

## EXERCISE 2

- (1) Simulate and plot 10 timesteps of the following model,

$$y_t = 0.1y_{t-1} + 0.9y_{t-2} + 0.5w_t, \quad y_1 = 0, \quad w_t \sim \mathcal{N}(0, 1) \text{ are iid.}$$

- (2) Write the model above in state space form. That is,

$$\begin{aligned} Y_{t+1} &= AY_t + HV_t, \\ y_t &= BY_t \end{aligned}$$

for some  $Y_t \in \mathbb{R}^3$ ,  $A \in \mathbb{R}^{3 \times 3}$ ,  $H \in \mathbb{R}^{3 \times 1}$ ,  $B \in \mathbb{R}^{1 \times 3}$  and  $V_t \in \mathbb{R}$ .

- (3) Implement the Kalman filter for this model and compute the conditional distribution of  $y_t$  given  $y_1, \dots, y_{10}$  for  $t = 11, 12, \dots, 20$ . (That is, the density of a multivariate Gaussian distribution with mean,

$$E[y_t | y_1, \dots, y_{10}],$$

and variance,

$$E[(y_t - E[y_t | y_1, \dots, y_{10}])^2 | y_1, \dots, y_{10}].$$

- (4) Plot,

$$E[y_t | y_1, \dots, y_{10}],$$

for  $t = 11, \dots, 20$ , with approximate 95%-confidence intervals. How are the confidence intervals for  $t$  close to 10? What happens when  $t$  becomes large?

*Hint.* Approximate 95% confidence intervals can be computed as,

$$E[y_t | y_1, \dots, y_{10}] \pm 2E[(y_t - E[y_t | y_1, \dots, y_{10}])^2 | y_1, \dots, y_{10}].$$

- (5) Assume now that the observations are given as,

$$z_t = y_t + v_t,$$

where  $v_t \sim \mathcal{N}(0, 1)$ . Plot  $z_t$  for  $t = 1, \dots, 10$ . Compute the conditional distribution of  $y_t$  given  $z_1, \dots, z_{10}$ , for  $t = 11, \dots, 20$ , and plot it with the observations. Include 95% confidence intervals.

- (6) Determine whether the following difference equation has a stationary and/or causal solution,

$$(1 - 0.6B + 0.05B^2)X_t = W_t.$$

- (7) Follow the instructions on the next page, and try to estimate the parameters of the following  $AR(1)$  model with observations noise:

$$\begin{aligned} X_{t+1} &= \begin{pmatrix} \alpha & 0 \\ 1 & 0 \end{pmatrix} X_t + \sigma_1 \begin{pmatrix} 1 \\ 0 \end{pmatrix} v_t \\ Y_t &= (1, 0)X_t + \sigma_2 w_t, \end{aligned}$$

where  $v_t, w_t \sim \mathcal{N}(0, 1)$  and are iid.

Do this as follows: simulate  $Y_t$  for  $t = 1, \dots, 20$ , where  $\alpha, \sigma_1, \sigma_2 = 0.5$ . Implement the likelihood computation, and use the true parameter values as the initial guess (this is the "par" argument in `optim`).

STATE SPACE MODELS, KALMAN FILTER, AND COMPUTING THE LIKELIHOOD OF  
 $Y_1, \dots, Y_T$

A Gaussian and linear state space model is given as follows,

$$\begin{aligned} X_{t+1} &= AX_t + HV_t, \\ Y_t &= BX_t + DW_t, \end{aligned}$$

where  $X_t, X_1 \in \mathbb{R}^N$  and  $Y_t \in \mathbb{R}^M$  has are zero mean multivariate Gaussian, and,  $H \in \mathbb{R}^{N \times K}$ ,  $D \in \mathbb{R}^{M \times L}$ , for integers  $N, M, K, L$ . Furthermore,  $W_t \sim \mathcal{N}(0, I)$  is  $L$ -dimensional multivariate Gaussian, and  $V_t \sim \mathcal{N}(0, I)$  is  $K$ -dimensional multivariate Gaussian, with  $W_t, V_t$  independent for each  $t$ .

We use the short notation,  $X_{k|m} \in \mathbb{R}^N$  as the best linear predictor of  $X_k$  using  $Y_1, \dots, Y_m$ . In the Gaussian case, this is just  $E[X_k | Y_1, \dots, Y_m]$ .

Recall that the best linear predictor of  $X$  in terms of  $Y$  can be found using the best-linear-predictor-condition: for some  $M \in \mathbb{R}^{N \times M}$ , we must have,

$$X = MY \iff E[(X - MY)Y^T] = 0,$$

where the expectation acts compenentwise on the the  $R^{N \times M}$  entries of  $(X - MY)Y^T$ . We need to solve this equation for some matrix  $M$ . And it turns out that  $M$  is given by,

$$M = E[XY^T](E[YY^T])^{-1},$$

(where  $(E[YY^T])^{-1}$  is the psuedo inverse of  $E[YY^T]$ ).

We define the  $k$ 'th innovation,

$$I_k := Y_k - Y_{k|k-1},$$

and we also introduce use the simplified notation,

$$S_{k|m} = E[(X_k - X_{k|m})(X_k - X_{k|m})^T | Y_1, \dots, Y_m].$$

Then we can compute iteratively  $I_t$  and  $C_t := E[I_t I_t^T]$  using the Kalman filter.

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**Algorithm 1** Kalman filter

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```
(Initial condition,  $X \sim \mathcal{N}(X_1, S_1)$ )
 $X_{1|0} \leftarrow X_1$ 
 $S_{1|0} \leftarrow S_1$ 
for  $t = 1, \dots, T$  do
  Compute innovation,  $I_t$ 
   $I_t = Y_t - BX_{t|t-1}$ 
  Compute projection matrix  $M$ 
   $M = S_{t|t-1}B^T C_t^{-1}$ ,  $C_t = BS_{t|t-1}B^T + DD^T$ 
  Condition on innovation/observation
   $X_{t|t} \leftarrow X_{t|t-1} + MI_t$ 
  Update  $X_{t|t}, S_{t|t-1}$ 
   $X_{t+1|t} \leftarrow AX_{t|t}$ 
   $S_{t+1|t} \leftarrow A(I - MB)S_{t|t-1}A^T + HH^T$ 
end for
```

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We can use the computed innovations and their variance to compute the likelihood in this case. Since the innovations are Gaussian and uncorrelated, they are independent, and we have,

$$L(\theta) = C \prod_{t=1}^T \det(C_t)^{-1/2} e^{-\frac{1}{2} I_t^T C_t^{-1} I_t},$$

for some constant  $C$  independent of  $\theta$ .

In practice, we need to optimize the logarithm of the likelihood, since this is computationally easier.

In R, you can minimize any function, "obj", that you have defined, by using optim as follows,

```
opt <- optim(par=initial_param, fn=obj, method="BFGS",
             control=list(maxit=300, reltol=1e-8)).
```

use ?optim in the console in Rstudio to get more info on the arguments of optim, and what is returned.

*Hint.* A skeleton of the likelihood implementation in R is,

```
n <- 3

# objective function to be minimized
obj <- function(param) {
  # param is a vector of length 3 with the parameters in the model
  alpha <- param[1]
  sigma1 <- param[2]
  sigma2 <- param[3]

  # define the matrix A
  A <- matrix(0,nrow=n,ncol=n)
  A[1,1] = ...
  A[2,1] = ...

  # observation matrix
  B <- matrix(0,nrow=1,ncol=n)
  B[1,1] = ...

  # system noise
  HH <- ...

  # observation noise
  DD <- ...

  # compute likelihood

  # initial condition
  x <- matrix(0,nrow=n,ncol=1)
  S <- diag(n) # just use something here :)

  # log likelihood
  ll <- 0.
  for (t in 1:tt) {
    # compute innovation

    # innovation variance

    # compute M matrix
```

```
      # update x, S

      # add to ll
      ll <- ll + ...

    }

    return(ll)
  }
```