t, \boldsymbol{\alpha}_{t-1}, \boldsymbol{\alpha}_{t+1}$. However, assume that $g_t(\mathbf{y}_t \mid \boldsymbol{\alpha}_t)$ is log-concave in $\boldsymbol{\alpha}_t$. If this is true then it is easy to show that all of the previous three conditional distributions are unimodal. Hence, we can make a multivariate normal approximation as in Pitt and Shephard (1999). To do so, we make a second order Taylor expansion around some value \mathbf{z} to get

$$\begin{aligned} k_t(\boldsymbol{\alpha}_t) &= \log g_t(\mathbf{y}_t \mid \boldsymbol{\eta}(\boldsymbol{\alpha}_t)), & \boldsymbol{\eta}(\boldsymbol{\alpha}_t) &= \mathbf{X}_t \boldsymbol{\alpha}_t + \mathbf{h}_t \\ \log g_t(\mathbf{y}_t \mid \boldsymbol{\alpha}_t) &\approx \text{D}k_t(\mathbf{z})(\boldsymbol{\alpha}_t - \mathbf{z}) + \frac{1}{2}(\boldsymbol{\alpha}_t - \mathbf{z})^\top \text{H}k_t(\mathbf{z})(\boldsymbol{\alpha}_t - \mathbf{z}) + \dots \\ &= \boldsymbol{\alpha}_t^\top \text{D}k_t(\mathbf{z})^\top - \frac{1}{2}(\boldsymbol{\alpha}_t - \mathbf{z})^\top (-\text{H}k_t(\mathbf{z}))(\boldsymbol{\alpha}_t - \mathbf{z}) + \dots \\ &= \boldsymbol{\alpha}_t^\top (-\text{H}k_t(\mathbf{z}))(\mathbf{z} - \text{H}k_t(\mathbf{z})^{-1} \text{D}k_t(\mathbf{z})^\top) - \frac{1}{2} \boldsymbol{\alpha}_t^\top (-\text{H}k_t(\mathbf{z})) \boldsymbol{\alpha}_t + \dots \end{aligned}$$

where \dots includes the zero order term, $\text{D}k_t$ is the Jacobian, and $\text{H}k_t$ denotes the Hessian. I add a subscript to D and H to indicate when Jacobian or Hessian is with respect to a given variable. We still assume that we use a first order state space model such that $r = p$. We notice that

$$\begin{aligned} \text{H}k_t(\mathbf{z}) &= \text{D}_{\boldsymbol{\alpha}_t} \boldsymbol{\eta}(\mathbf{z})^\top \text{H}_{\boldsymbol{\eta}} \log g_t(\mathbf{y}_t \mid \boldsymbol{\eta}(\mathbf{z})) \text{D}_{\boldsymbol{\alpha}_t} \boldsymbol{\eta}(\mathbf{z}) \\ &= \mathbf{X}_t^\top (-\mathbf{G}_t(\mathbf{z})) \mathbf{X}_t, & \mathbf{G}_t(\mathbf{z}) &= -\text{H}_{\boldsymbol{\eta}} \log g_t(\mathbf{y}_t \mid \boldsymbol{\eta}(\mathbf{z})) \end{aligned}$$

which follows from the chain rule and we use that $\text{H}_{\boldsymbol{\alpha}_t} \boldsymbol{\eta}(\mathbf{z}) = \mathbf{0}$. Thus,

$$\begin{aligned} \log g_t(\mathbf{y}_t \mid \boldsymbol{\alpha}_t) &\approx \boldsymbol{\alpha}_t^\top \mathbf{X}_t^\top \mathbf{G}_t(\mathbf{z}) \mathbf{u}_t(\mathbf{z}) - \frac{1}{2} \boldsymbol{\alpha}_t^\top \mathbf{X}_t^\top \mathbf{G}_t(\mathbf{z}) \mathbf{X}_t \boldsymbol{\alpha}_t \\ \mathbf{u}_t(\mathbf{z}) &= \mathbf{X}_t \mathbf{z} - \mathbf{X}_t \text{H}k_t(\mathbf{z})^{-1} \text{D}k_t(\mathbf{z})^\top \end{aligned}$$

This yields the following multivariate normal approximation

$$g_t(\mathbf{y}_t \mid \boldsymbol{\alpha}_t) \approx \phi(\mathbf{X}_t \boldsymbol{\alpha}_t \mid \mathbf{u}_t(\mathbf{z}), \mathbf{G}_t(\mathbf{z})^{-1})$$

The Taylor approximation is easily used in the proposal distributions. E.g., for given \mathbf{z} , we get the following mean and covariance matrix analogues to Equation (3) and (4) in the proposal distribution in the forward particle filter

$$\begin{aligned} \boldsymbol{\Sigma}_t(\mathbf{z}) &= (\mathbf{Q}^{-1} + \mathbf{X}_t^\top \mathbf{G}_t(\mathbf{z}) \mathbf{X}_t)^{-1} \\ \boldsymbol{\mu}_t(\mathbf{x}, \mathbf{z}) &= \boldsymbol{\Sigma}_t(\mathbf{z}) (\mathbf{Q}^{-1} \mathbf{F} \mathbf{x} + \mathbf{X}_t^\top \mathbf{G}_t(\mathbf{z}) \mathbf{u}_t(\mathbf{z})) \end{aligned}$$

As for the re-sampling weights, we can use

$$\begin{aligned}
\hat{\boldsymbol{\alpha}} &= \boldsymbol{\mu}_t(\boldsymbol{\alpha}_{t-1}^{(j)}, \mathbf{z}) \\
\beta_t^{(j)} &\propto p\left(\mathbf{y}_t \mid \boldsymbol{\alpha}_{t-1}^{(j)}\right) w_{t-1}^{(j)} \\
&= \frac{p\left(\mathbf{y}_t \mid \boldsymbol{\alpha}_{t-1}^{(j)}\right)}{g_t\left(\mathbf{y}_t \mid \hat{\boldsymbol{\alpha}}\right) f\left(\hat{\boldsymbol{\alpha}} \mid \boldsymbol{\alpha}_{t-1}^{(j)}\right)} g_t\left(\mathbf{y}_t \mid \hat{\boldsymbol{\alpha}}\right) f\left(\hat{\boldsymbol{\alpha}} \mid \boldsymbol{\alpha}_{t-1}^{(j)}\right) w_{t-1}^{(j)} \\
&= \frac{g_t\left(\mathbf{y}_t \mid \hat{\boldsymbol{\alpha}}\right) f\left(\hat{\boldsymbol{\alpha}} \mid \boldsymbol{\alpha}_{t-1}^{(j)}\right) w_{t-1}^{(j)}}{p\left(\hat{\boldsymbol{\alpha}} \mid \boldsymbol{\alpha}_{t-1}^{(j)}, \mathbf{y}_t\right)} \\
&\approx \frac{g_t\left(\mathbf{y}_t \mid \hat{\boldsymbol{\alpha}}\right) f\left(\hat{\boldsymbol{\alpha}} \mid \boldsymbol{\alpha}_{t-1}^{(j)}\right) w_{t-1}^{(j)}}{q\left(\hat{\boldsymbol{\alpha}} \mid \boldsymbol{\alpha}_{t-1}^{(j)}, \mathbf{y}_t\right)}
\end{aligned}$$

as in Fearnhead et al. (2010). We can approximate the backwards particle filter re-sampling weights in equation (10) in a similar way

$$\begin{aligned}
\tilde{\beta}_t^{(k)} &\propto \tilde{p}\left(\mathbf{y}_t \mid \tilde{\boldsymbol{\alpha}}_{t+1}^{(k)}\right) \tilde{w}_{t+1}^{(k)} \\
&\approx \frac{g_t\left(\mathbf{y}_t \mid \hat{\boldsymbol{\alpha}}\right) \tilde{p}\left(\hat{\boldsymbol{\alpha}} \mid \tilde{\boldsymbol{\alpha}}_{t+1}^{(k)}\right) \tilde{w}_{t+1}^{(k)}}{\tilde{q}\left(\hat{\boldsymbol{\alpha}} \mid \mathbf{y}_t, \tilde{\boldsymbol{\alpha}}_{t+1}^{(k)}\right)} \\
&= \frac{g_t\left(\mathbf{y}_t \mid \hat{\boldsymbol{\alpha}}\right) f\left(\tilde{\boldsymbol{\alpha}}_{t+1}^{(k)} \mid \hat{\boldsymbol{\alpha}}\right) \gamma_t\left(\hat{\boldsymbol{\alpha}}\right) \tilde{w}_{t+1}^{(k)}}{\tilde{q}\left(\hat{\boldsymbol{\alpha}} \mid \mathbf{y}_t, \tilde{\boldsymbol{\alpha}}_{t+1}^{(k)}\right) \gamma_{t+1}\left(\tilde{\boldsymbol{\alpha}}_{t+1}^{(k)}\right)} \tag{20}
\end{aligned}$$

$$\hat{\boldsymbol{\alpha}} = \overleftarrow{\boldsymbol{\mu}}\left(\tilde{\boldsymbol{\alpha}}_{t+1}^{(k)}, \mathbf{z}\right)$$

$$\overleftarrow{\boldsymbol{\mu}}(\mathbf{x}, \mathbf{z}) = \overleftarrow{\boldsymbol{\Sigma}}_t(\mathbf{z}) \left(\mathbf{P}_t^{-1} \mathbf{m}_t + \mathbf{F}^\top \mathbf{Q}^{-1} \mathbf{x} + \mathbf{X}_t^\top \mathbf{G}_t(\mathbf{z}) \mathbf{u}_t(\mathbf{z})\right) \tag{21}$$

$$\overleftarrow{\boldsymbol{\Sigma}}_t(\mathbf{z}) = \left(\mathbf{P}_t^{-1} + \mathbf{F}^\top \mathbf{Q}^{-1} \mathbf{F} + \mathbf{X}_t^\top \mathbf{G}_t(\mathbf{z}) \mathbf{X}_t\right)^{-1} \tag{22}$$

We may also use a multivariate t -distribution for the proposal distribution to get heavier tails than we do with the multivariate normal distribution.

2.1 Where to make the expansion

An options is to make the Taylor expansion at a mode for each particle or particle pair in the smoothing step. This yields

$$\begin{aligned}
\mathbf{z}^{(j)} &= \arg \max_{\mathbf{z}} g_t\left(\mathbf{y}_t \mid \mathbf{z}\right) f\left(\mathbf{z} \mid \boldsymbol{\alpha}_{t-1}^{(j)}\right) \\
\mathbf{z}^{(k)} &= \arg \max_{\mathbf{z}} g_t\left(\mathbf{y}_t \mid \mathbf{z}\right) \gamma_t(\mathbf{z}) f\left(\tilde{\boldsymbol{\alpha}}_{t+1}^{(k)} \mid \mathbf{z}\right) \\
\mathbf{z}^{(i)} &= \arg \max_{\mathbf{z}} f\left(\mathbf{z} \mid \boldsymbol{\alpha}_{t-1}^{(j_i)}\right) g_t\left(\mathbf{y}_t \mid \mathbf{z}\right) f\left(\tilde{\boldsymbol{\alpha}}_{t+1}^{(k_i)} \mid \mathbf{z}\right)
\end{aligned}$$

for respectively the forward particle filter, backward particle filter, and smoother. The downside is a $\mathcal{O}(r^2 n_{\max} N_S)$ or $\mathcal{O}(r^2 n_{\max} N)$ computational complexity at each time point as we have to evaluate $\mathbf{X}_t^\top \mathbf{G}_t(\mathbf{z}) \mathbf{X}_t$ for each particle or particle pair. Instead we can make the approximation once at each time point at respectively $\sum_{i=1}^N w_{t-1}^{(j)} \boldsymbol{\alpha}_{t-1}^{(j)}$, $\sum_{i=1}^N \tilde{w}_{t+1}^{(j)} \tilde{\boldsymbol{\alpha}}_{t+1}^{(j)}$, and

$$(\mathbf{Q}^{-1} + \mathbf{F}^\top \mathbf{Q}^{-1} \mathbf{F})^{-1} \left(\mathbf{Q}^{-1} \mathbf{F} \sum_{i=1}^N w_{t-1}^{(j)} \boldsymbol{\alpha}_{t-1}^{(j)} + \mathbf{F}^\top \mathbf{Q}^{-1} \sum_{i=1}^N \tilde{w}_{t+1}^{(j)} \tilde{\boldsymbol{\alpha}}_{t+1}^{(j)} \right)$$

which will reduce the computational complexity at each time point to $\mathcal{O}(r n_{\max} N_S + r p n_{\max})$ or $\mathcal{O}(r n_{\max} N + r p n_{\max})$.

3 Higher order state-space models

Now, we will consider the case where $p > r$. Currently this is not supported in the package but may be in the future. An example is a second order vector auto-regression, VAR(2), with $2r = p$ and

$$\mathbf{F} = \begin{pmatrix} \mathbf{F}_1 & \mathbf{F}_2 \\ \mathbf{I}_r & \mathbf{0} \end{pmatrix}, \quad \mathbf{F}_i \in \mathbb{R}^{r \times r}$$

$$\boldsymbol{\alpha}_t = (\boldsymbol{\xi}_t^\top \quad \boldsymbol{\xi}_{t-1}^\top)^\top$$

Here

$$f(\boldsymbol{\alpha}_t | \boldsymbol{\alpha}_{t-1}) = \delta_{\mathbf{K}\mathbf{F}\boldsymbol{\alpha}_{t-1}}(\mathbf{K}\boldsymbol{\alpha}_t) \phi(\mathbf{R}^+ \boldsymbol{\alpha}_t | \mathbf{R}^+ \mathbf{F} \boldsymbol{\alpha}_{t-1}, \mathbf{Q}), \quad \mathbf{K} = \begin{pmatrix} \mathbf{0} \\ \mathbf{I}_{p-r} \end{pmatrix}$$

$$= \delta_{\boldsymbol{\xi}_{t-1}}(\mathbf{K}\boldsymbol{\alpha}_t) \phi(\boldsymbol{\xi}_t | \mathbf{R}^+ \mathbf{F} \boldsymbol{\alpha}_{t-1}, \mathbf{Q})$$

the second equality follows in a second order vector auto-regression. This is easily implemented in the forward particle filter by sampling $\boldsymbol{\xi}_t | \boldsymbol{\alpha}_{t-1}$ and setting the remaining $p - r$ variables of $\boldsymbol{\alpha}_t$ to $\mathbf{K}\mathbf{F}\boldsymbol{\alpha}_{t-1}$ (the last $p - r$ rows of $\mathbf{F}\boldsymbol{\alpha}_{t-1}$). It is not as easy for the backward filter. To see this, consider the artificial transition density in Equation (7)

$$\tilde{p}(\boldsymbol{\alpha}_t | \boldsymbol{\alpha}_{t+1}) = \frac{f(\boldsymbol{\alpha}_{t+1} | \boldsymbol{\alpha}_t) \gamma_t(\boldsymbol{\alpha}_t)}{\gamma_{t+1}(\boldsymbol{\alpha}_{t+1})}$$

$$= \frac{\delta_{\mathbf{K}\mathbf{F}\boldsymbol{\alpha}_t}(\mathbf{K}\boldsymbol{\alpha}_{t+1}) \phi(\mathbf{R}^+ \boldsymbol{\alpha}_{t+1} | \mathbf{R}^+ \mathbf{F} \boldsymbol{\alpha}_t, \mathbf{Q}) \gamma_t(\boldsymbol{\alpha}_t)}{\gamma_{t+1}(\boldsymbol{\alpha}_{t+1})}$$

where the last equality holds in this example. This distribution is obviously degenerate as it only has mass at the point $\boldsymbol{\xi}_t = \mathbf{K}\mathbf{F}\boldsymbol{\alpha}_t$ in the VAR(2) model.

3.1 Backward particle filter

What we can do in the backward particle filter in a VAR(o) model with $o > 1$ is to sample $\boldsymbol{\xi}_t$ at time t given $\tilde{\boldsymbol{\alpha}}_{t+o}^{(k)}$. Thus, we start Algorithm 3 by sampling $\boldsymbol{\alpha}_{d+o}$. Further, we change

the artificial prior distribution density in Equation (6) to

$$\begin{aligned}\gamma_t(\boldsymbol{\alpha}_t) &= \phi\left(\boldsymbol{\alpha}_t \mid \overleftarrow{\mathbf{m}}_t, \overleftarrow{\mathbf{P}}_t\right) \\ \gamma_t(\boldsymbol{\xi}_t) &= \phi\left(\boldsymbol{\alpha}_t \mid \mathbf{R}^+ \overleftarrow{\mathbf{m}}_t, \mathbf{R}^+ \overleftarrow{\mathbf{P}}_t \mathbf{R}^{+\top}\right) \\ \overleftarrow{\mathbf{m}}_t &= \mathbf{F}^t \boldsymbol{\mu}_0 \\ \overleftarrow{\mathbf{P}}_t &= \begin{cases} \mathbf{Q}_0 & t = 0 \\ \mathbf{F} \overleftarrow{\mathbf{P}}_{t-1} \mathbf{F}^\top + \mathbf{R} \mathbf{Q} \mathbf{R}^\top & t > 0 \end{cases}\end{aligned}$$

Next, we find the conditional distribution $\tilde{\boldsymbol{\alpha}}_{t+o}^{(k)} \mid \boldsymbol{\xi}_t$. To do so, we first find the joint distribution of $(\boldsymbol{\alpha}_{t+o}^\top, \boldsymbol{\xi}_t^\top)^\top$. We can find that

$$\begin{aligned}\boldsymbol{\xi}_k &= \mathbf{R}^+ \mathbf{F}^k \boldsymbol{\alpha}_0 + \sum_{i=1}^k \mathbf{G}(k-i) \boldsymbol{\epsilon}_i \\ \mathbf{G}(j) &= \begin{cases} \mathbf{I}_r & j = 0 \\ \sum_{i=1}^{\min(j,o)} \mathbf{F}_i \mathbf{G}(j-i) & j > 0 \end{cases}\end{aligned}$$

Thus,

$$\begin{aligned}\mathbf{E}(\boldsymbol{\xi}_t) &= \mathbf{R}^+ \mathbf{F}^t \boldsymbol{\mu}_0 = \mathbf{R}^+ \mathbf{m}_t \\ \mathbf{E}(\boldsymbol{\alpha}_t) &= \mathbf{m}_t \\ \text{Var}(\boldsymbol{\xi}_t) &= \sum_{i=1}^t \mathbf{G}(t-i) \mathbf{Q} \mathbf{G}(t-i)^\top + \mathbf{R}^+ \mathbf{F}^t \mathbf{Q}_0 \mathbf{F}^{t\top} \mathbf{R}^{+\top} \\ &= \mathbf{R}^+ \mathbf{P}_t \mathbf{R}^{+\top} \\ \text{Cov}(\boldsymbol{\xi}_k, \boldsymbol{\xi}_l) &= \sum_{i=1}^l \mathbf{G}(k-i) \mathbf{Q} \mathbf{G}(l-i)^\top + \mathbf{R}^+ \mathbf{F}^k \mathbf{Q}_0 \mathbf{F}^{l\top} \mathbf{R}^{+\top}, \quad k > l \\ \text{Var}(\boldsymbol{\alpha}_t) &= \mathbf{P}_t\end{aligned}$$

with which we can find that the joint distribution is

$$\begin{pmatrix} \boldsymbol{\alpha}_{t+o} \\ \boldsymbol{\xi}_t \end{pmatrix} \sim \mathcal{N}\left(\begin{pmatrix} \mathbf{m}_{t+o} \\ \mathbf{R}^+ \mathbf{m}_t \end{pmatrix}, \begin{pmatrix} \mathbf{P}_{t+o} & \text{Cov}(\boldsymbol{\alpha}_{t+o}, \boldsymbol{\xi}_t) \\ \text{Cov}(\boldsymbol{\xi}_t, \boldsymbol{\alpha}_{t+o}) & \mathbf{R}^+ \mathbf{P}_t \mathbf{R}^{+\top} \end{pmatrix}\right)$$

We are now able to compute

$$\begin{aligned}\mathbf{E}(\boldsymbol{\alpha}_{t+o} \mid \boldsymbol{\xi}_t) &= \boldsymbol{\mu}_{\boldsymbol{\alpha}_{t+o} \mid \boldsymbol{\xi}_t} = \mathbf{m}_{t+o} + \text{Cov}(\boldsymbol{\alpha}_{t+o}, \boldsymbol{\xi}_t) \mathbf{R}^\top \mathbf{P}_t^{-1} \mathbf{R} (\boldsymbol{\xi}_t - \mathbf{R}^+ \mathbf{m}_t) \\ &= \mathbf{v}_t + \mathbf{V}_t \boldsymbol{\xi}_t \\ \mathbf{V}_t &= \text{Cov}(\boldsymbol{\alpha}_{t+o}, \boldsymbol{\xi}_t) \mathbf{R}^\top \mathbf{P}_t^{-1} \mathbf{R} \\ \mathbf{v}_t &= \mathbf{m}_{t+o} - \mathbf{V}_t \mathbf{R}^+ \mathbf{m}_t\end{aligned}$$

$$\text{Var}(\boldsymbol{\alpha}_{t+o} \mid \boldsymbol{\xi}_t) = \boldsymbol{\Sigma}_{\boldsymbol{\alpha}_{t+o} \mid \boldsymbol{\xi}_t} = \mathbf{P}_{t+o} - \text{Cov}(\boldsymbol{\alpha}_{t+o}, \boldsymbol{\xi}_t) \mathbf{R}^\top \mathbf{P}_t^{-1} \mathbf{R} \text{Cov}(\boldsymbol{\xi}_t, \boldsymbol{\alpha}_{t+o})$$

Having found this conditional distribution then we can apply similar arguments as we did before and find the following mean and covariance matrix in the proposal distribution

$$\begin{aligned}\overleftarrow{\boldsymbol{\mu}}(\tilde{\boldsymbol{\alpha}}_{t+o}^{(k)}, \mathbf{z}) &= \overleftarrow{\boldsymbol{\Sigma}}_t(\mathbf{z}) \left(\mathbf{R}^\top \mathbf{P}_t^{-1} \mathbf{m}_t + \mathbf{V}_t^\top \boldsymbol{\Sigma}_{\boldsymbol{\alpha}_{t+o} \mid \boldsymbol{\xi}_t}^{-1} (\tilde{\boldsymbol{\alpha}}_{t+o}^{(k)} - \mathbf{v}_t) + \mathbf{X}_t^\top \mathbf{G}_t(\mathbf{z}) \mathbf{u}_t(\mathbf{z}) \right) \\ \overleftarrow{\boldsymbol{\Sigma}}_t(\mathbf{z}) &= \left(\mathbf{R}^\top \mathbf{P}_t^{-1} \mathbf{R} + \mathbf{V}_t^\top \boldsymbol{\Sigma}_{\boldsymbol{\alpha}_{t+o} \mid \boldsymbol{\xi}_t}^{-1} \mathbf{V}_t + \mathbf{X}_t^\top \mathbf{G}_t(\mathbf{z}) \mathbf{X}_t \right)^{-1}\end{aligned}$$

which replaces the mean and covariance matrix in Equation (21) and (22), $\mathbf{z}_t \in \mathbb{R}^r$ is the value of $\boldsymbol{\xi}_t$ at which we make the Taylor expansion, and \mathbf{G}_t and \mathbf{u}_t are defined in terms of $\boldsymbol{\xi}_t$. The new proposal distribution is used in Equation (20) when we compute the re-sampling weights where we change the γ_t and γ_{t+1} function. The two densities are changed to be $\gamma_t(\widehat{\boldsymbol{\xi}})$ in the numerator and $\gamma_{d+1}(\widetilde{\boldsymbol{\alpha}}_{d+1}^{(k)})$ or $\gamma_{t+1}(\widetilde{\boldsymbol{\xi}}_{t+1}^{(k)})$ at $t < d$ in the denominator where γ_t is understood as marginal density of $\boldsymbol{\xi}_t$ when xi is used instead of alpha.

3.2 Combining/smoothing

Similarly, in Algorithm 1 we sample pairs of particles $(\boldsymbol{\alpha}_{t-1}^{(j_i)}, \widetilde{\boldsymbol{\alpha}}_{t+o}^{(k_i)})$, and sample $\boldsymbol{\xi}_t$ given each pair. We can replace the mean and covariance matrix in the proposal distribution in Equation (13) and (12) with the approximation

$$\begin{aligned} \overleftrightarrow{\boldsymbol{\Sigma}}_t(\mathbf{z}) &= \left(\mathbf{Q}^{-1} + \mathbf{V}_t^\top \boldsymbol{\Sigma}_{\alpha_{t+o}|\boldsymbol{\xi}_t}^{-1} \mathbf{V}_t + \mathbf{X}_t^\top \mathbf{G}_t(\mathbf{z}) \mathbf{X}_t \right)^{-1} \\ \overleftrightarrow{\boldsymbol{\mu}}_t(\boldsymbol{\alpha}_{t-1}^{(j_i)}, \widetilde{\boldsymbol{\alpha}}_{t+o}^{(k_i)}, \mathbf{z}) &= \boldsymbol{\Sigma}_t(\mathbf{z}) \left(\mathbf{Q}^{-1} \mathbf{R}^+ \mathbf{F} \boldsymbol{\alpha}_{t-1}^{(j_i)} + \mathbf{V}_t^\top \boldsymbol{\Sigma}_{\alpha_{t+o}|\boldsymbol{\xi}_t}^{-1} (\widetilde{\boldsymbol{\alpha}}_{t+o}^{(k_i)} - \mathbf{v}_t) + \mathbf{X}_t^\top \mathbf{G}_t(\mathbf{z}) \mathbf{X}_t \right) \end{aligned}$$

where \mathbf{z} is the value of $\boldsymbol{\xi}_t$ that we make the Taylor expansion at.

4 Log likelihood evaluation

We can evaluate the log likelihood for a particular value of $\boldsymbol{\theta} = \{\mathbf{Q}, \mathbf{Q}_0, \boldsymbol{\mu}_0, \mathbf{F}\}$ as described in Doucet and Johansen (2009, page 5) and Malik and Pitt (2011, page 193) using the forward particle filter shown in Algorithm 2.

5 Parameter inference

In this section I first show an example of parameter estimation in the first order random walk using the EM-algorithm (Dempster et al., 1977). Then I cover the general vector auto-regression model and how one can estimate the fixed effects. Lastly, I will turn to estimation of the observed information matrix.

The formulas for parameter estimation for the first order random are particularly simple. We need to estimate \mathbf{Q} and \mathbf{a}_0 elements of $\boldsymbol{\varphi} = \{\mathbf{Q}, \mathbf{Q}_0, \boldsymbol{\mu}_0\}$. We do this by running Algorithm 1 for the current $\boldsymbol{\varphi}$. This yields the following quantities from the E-step

$$\begin{aligned} \mathbf{t}_t^{(\boldsymbol{\varphi})} &\approx \sum_{i=1}^{N_s} \widehat{\boldsymbol{\alpha}}_t^{(i)} \widehat{w}_t^{(i)} \\ \mathbf{T}_t^{(\boldsymbol{\varphi})} &\approx \sum_{i=1}^{N_s} \left(\widehat{\boldsymbol{\alpha}}_t^{(i)} - \mathbf{F} \boldsymbol{\alpha}_{t-1}^{(j_{it})} \right) \left(\widehat{\boldsymbol{\alpha}}_t^{(i)} - \mathbf{F} \boldsymbol{\alpha}_{t-1}^{(j_{it})} \right)^\top \widehat{w}_t^{(i)} \end{aligned} \tag{23}$$

where we have extended the notation in Algorithm 1 such that superscript j_{it} is the index from forward cloud at time $t-1$ matching with i th smoothed particle at time t . Then we carry out the M-step by updating $\boldsymbol{\mu}_0$ and \mathbf{Q} given the summary statistics above

$$\boldsymbol{\mu}_0 = \mathbf{t}_1^{(\boldsymbol{\varphi})} \quad \mathbf{Q} = \frac{1}{d-1} \sum_{t=2}^d \mathbf{R}^+ \mathbf{T}_t^{(\boldsymbol{\varphi})} \mathbf{R}^{+\top} \tag{24}$$

We then take another iteration of the EM algorithm with the new $\boldsymbol{\mu}_0$ and \mathbf{Q} till a convergence criteria is satisfied. See Kantas et al. (2015), Del Moral et al. (2010) and Schön et al. (2011) for further details on parameter estimation with particle filters.

5.1 Vector auto-regression models

We start by defining the following matrices to cover estimation in general vector auto-regression models for the latent space variable

$$\begin{aligned}\mathbf{N} &= \left(\hat{\boldsymbol{\alpha}}_2^{(1)}, \dots, \hat{\boldsymbol{\alpha}}_2^{(N_s)}, \hat{\boldsymbol{\alpha}}_3^{(1)}, \dots, \hat{\boldsymbol{\alpha}}_3^{(N_s)}, \hat{\boldsymbol{\alpha}}_4^{(1)}, \dots, \hat{\boldsymbol{\alpha}}_d^{(N_s)} \right)^\top \mathbf{R}^{+\top} \\ \mathbf{M} &= \left(\boldsymbol{\alpha}_1^{(j_{12})}, \dots, \boldsymbol{\alpha}_1^{(j_{N_s 2})}, \boldsymbol{\alpha}_2^{(j_{13})}, \dots, \boldsymbol{\alpha}_2^{(j_{N_s 3})}, \boldsymbol{\alpha}_3^{(j_{14})}, \dots, \boldsymbol{\alpha}_{d-1}^{(j_{N_s d})} \right)^\top \\ \mathbf{W} &= \text{diag} \left(\hat{w}_2^{(1)}, \dots, \hat{w}_2^{(N_s)}, \hat{w}_3^{(1)}, \dots, \hat{w}_3^{(N_s)}, \hat{w}_4^{(1)}, \dots, \hat{w}_d^{(N_s)} \right)\end{aligned}$$

where $\text{diag}(\cdot)$ is a diagonal matrix. We suppress the dependence above on the result of the E-step in a given iteration of the EM algorithm to ease the notation. The goal is to estimate \mathbf{F} and \mathbf{Q} in Equation (1). We can show that the M-step maximizers are

$$\hat{\mathbf{F}}^\top \mathbf{R}^{+\top} = (\mathbf{M}^\top \mathbf{W} \mathbf{M})^{-1} \mathbf{M}^\top \mathbf{W} \mathbf{N} \quad (25)$$

$$\hat{\mathbf{Q}} = \frac{1}{d-1} \left(\mathbf{N} - \mathbf{R}^+ \hat{\mathbf{F}} \mathbf{M} \right)^\top \mathbf{W} \left(\mathbf{N} - \mathbf{R}^+ \hat{\mathbf{F}} \mathbf{M} \right) \quad (26)$$

which are the typical vector auto-regression estimators with weights. Equation (25) and (26) can easily be computed in parallel using QR decompositions as in the `bam` function in the `mgcv` package with a low memory footprint (see Wood et al., 2014). This is currently implemented. Though, the gains from a parallel implementation may be small as the computation here have a computation time which is independent of the number of observations. Thus, the computation involved here is often fast as the dimension of the state vector is small.

5.2 Restricted vector auto-regression models

Suppose that we want to restrict some of the parameters of \mathbf{F} and \mathbf{Q} . Let

$$\begin{aligned}(s_1, s_2, \dots, s_r)^\top &= \mathbf{J}\boldsymbol{\psi} \\ (o_{21}, o_{31}, \dots, o_{r1}, o_{32}, \dots, o_{r2}, o_{43}, \dots, o_{r,r-1})^\top &= \mathbf{K}\boldsymbol{\phi}\end{aligned}$$

Then we can restrict the model such that

$$\begin{aligned}\sigma_i &= \exp(s_i) & \rho_{ij} &= \frac{2}{1 + \exp(-o_{ij})} - 1 \\ \text{vec}(\mathbf{R}^+ \mathbf{F}) &= \mathbf{G}\boldsymbol{\theta} & \mathbf{Q} &= \mathbf{V}\mathbf{C}\mathbf{V} \\ \mathbf{V} &= \begin{pmatrix} \sigma_1 & 0 & \cdots & 0 \\ 0 & \sigma_2 & \ddots & 0 \\ \vdots & \ddots & \ddots & \vdots \\ 0 & \cdots & 0 & \sigma_r \end{pmatrix} & \mathbf{C} &= \begin{pmatrix} 1 & \rho_{21} & \cdots & \rho_{r1} \\ \rho_{21} & 1 & \ddots & \rho_{r2} \\ \vdots & \ddots & \ddots & \vdots \\ \rho_{r1} & \cdots & \rho_{r,r-1} & 1 \end{pmatrix}\end{aligned}$$

and where $\text{vec}(\cdot)$ is the vectorization function which stacks the the columns of a matrix from left to right. E.g.,

$$\text{vec}(\mathbf{A}) = (a_{11}, a_{21}, a_{31}, a_{12}, a_{22}, a_{32}, a_{13}, a_{23}, a_{33})^\top$$

$$\mathbf{A} = \begin{pmatrix} a_{11} & a_{12} & a_{13} \\ a_{21} & a_{22} & a_{23} \\ a_{31} & a_{32} & a_{33} \end{pmatrix}$$

$\mathbf{G} \in \mathbb{R}^{rp \times g}$ is a known matrix with $g \leq rp$ and we assume that it has full column rank. Similarly, $\mathbf{J} \in \mathbb{R}^{r \times l}$ with $l \leq r$ and $\mathbf{K} \in \mathbb{R}^{r(r-1)/2 \times k}$ with $k \leq r(r-1)/2$. Both are known and have full column rank. We assume that \mathbf{G} is such that \mathbf{F} is non-singular for some $\boldsymbol{\theta}$ since Equation (9) used. Further, we assume that \mathbf{J} and \mathbf{K} are such that \mathbf{Q} is a positive definite matrix for some $\boldsymbol{\psi}$ and $\boldsymbol{\phi}$ pair. \mathbf{V} is a diagonal matrix containing the standard deviations and \mathbf{C} is the correlation matrix.

We cannot jointly maximize $\boldsymbol{\theta}$, $\boldsymbol{\psi}$, and $\boldsymbol{\phi}$ analytically but we can maximize $\boldsymbol{\theta}$ analytically conditional on $\boldsymbol{\psi}$ and $\boldsymbol{\phi}$. Hence, we can employ a Monte Carlo expectation conditional maximization algorithm in which we take two so-called conditional maximization steps (see Meng and Rubin, 1993, on the, non-Monte Carlo, expectation maximization algorithm). The first conditional maximization step is

$$\boldsymbol{\theta}^{(i+1)} = \mathbf{G}^+ \left(\mathbf{Q}^{(i)} \otimes (\mathbf{M}^\top \mathbf{W} \mathbf{M})^{-1} \right) \mathbf{G}^{+\top} \mathbf{G}^\top \text{vec} \left(\mathbf{M}^\top \mathbf{W} \mathbf{N} \mathbf{Q}^{-(i)} \right) \quad (27)$$

where \otimes is the Kronecker product and $\mathbf{Q}^{-(i)}$ is the inverse of $\mathbf{Q}^{(i)}$. Equation (27) is easily computed with the QR decomposition we make to compute for Equation (25). Having obtained the new $\boldsymbol{\theta}^{(i+1)}$, we update the variable elements of \mathbf{F} (those columns "picked out" by \mathbf{R} in $\mathbf{R}^+ \mathbf{F}$) and denote the new estimate $\widehat{\mathbf{F}}^{(i+1)}$. The second conditional maximization step which updates $\boldsymbol{\psi}$ and $\boldsymbol{\phi}$ is

$$\mathbf{Z} = \left(\mathbf{N} - \mathbf{R}^+ \widehat{\mathbf{F}}^{(i+1)} \mathbf{M} \right)^\top \mathbf{W} \left(\mathbf{N} - \mathbf{R}^+ \widehat{\mathbf{F}}^{(i+1)} \mathbf{M} \right)$$

$$\boldsymbol{\psi}^{(i+1)}, \boldsymbol{\phi}^{(i+1)} = \arg \max_{\boldsymbol{\psi}, \boldsymbol{\phi}} -(d-1) \log |\mathbf{Q}(\boldsymbol{\psi}, \boldsymbol{\phi})| - \text{tr} \left(\mathbf{Q}(\boldsymbol{\psi}, \boldsymbol{\phi})^{-1} \mathbf{Z} \right)$$

which can be done numerically. We have made \mathbf{Q} 's dependence on $\boldsymbol{\psi}$ and $\boldsymbol{\phi}$ explicit to emphasize which factors are affected. \mathbf{C} may not be a valid correlation matrix for all $\boldsymbol{\phi} \in \mathbb{R}^k$ for some choices of \mathbf{K} . The invalid values are ruled out excluded doing the numerical optimization. This completes the two conditional maximization steps. The next E-step is then performed using $\boldsymbol{\theta}^{(i+1)}$, $\boldsymbol{\psi}^{(i+1)}$, $\boldsymbol{\phi}^{(i+1)}$. Meng and Rubin (1993, see the discussion) comments that it may be beneficial to perform an E-step between each conditional maximization step when the E-step is relatively cheap. This is not the case here since all the above computations are independent of n_{\max} . Thus, if we have a moderately large number of observations at each time point relative to the dimension of the state vector, then we will use most of the computation time performing the E-step.

5.3 Estimating fixed effect coefficients

Next, we turn to estimating the fixed effects coefficients, $\boldsymbol{\omega}$, in Equation (1). If we assume that observations, y_{it} s, are from an exponential family then it is easy to show that the M-step estimator amounts to generalized linear model with N_s observations for each y_{it}

which differ only by an offset term and a weight. The offset term comes from the $\mathbf{x}_{it}^\top \mathbf{R} + \hat{\boldsymbol{\alpha}}_j^{(t)}$ term in Equation (1) for each of the $j = 1, \dots, N_s$ smoothed particles. The corresponding weights are the smoothed weights, $\hat{w}_j^{(t)}$. The problem can be solved in parallel using QR decompositions as in Section 5.1. This is what is done in the current implementation.

Currently, I only take one iteration of the iteratively re-weighted least squares. I gather I have to repeat till convergence though... This is however not nice computationally and the difference in the estimate from one M-step iteration to the next is very minor when you only take one iteratively re-weighted least square iteration...

5.4 Observed information matrix

Computing the observed information matrix requires an application of the missing information principle (Louis, 1982). However we cannot evaluate the quantities we need with the output from Algorithm 1 since we only have discrete approximation of the smoothed distribution of triplet of particles, $\boldsymbol{\alpha}_{t-1:t+1}$, but need an approximation for the entire path, $\boldsymbol{\alpha}_{1:d}$. One solution is to use the methods covered in e.g., Cappé et al. (2005, section 8.3 and chapter 11) and Poyiadjis et al. (2011). The method in Cappé et al. (2005) only requires the forward particle filter output. It is not implemented, though.

6 Other filter and smoother options

The $\mathcal{O}(N^2)$ two-filter smoother in Fearnhead et al. (2010) is going to be computationally expensive as an approximation is going to be needed for Equation (8) in their article. The non-auxiliary version in Briers et al. (2009) is more feasible as it only requires evaluation of f in the smoothing part of the generalized two-filter smoother (see Equation (46) in their paper). Similar conclusions applies to the forward smoother in Del Moral et al. (2010) and the backward smoother as presented in Kantas et al. (2015). Both have a $\mathcal{O}(N^2)$ computational cost.

Despite the $\mathcal{O}(N^2)$ cost of the method in Briers et al. (2009) and Del Moral et al. (2010) they are still worthy candidates as the computational cost is independent of the number of observations, n . Further, the computational cost can be reduced to $\mathcal{O}(N \log(N))$ run times with the approximations in Klaas et al. (2006).

The method in Malik and Pitt (2011, see particularly section 6.2 on page 203) can be used to do continuous likelihood evaluation. I am not sure how well this method scale with higher state dimensions, p .

Kantas et al. (2015) show empirically that it may be worth just using a forward filter. However, the example is with a univariate outcome ($n = 1$ – not to be confused with the number of periods d). The cost here of the forward filter is at least $\mathcal{O}(dNn_{\max}p)$. Every new particle yields an $\mathcal{O}(dn_{\max}p)$ cost which is expensive due to the large number of outcomes, n . Thus, the considerations are different and a $\mathcal{O}(dNn_{\max}p + N^2)$ method will not make a big difference unless N is large. Another alternative is to add noise to the parameters $\boldsymbol{\theta}$ at each time t and use the methods in Andrieu and Doucet (2002) or similar ideas to perform online estimation.

7 Briers et al. (2009)

The $\mathcal{O}(N^2)$ smoother from Briers et al. (2009) is also implemented as it is feasible for a moderate number of particles (though, we can use the approximations in Kantas et al., 2015 to reduce the computational complexity). It is shown in Algorithm 4. The weights in Equation (30) comes from the generalized two-filter formula. To motivate the smoother, we use

$$p(\mathbf{y}_{t:d}|\boldsymbol{\alpha}_t) = \tilde{p}(\mathbf{y}_{t:d}) \frac{\tilde{p}(\boldsymbol{\alpha}_t|\mathbf{y}_{t:d})}{\gamma_t(\boldsymbol{\alpha}_t)}$$

to generalize the two-filter formula in Kitagawa (1994) as follows

$$\begin{aligned} p(\boldsymbol{\alpha}_t|\mathbf{y}_{1:d}) &= \frac{p(\boldsymbol{\alpha}_t|\mathbf{y}_{1:t-1}) p(\mathbf{y}_{t:d}|\boldsymbol{\alpha}_t)}{p(\mathbf{y}_{t:d}|\mathbf{y}_{1:t-1})} \\ &\propto p(\boldsymbol{\alpha}_t|\mathbf{y}_{1:t-1}) p(\mathbf{y}_{t:d}|\boldsymbol{\alpha}_t) \\ &= p(\boldsymbol{\alpha}_t|\mathbf{y}_{1:t-1}) \tilde{p}(\mathbf{y}_{t:d}) \frac{\tilde{p}(\boldsymbol{\alpha}_t|\mathbf{y}_{t:d})}{\gamma_t(\boldsymbol{\alpha}_t)} \\ &\propto p(\boldsymbol{\alpha}_t|\mathbf{y}_{1:t-1}) \frac{\tilde{p}(\boldsymbol{\alpha}_t|\mathbf{y}_{t:d})}{\gamma_t(\boldsymbol{\alpha}_t)} \\ &= \tilde{p}(\boldsymbol{\alpha}_t|\mathbf{y}_{t:d}) \frac{[\int p(\boldsymbol{\alpha}_{t-1}|\mathbf{y}_{1:t-1}) f(\boldsymbol{\alpha}_t|\boldsymbol{\alpha}_{t-1}) \partial\boldsymbol{\alpha}_{t-1}]}{\gamma_t(\boldsymbol{\alpha}_t)} \\ &\approx \sum_{i=1}^N \tilde{w}_t^{(i)} \delta_{\tilde{\boldsymbol{\alpha}}_t^{(i)}}(\boldsymbol{\alpha}_t) \frac{[\sum_{j=1}^N w_{t-1}^{(j)} f(\tilde{\boldsymbol{\alpha}}_t^{(i)}|\boldsymbol{\alpha}_{t-1}^{(j)})]}{\gamma_t(\tilde{\boldsymbol{\alpha}}_t^{(i)})} \end{aligned} \tag{28}$$

where \approx means approximately proportional. Similar arguments leads to

$$\begin{aligned} p(\boldsymbol{\alpha}_{t-1:t}|\mathbf{y}_{1:d}) &\propto p(\boldsymbol{\alpha}_{t-1:t}|\mathbf{y}_{1:t-1}) p(\mathbf{y}_{t:d}|\boldsymbol{\alpha}_{t-1:t}) \\ &= f(\boldsymbol{\alpha}_t|\boldsymbol{\alpha}_{t-1}) p(\boldsymbol{\alpha}_{t-1}|\mathbf{y}_{1:t-1}) p(\mathbf{y}_{t:d}|\boldsymbol{\alpha}_t) \\ &\propto f(\boldsymbol{\alpha}_t|\boldsymbol{\alpha}_{t-1}) p(\boldsymbol{\alpha}_{t-1}|\mathbf{y}_{1:t-1}) \frac{\tilde{p}(\boldsymbol{\alpha}_t|\mathbf{y}_{t:d})}{\gamma_t(\boldsymbol{\alpha}_t)} \\ &\approx \sum_{i=1}^N \sum_{j=1}^N \tilde{w}_t^{(i)} \delta_{\tilde{\boldsymbol{\alpha}}_t^{(i)}}(\boldsymbol{\alpha}_t) \frac{[\sum_{k=1}^N w_{t-1}^{(k)} f(\tilde{\boldsymbol{\alpha}}_t^{(i)}|\boldsymbol{\alpha}_{t-1}^{(k)})]}{\gamma_t(\tilde{\boldsymbol{\alpha}}_t^{(i)})} \\ &\quad \cdot \frac{w_{t-1}^{(j)} \delta_{\boldsymbol{\alpha}_{t-1}^{(j)}}(\boldsymbol{\alpha}_{t-1}) f(\tilde{\boldsymbol{\alpha}}_t^{(i)}|\boldsymbol{\alpha}_{t-1}^{(j)})}{[\sum_{k=1}^N w_{t-1}^{(k)} f(\tilde{\boldsymbol{\alpha}}_t^{(i)}|\boldsymbol{\alpha}_{t-1}^{(k)})]} \\ &= \sum_{i=1}^N \sum_{j=1}^N \hat{w}_t^{(i,j)} \delta_{\tilde{\boldsymbol{\alpha}}_t^{(i)}}(\boldsymbol{\alpha}_t) \delta_{\boldsymbol{\alpha}_{t-1}^{(j)}}(\boldsymbol{\alpha}_{t-1}) \end{aligned}$$

where

$$\hat{w}_t^{(i,j)} = \tilde{w}_t^{(i)} \frac{w_{t-1}^{(j)} f(\tilde{\boldsymbol{\alpha}}_t^{(i)}|\boldsymbol{\alpha}_{t-1}^{(j)})}{[\sum_{j=1}^N w_{t-1}^{(j)} f(\tilde{\boldsymbol{\alpha}}_t^{(i)}|\boldsymbol{\alpha}_{t-1}^{(j)})]} \tag{29}$$

We need the latter for the EM-algorithm.

Algorithm 4 $\mathcal{O}(N^2)$ generalized two-filter smoother using the method in Briers et al. (2009).

Input:

- $\mathbf{Q}, \mathbf{Q}_0, \mathbf{a}_0, \mathbf{X}_1, \dots, \mathbf{X}_d, \mathbf{y}_1, \dots, \mathbf{y}_d, R_1, \dots, R_d, \boldsymbol{\omega}$
- 1: **procedure** FILTER FORWARD
 - 2: Run a forward particle filter to get particle clouds $\left\{ \boldsymbol{\alpha}_t^{(j)}, w_t^{(j)}, \beta_{t+1}^{(j)} \right\}_{j=1, \dots, N}$ approximating $p(\boldsymbol{\alpha}_t | \mathbf{y}_{1:t})$ for $t = 0, 1, \dots, d$. See Algorithm 2.
 - 3: **procedure** FILTER BACKWARDS
 - 4: Run a similar backward filter to get $\left\{ \tilde{\boldsymbol{\alpha}}_t^{(k)}, \tilde{w}_t^{(k)}, \tilde{\beta}_{t-1}^{(k)} \right\}_{k=1, \dots, N}$ approximating $\tilde{p}(\boldsymbol{\alpha}_t | \mathbf{y}_{t:d})$ for $t = d+1, d, d-1, \dots, 1$. See Algorithm 3.
 - 5: **procedure** SMOOTH (COMBINE)
 - 6: **for** $t = 1, \dots, d$ **do**
 - 7: Assign each backward filter particle a smoothing weight given by

$$\hat{w}_t^{(i)} \propto \tilde{w}_t^{(i)} \frac{\left[\sum_{j=1}^N w_{t-1}^{(j)} f\left(\tilde{\boldsymbol{\alpha}}_t^{(i)} | \boldsymbol{\alpha}_{t-1}^{(j)}\right) \right]}{\gamma_t\left(\tilde{\boldsymbol{\alpha}}_t^{(i)}\right)} \quad (30)$$

With the result above, we can show the arguments behind the smoother from Fearnhead et al. (2010). Similar to Equation (28) we find that

$$\begin{aligned} p(\boldsymbol{\alpha}_t | \mathbf{y}_{1:d}) &\propto p(\boldsymbol{\alpha}_t | \mathbf{y}_{1:t-1}) p(\mathbf{y}_{t:d} | \boldsymbol{\alpha}_t) \\ &= p(\boldsymbol{\alpha}_t | \mathbf{y}_{1:t-1}) g_t(\mathbf{y}_t | \boldsymbol{\alpha}_t) p(\mathbf{y}_{t+1:d} | \boldsymbol{\alpha}_t) \\ &= \int f(\boldsymbol{\alpha}_t | \boldsymbol{\alpha}_{t-1}) p(\boldsymbol{\alpha}_{t-1} | \mathbf{y}_{1:t-1}) d\boldsymbol{\alpha}_{t-1} g_t(\mathbf{y}_t | \boldsymbol{\alpha}_t) \\ &\quad \cdot \int f(\boldsymbol{\alpha}_{t+1} | \boldsymbol{\alpha}_t) p(\mathbf{y}_{t+1:d} | \boldsymbol{\alpha}_{t+1}) d\boldsymbol{\alpha}_{t+1} \\ &\propto \int f(\boldsymbol{\alpha}_t | \boldsymbol{\alpha}_{t-1}) p(\boldsymbol{\alpha}_{t-1} | \mathbf{y}_{1:t-1}) d\boldsymbol{\alpha}_{t-1} g_t(\mathbf{y}_t | \boldsymbol{\alpha}_t) \\ &\quad \cdot \int f(\boldsymbol{\alpha}_{t+1} | \boldsymbol{\alpha}_t) \frac{\tilde{p}(\boldsymbol{\alpha}_{t+1} | \mathbf{y}_{t+1:d})}{\gamma_{t+1}(\boldsymbol{\alpha}_{t+1})} d\boldsymbol{\alpha}_{t+1} \\ &\propto \sum_{i=1}^N \sum_{j=1}^N f\left(\boldsymbol{\alpha}_t | \boldsymbol{\alpha}_{t-1}^{(j)}\right) w_{t-1}^{(j)} g_t(\mathbf{y}_t | \boldsymbol{\alpha}_t) f\left(\tilde{\boldsymbol{\alpha}}_{t+1}^{(i)} | \boldsymbol{\alpha}_t\right) \frac{\tilde{w}_{t+1}^{(i)}}{\gamma_{t+1}(\tilde{\boldsymbol{\alpha}}_{t+1}^{(i)})} \end{aligned}$$

Thus, we can sample $\boldsymbol{\alpha}_t$ from a proposal distribution given the time $t-1$ forward filter particle, $\boldsymbol{\alpha}_{t-1}^{(j)}$, and time $t+1$ backward filter particle, $\tilde{\boldsymbol{\alpha}}_{t+1}^{(i)}$, for all N^2 particle pairs. Alternatively, we can sample the $t-1$ and $t+1$ particles independently which yields Algorithm 1.

Further, we can find that

$$\begin{aligned}
p(\boldsymbol{\alpha}_{t-1:t}|\mathbf{y}_{1:d}) &= p(\boldsymbol{\alpha}_{t-1:t}|\mathbf{y}_{1:t-1}) g_t(\mathbf{y}_t|\boldsymbol{\alpha}_{t-1:t}) p(\mathbf{y}_{t+1:d}|\boldsymbol{\alpha}_{t-1:t}) \\
&\propto f(\boldsymbol{\alpha}_t|\boldsymbol{\alpha}_{t-1}) p(\boldsymbol{\alpha}_{t-1}|\mathbf{y}_{1:t-1}) g_t(\mathbf{y}_t|\boldsymbol{\alpha}_t) \\
&\quad \cdot \int f(\boldsymbol{\alpha}_{t+1}|\boldsymbol{\alpha}_t) \frac{\tilde{p}(\boldsymbol{\alpha}_{t+1}|\mathbf{y}_{t+1:d})}{\gamma_{t+1}(\boldsymbol{\alpha}_{t+1})} d\boldsymbol{\alpha}_{t+1} \\
&\approx \sum_{i=1}^{N_s} \delta_{\hat{\boldsymbol{\alpha}}_t^{(i)}}(\boldsymbol{\alpha}_t) \delta_{\boldsymbol{\alpha}_{t-1}^{(j_i)}}(\boldsymbol{\alpha}_t) f(\hat{\boldsymbol{\alpha}}_t^{(i)}|\boldsymbol{\alpha}_{t-1}^{(j_i)}) w_{t-1}^{(j_i)} g_t(\mathbf{y}_t|\hat{\boldsymbol{\alpha}}_t^{(i)}) \\
&\quad \cdot \int f(\boldsymbol{\alpha}_{t+1}|\hat{\boldsymbol{\alpha}}_t^{(i)}) \frac{\tilde{p}(\boldsymbol{\alpha}_{t+1}|\mathbf{y}_{t+1:d})}{\gamma_{t+1}(\boldsymbol{\alpha}_{t+1})} d\boldsymbol{\alpha}_{t+1} \\
&\approx \sum_{i=1}^{N_s} \hat{w}_t^{(i)} \delta_{\hat{\boldsymbol{\alpha}}_t^{(i)}}(\boldsymbol{\alpha}_t) \delta_{\boldsymbol{\alpha}_{t-1}^{(j_i)}}(\boldsymbol{\alpha}_t)
\end{aligned}$$

where superscripts j_i are used as in Algorithm 1 implicitly dependent on t .

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A Vector and matrix products

The methods described above involves many vector and matrix products. These are implemented with C++ abstract classes with specialized `map` member functions for particular problems to decrease the computation. An alternative would have been to use a sparse matrix implementation. It does not matter much in terms of computation time if n_t is much larger then the state vector's dimension. However, this section for people who looks through the code. As an example we have a mapping matrix \mathbf{A} which C++ abstract member on the main data object used in the package called `xyz`. E.g., this could be \mathbf{F} with name `state_trans`. Then the following operations are implemented

- `map()`: Returns \mathbf{A} .
- `map(const arma::vec &x, bool tranpose)`: Returns $\mathbf{A}^\top \mathbf{x}$ if `tranpose == true` and $\mathbf{A}\mathbf{x}$ otherwise.
- `map(const arma::mat &X, side s, bool tranpose)`: Let $\mathbf{B} = \mathbf{A}^\top$ if `tranpose == true` and otherwise $\mathbf{B} = \mathbf{A}$. Then the result is $\mathbf{B}\mathbf{X}\mathbf{B}^\top$ if `s == both`, $\mathbf{B}\mathbf{X}$ if `s == left`, and $\mathbf{X}\mathbf{B}^\top$ if `s == right`.

These are implemented for

C++ member name	Matrix \mathbf{A}
<code>err_state</code>	\mathbf{R}
<code>err_state_inv</code>	$\mathbf{R}^+ = \mathbf{R}^\top$
<code>state_trans</code>	\mathbf{F}
<code>state_trans_err</code>	$\mathbf{R}^+\mathbf{F} = \mathbf{R}^\top\mathbf{F}$
<code>state_trans_inv</code>	\mathbf{F}^{-1}

Further, we will need function to compute terms

$$\hat{\mathbf{P}}_t \mathbf{F}^\top \hat{\mathbf{P}}_{t+1}^{-1} \boldsymbol{\alpha}_{t+1} + \hat{\mathbf{S}}_t \hat{\mathbf{P}}_t^{-1} \hat{\mathbf{m}}_t$$

for an arbitrary $\boldsymbol{\alpha}_{t+1}$, $\hat{\mathbf{S}}_t = \hat{\mathbf{P}}_t \mathbf{F}^\top \hat{\mathbf{P}}_{t+1}^{-1} \mathbf{R} \mathbf{Q} \mathbf{R}^\top \mathbf{F}^{-\top}$, $\hat{\mathbf{P}}_t^{-1} \hat{\mathbf{m}}_t$, and $\hat{\mathbf{P}}_t^{-1}$. This is done with the methods `bw_mean(signed int, const arma::vec&)`, `bw_covar(signed int)`, `uncond_mean_term(signed int)`, and `uncond_covar_inv(signed int)`. The terms and factors that can be computed once are computed once.