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**Fast parameter and confidence interval estimation for Hidden Markov Models using Template Model Builder**

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Hidden Markov Models (HMMs) are a class of models widely used in speech recognition (See e.g. [Gales](#_bookmark26) [and Young](#_bookmark26), [2008](#_bookmark26); [Fredkin and Rice](#_bookmark24), [1992](#_bookmark24)). There are straightforward ways to compute Maximum Likeli- hood Estimates (MLE) of their parameters, but getting confidence intervals can be more difficult (See e.g. [Zucchini et al.](#_bookmark38), [2016](#_bookmark38); [Lystig and Hughes](#_bookmark32), [2002](#_bookmark32)). In addition, computing MLEs can be time-consuming when the dataset and the complexity become very large. We show in this paper a way to speed up compu- tation by up to a few hundred times in the language R when compared to usual optimization approaches, and at the same time retrieve standard errors easily. Here, we show in a first part how to optimize HMMs with the TMB package in R and how to retrieve confidence intervals. In a second part, we compare different optimizers such as nlminb, and minimize the negative log-likelihood directly on different datasets: a small one (240 data points) from [Leroux and Puterman](#_bookmark31) ([1992](#_bookmark31)), a medium sized simulated one (2000 data points), and a large one (87648) from a hospital.

*Key words:* Hidden Markov Model; TMB; Confidence intervals

*(Up to five keywords are allowed and should be given in alphabetical order. Please capitalize the key words)*

Supporting Information for this article is available from the author or on the WWW under<http://dx.doi.org/10.1022/bimj.XXXXXXX> (please delete if not applicable)

# 1 Introduction

Hidden Markov models (HMM) are a versatile type of model that have used in many different situations since their introduction by [Baum and Petrie](#_bookmark21) ([1966](#_bookmark21)). As an example, [Fredkin and Rice](#_bookmark24) ([1992](#_bookmark24)) applied them in speech recognition, [Lystig and Hughes](#_bookmark32) ([2002](#_bookmark32)) to rainfall occurence data, and [Schadt et al.](#_bookmark35) ([1998](#_bookmark35)) to phylogenetetic trees. [Fredkin and Rice](#_bookmark24) ([1992](#_bookmark24)) and [Zucchini et al.](#_bookmark38) ([2016](#_bookmark38)) are references in the theory of HMMs. Evaluating uncertainty and getting confidence intervals in HMMs is not easy. Although [Cappe´](#_bookmark22) [et al.](#_bookmark22) ([2006](#_bookmark22), Chapter 12) showed it can be achieved using asymptotic normality of the MLEs of the param- eters under certain conditions, [Fru¨hwirth-Schnatter](#_bookmark25) ([2006](#_bookmark25)) points out that in indepedent mixture models, “The regularity conditions are often violated”, and [McLachlan and Peel](#_bookmark33) ([2004](#_bookmark33), p. 68) adds that “In partic- ular for mixture models, it is well known that the sample size *n* has to be very large before the asymptotic theory of maximum likelihood applies.” [Lystig and Hughes](#_bookmark32) ([2002](#_bookmark32)) shows a way to compute the exact Hessian, and [Zucchini et al.](#_bookmark38) ([2016](#_bookmark38)) shows another way to compute the approximate Hessian and thus con- fidence intervals but admits that “the use of the Hessian to compute standard errors (and thence confidence intervals) is unreliable if some of the parameters are on or near the boundary of their parameter space”.

TMB (Template Model Builder) is an R package for efficient fitting of complex statistical random effect models to data, as described by [Kristensen et al.](#_bookmark30) ([2015](#_bookmark30)). It provides exact calculations of first and second

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2 First Author*et al.* and dd: Running title

order derivatives of the likelihood of a model by automatic differentiation, which allows for efficient gra- dient and/or Hessian based optimization of the likelihood as well as uncertainty estimates by means of the Hessian. The Hessian is not necessarily directly applicable for evaluating parameter uncertainty in HMMs as there are several parameter constraints in these models. This can be adressed by constraint optimiza- tion, and subsequently combining the Hessian with the Jacobian of the constraints to obtain the covariance matrix as shown by [Visser et al.](#_bookmark37) ([2000](#_bookmark37)). Alternatively, [Zucchini et al.](#_bookmark38) ([2016](#_bookmark38)) shows how the constraints can be imposed by suitable transformations of the parameters. The covariance matrix of the untransformed (original) parameters can then be obtained by the delta method, a feature implemented in TMB.

In this paper, we will show how to optimize a Poisson HMM using TMB in the language R. Afterwards, we will use an example to show how to make a nested model, easily compute confidence intervals, retrieve interesting probabilities, we will apply a Poisson HMM on a hospital dataset. Eventually, we will see that TMB can accelerate traditional optimizers in R by up to a few hundred times on a large dataset, simplifies evaluating confidence intervals, and gives similar parameter uncertainty to bootstrap and profile methods via its Hessian based approach.

We decided to use a direct maximization approach instead of the Expectation-Maximization (EM) algo- rithm. The reason is that the direct maximization approach is easier to adapt if one wants to fit different and more complex models. It also deals easily with missing observations whereas the EM approach is more complex.

# Principles of using TMB for Maximum Likelihood Estimation (MLE)

## Setup

In order to use TMB in R, it is necessary to use the R-package TMB and have the software Rtools

installed. The latter is used for compiling the C++ code.

In this paper, we will use R version 3.6.0 and the code was run on Windows 10 Enterprise version 1809, with an Intel(R) Core(TM) i7-8700 processor.

The C++ code can be difficult to debug as it operates using a specific template, see [Section 2.2](#_bookmark1) below for an example. TMB provides a debugging feature available with RStudio which can be useful to retrieve diagnostic error messages. Enabling it is optional. The command TMB:::setupRStudio() enables this feature and requires manual input to confirm.

## Linear regression example

We start by illustrating how TMB can be used to fit a simple linear model and how to handle parameters subject to constraints, which is particularly relevant for HMM models. A more comprehensive tutorial on how to use TMB can be found at https://kaskr.github.io/adcomp/ book/Tutorial.html.

Let ***x*** and ***y*** be datasets of size n, then the negative log-likelihood corresponding to a simple linear model is given by

*n*

−*l*(*a, b, σ*2) = − ) log(*φ*(*yi*; *a* + *bxi, σ*2)))*,* (1)

*i*=1

where *φ*(·; *µ, σ*2) is the density function of the univariate normal distribution with mean *µ* and variance *σ*2.

TMB requires the (negative) log-likelihood function to be written in C++ under a specific template, and then loaded into R. Minimization and other post-processing procedures are all carried out in R.

Therefore, we require 2 files. The first file named *linreg.cpp* is written in C++,*.* and defines the negative log-likelihood (nll) function of the linear model as follows

Biometrical Journal **52** (2020) 61 3

**# i n c l u d e** *<*TMB. hpp*> / / import t h e TMB t e m p l a t e*

**template** *<***c l a s s** Type*>*

Type o b j e c t i v e f u n c t i o n *<*Type *>* : : **operator** ( ) ( )

*{*

DATA VECTOR( y ) ; */ / Data v e c t o r y t r a n s m i t t e d from R*

DATA VECTOR( x ) ; */ / Data v e c t o r x t r a n s m i t t e d from R*

PARAMETER( a ) ; */ / Parameter a t r a n s m i t t e d from R* PARAMETER( b ) ; */ / Parameter b t r a n s m i t t e d from R* PARAMETER( log Sigma ) ; */ / Sigma on log−s c a l e t r a n s m i t t e d from R*

*/ / Transform log Sigma back t o n a t u r a l s c a l e*

Type sigma = exp ( log Sigma ) ;

*/ / Declare n e g a t i v e log−l i k e l i h o o d*

Type n l l = *−* sum ( dnorm ( y ,

a + b \* x , sigma ,

**t ru e** ) ) ;

*/ / This l e t s t h e user r e t r i e v e any v a r i a b l e s along with t h e i r s t andard e r r o r s*

ADREPORT( a ) ; ADREPORT( b ) ;

ADREPORT( sigma ) ;

**return** n l l ;

*}*

Note that we define data inputs  *x* and *y*  using the DATA\_VECTOR() declaration in the above code. Furthermore, we declare the nll as a function of the three parameters using the PARAMETER() declaration.

Since the parameter *σ* has to be non-negative, the nll is parameterized in terms of log(*σ*) which can be freely estimated. This approach avoids the need of using constraint optimization methods. By subsequently calling the function ADREPORT on the variables, TMB can apply the generalized delta method described by [Kass and Steffey](#_bookmark29) ([1989](#_bookmark29)) to cal- culate the corresponding standard errors. This feature is particularly useful when the likelihood has been reparameterized as above, and is especially relevant for HMMs.

The second file needed is written in R and serves for compiling the nll function defined above, carrying out the estimation procedure by numerical optimization of the nll function and various post-processing of the results.

Once the C++ file with the negative log-likelihood is established it can be compiled and the objective function can be minimized as shown in the .R file below:

*# Setting up TMB*

**library**(TMB) TMB::**compile**("code/linreg.cpp")

## [1] 0

**dyn.load**(**dynlib**("code/linreg"))

*# Generate the data*

**set.seed**(123)

4 First Author*et al.* and dd: Running title

data <- **list**(y = **rnorm**(20) + 1:20, x = 1:20) parameters <- **list**(a = 0, b = 0, logSigma = 0)

*# Tell TMB to create the likelihood function*

obj\_linreg <- **MakeADFun**(data, parameters, DLL = "linreg", silent = TRUE)

*# Optimization of the objective function with TMB* fit\_linreg <- **nlminb**(obj\_linreg$par, obj\_linreg$fn) fit\_linreg$par

##

a

b

logSigma

## 0.31009240270 0.98395534754 -0.05814659321

The function sdreport will by default give the estimates and standard errors of the variables that the negative log-likelihood is parameterized in terms of:

**sdreport**(obj\_linreg)

##

##

##

##

##

##

sdreport(.) result

Estimate

a b

logSigma

0.31009240270

0.98395534754

-0.05814659321

Std. Error 0.43829082894

0.03658781149

0.15811381455

Maximum gradient component: 6.931683128e-05

To get the estimates and standard errors of the variables called with ADREPORT() in the C++ file, we can run a summary call with the select = "report" option:

|  |  |  |  |
| --- | --- | --- | --- |
| ## |  | Estimate | Std. Error |
| ## | a | 0.3100924027 | 0.43829082894 |
| ## | b | 0.9839553475 | 0.03658781149 |
| ## | sigma | 0.9435116249 | 0.14918222209 |

And we can see the same estimates from a regular linear regression.

**summary**(**sdreport**(obj\_linreg), select = "report")

**lm**(y ˜ x, data = data)$coefficients

## (Intercept)

x

## 0.3100925100 0.9839553612

# Estimation TMB/HMM

## Notation

A Hidden Markov Model (HMM) is a model where the data is assumed to follow a mixture of distributions, going from one to the other as time progresses according to probabilities set by an underlying homogeneous stationary Markov chain. In this paper we will focus on a Poisson-HMM, but only small changes to the code are needed to obtain models with other conditional distributions than the Poisson. The data denoted

as {*Xn* : *n* = 1*, . . . , N* } is assumed to follow a mixture of *m* Poisson distributions with parameters {*λi* :

*i* = 1*, . . . , m*}.

Biometrical Journal **52** (2020) 61 5

Let {*Ct* : *t* = 1*, . . . , N* } be the underlying Markov chain. The Markov chain is assumed irreducible and aperiodic. [Grimmett et al.](#_bookmark27) ([2001](#_bookmark27), Lemma 6.3.5 on p. 225 and Theorem 6.4.3 on p. 227) has shown that irreducibility ensures the existence of the stationary distribution. As has been shown by [Feller](#_bookmark23) ([1968](#_bookmark23),

p. 394), aperiodicity implies that a unique limiting distribution exists and is the stationary distribution.

*p*1(*x*) 0 

Let *pi*(*x*) = *P* (*Xt* = *x*|*Ct* = *i*)*,* ∀*i* = 1*, . . . , m* and let **P**(*x*) = 







*p*2(*x*)



. . .



.



*p*1(*x*1) *p*2(*x*1) *. . . pm*(*x*1)

We will call *p*1(*x*2) *p*2(*x*2) *. . . pm*(*x*2)

0 *pm*(*x*)

 ..





.



...

. . .

...

 the emission probability matrix. If *x* is a missing data,



*p*1(*xn*) *p*2(*xn*) *. . . pm*(*xn*)

we will define *pi*(*x*) = 1 and therefore the emission matrix of the dataset will contain a line filled with ones.

Furthermore, we let **Γ** = {*γij* } denote the transition matrix of the Markov chain, and ***δ*** be its stationary distribution.

Finally, we let *X*(*t*) = {*X*1*, . . . , Xt*} and *x*(*t*) = {*x*1*, . . . , xt*} denote the history up to time *t*.

In this paper we focus on Poisson HMMs, and so *pi*(*x*) = *e*

This can however be easily changed to other distributions.

*−λi x i*

*λ*

.

*x*!

## Hmm Likelihood

It can be shown [(Zucchini et al.](#_bookmark38), [2016](#_bookmark38), p. 36) that the likelihood of an HMM is given by

*LN* = **P**(*X*(*t*) = *x*(*t*)) = ***δ*P**(*x*1)**ΓP**(*x*2)**ΓP**(*x*3) *. . .* **ΓP**(*xN* )**1**!

It is usual for HMMs to estimate the parameters ***θ*** (where ***θ*** = (*γ*11*, . . . , γ*1*m, . . . , γm*1*, . . . , γmm, λ*1*, . . . , λm*)) subject to estimation in the case of a Poisson HMM) by optimizing a reparametrized version of the log-likelihood log *L*(***ψ***), where ***ψ*** = *g*−1(***θ***) represent a set of unconstrained parameters. That transfor- mation will be detailed below. As a result of the invariance principle, the maximum likelihood estimate of

***θ*** is given by ***θ***ˆ = *g*(***ψ***ˆ ) where ***ψ***ˆ is the minimizer of − log *L*(***ψ***). To avoid constraints on the parameters, we transform the natural parameters {**Γ***,* ***λ***} into working parameters {**T***,* ***η***}.

In practice, as explained in [Section 2](#_bookmark0), we first create a set of natural parameters and turn them into

working parameters before feeding those in an optimizer. One possible transformation of **Γ** is done this way:

exp(*τij* )

*γij* = 1 +

*k*j=*i τik*

*,* for *i* /= *j*

Each row must add to 1, so we can get the diagonal elements of **Γ** easily.

*ηi* = log(*λi*).

The reverse transformation is

And *λi* = exp(*ηi*).

*τij* = log

/ *γij* \

1 − *k*j=*i γik*

= log(*γij/γii*)*,*

for *i* /= *j*

The relevant R function to do so is given by [Zucchini et al.](#_bookmark38) ([2016](#_bookmark38), p. 52):

6 First Author*et al.* and dd: Running title

*# Transform Poisson natural parameters to working parameters*

pois.HMM.pn2pw <- **function**(m, lambda, gamma, delta = **NULL**,

stationary = TRUE) {

tlambda <- **log**(lambda)

**if** (m == 1) **return**(tlambda)

foo <- **log**(gamma / **diag**(gamma)) tgamma <- **as.vector**(foo[!**diag**(m)])

**if** (stationary) {

*# If tdelta is set to NULL and returned in the list,*

*# it will cause issues when optimizing with TMB*

**return**(**list**(tlambda = tlambda, tgamma = tgamma))

} **else** {

tdelta <- **log**(delta[- 1] / delta[1])

**return**(**list**(tlambda = tlambda, tgamma = tgamma, tdelta = tdelta))

}

}

These can be adapted to other distributions and transformations. Those working parameters are fed into TMB, which will then turn them back to natural parameters before calculating the negative log-likelihood. That transformation from working parameters to natural parameters is done in the C++ likelihood function.

The C++ function turning the working transition probability matrix into a natural parameter is the func- tion Gamma w2n available in the file utils.cpp. See Appendix A.2 [item (iii)](#_bookmark19).

If necessary, we can define the C++ function turning the working stationary distribution vector into a natural parameter with Delta w2n available in the file utils.cpp. See Appendix A.2 [item (ii)](#_bookmark18). In practice, we didn’t use it.

Given a working parameter vector log lambda, the C++ function turning the working vector of the Poisson means into a natural parameter is much simpler:

v e c t o r *<*Type*>* lambda = t l ambda . exp ( ) ;

The “forward algorithm” allows for a recursive computation of the likelihood.

To state the forward algorithm let us define the vector ***α****t* for *t* = 1*,* 2*, . . . , N* so that

***α****t* = ***δ*P**(*x*1)**ΓP**(*x*2)**ΓP**(*x*3) *. . .* **ΓP**(*xt*)

*t*

= ***δ*P**(*x*1) n **ΓP**(*xs*)

*s*=2

= (*αt*(1)*, . . . , αt*(*m*))

where ***δ*** denotes the initial distribution of the Markov chain. By convention, the empty product is the identity matrix. It is called the forward algorithm because of the way the ***α****t* values are calculated

***α***0 = ***δ***

***α****t* = ***α****t*−1**ΓP**(*xt*) for *t* = 1*,* 2*, . . . , N LT* = ***α****T* **1**!

Biometrical Journal **52** (2020) 61 7

In the same way, the “backward algorithm” allows for a recursive computation of the likelihood.

To state the backward algorithm let us define the vector ***β****t* for *t* = 1*,* 2*, . . . , N* so that

***β***! !

*t* = **ΓP**(*xt*+1)**ΓP**(*xt*+2) *. . .* **ΓP**(*xT* ) *. . .* **1**

/ *N*

= n

*s*=*t*+1

\

**ΓP**(*xs*) **1**!

= (*βt*(1)*, . . . , βt*(*m*))

It is called the backward algorithm because of the way the ***β****t* values are calculated

***β****T* = **1**!

***β****t* = **ΓP**(*xt*+1)***β****t*+1 for *t* = *N* − 1*, N* − 2*, . . . ,* 1

*LT* = ***δβ***1

Computing those probabilities directly can lead to underflow errors, it is therefore better to maximize the log-likelihood instead. We chose to minimize the negative log-likelihood for convenience purposes. In addition, in our case **Γ** and ***λ*** are subject to some constraints:

1. The means *λi* of the state-dependent distributions must be non-negative, for *i* = 1*, . . . , m*
2. The parameters *γi,j* of the transition probability matrix **Γ** must be non-negative, and the rows of **Γ**

must add to one

It should be noted that we use a scaled version of the forward algorithm in order to calculate the likeli- hood, as suggested by [Zucchini et al.](#_bookmark38) ([2016](#_bookmark38), p. 48).

In summary, we need to define the state-dependent probabilities (we used Poisson distributions, but it can easily be changed) and the transformations in R and C++ (they depend on the distributions used, but can be easily adapted). We show below the next step: writing the likelihood function in C++ and using it to model a dataset.

## Using TMB

* + 1. **Likelihood function**

As we did with the linear regression example, we first need to define our objective function: the negative log-likelihood, in a C++ file. Let’s name it poi hmm.cpp. Its code can be found in the Appendix A.2 [item (i)](#_bookmark17).

## Optimization

Once the objective function is written in C++, we can tell R to optimize its parameters.

1. Loading packages

*# Load TMB and optimization packages*

**library**(TMB) **library**(optimr)

*# Run the C++ file containing the TMB code*

TMB::**compile**("code/poi\_hmm.cpp", "-O1 -g", DLLFLAGS = "")

8 First Author*et al.* and dd: Running title

## [1] 0

*# Load it*

**dyn.load**(**dynlib**("code/poi\_hmm"))

*# Load the parameter transformation function*

**source**("functions/utils\_no\_tmb.R")

1. Loading the dataset, see [Section 6.1](#_bookmark8) for more details.

**load**("data/fetal-lamb.RData") lamb\_data <- lamb

1. Creating natural parameters as initial values for the optimizer.

We use the best amount of hidden states (2) according to the BIC that [Leroux and Puterman](#_bookmark31) ([1992](#_bookmark31)) reported. Although the AIC selects 3 hidden states, our goal is to show a simple example.

*# Model with 2 states*

m <- 2

TMB\_data <- **list**(x = lamb\_data, m = m)

*# Generate initial set of parameters for optimization*

lambda <- **c**(1, 3)

gamma <- **matrix**(**c**(0.8, 0.2,

0.2, 0.8), byrow = TRUE, nrow = m)

1. Transforming them to working parameters

*# Turn them into working parameters*

parameters <- **pois.HMM.pn2pw**(m, lambda, gamma)

1. Creating the TMB likelihood function object

obj\_tmb <- **MakeADFun**(TMB\_data, parameters, DLL = "poi\_hmm",

silent = TRUE)

1. Optimizing

mod\_tmb <- **nlminb**(start = obj\_tmb$par, objective = obj\_tmb$fn)

We are using nlminb because it is the fastest we have tested. Details are available in [item (iii)](#_bookmark13).

1. Getting the ML estimates of the natural parameters and their standard errors, sent by ADREPORT

It should be noted that the gamma parameter is shown column-wise below.

**summary**(**sdreport**(obj\_tmb), "report")

Biometrical Journal **52** (2020) 61 9

|  |  |  |  |
| --- | --- | --- | --- |
| ## |  | Estimate | Std. Error |
| ## | lambda | 0.25636529825 | 0.04016451459 |
| ## | lambda | 3.11475069040 | 1.02131434279 |
| ## | gamma | 0.98872130149 | 0.01063573157 |
| ## | gamma | 0.31033873770 | 0.18468644743 |
| ## | gamma | 0.01127869851 | 0.01063573157 |
| ## | gamma | 0.68966126230 | 0.18468644743 |
| ## | delta | 0.96493132139 | 0.03181447356 |
| ## | delta | 0.03506867861 | 0.03181447356 |

1. It is possible to tell the optimizer to use the gradient and/or hessian computed by TMB:

*# The negative log-likelihood is accessed by the objective*

*# attribute of the optimized object*

mod\_tmb$objective

## [1] 177.5188369

mod\_tmb <- **nlminb**(start = obj\_tmb$par, objective = obj\_tmb$fn, gradient = obj\_tmb$gr, hessian = obj\_tmb$he)

mod\_tmb$objective

## [1] 177.5188369

The use of an exact gradient and hessian provided by TMB can make the model fit better, although in our case it doesn’t make a difference.

The dataset used is provided by [Leroux and Puterman](#_bookmark31) ([1992](#_bookmark31)). Although our estimates are close to the paper’s estimates, they’re not exactly the same. The reason is that they use a modified likelihood and we don’t. This is clear when comparing the likelihoods with only 1 state.

A 1 state Poisson HMM is the same as a Poisson regression model, for which the log-likelihood has the expression

*l*(*λ*) = log

/ *n xi*

n

*λ*

*e*−*λ* \

*i*=1

*xi*!

/ *n* \ *n*

)

= −*nλ* + log(*λ*)

) *xi*

*i*=1

— log(*xi*!)*.*

*i*=1

They find a ML estimate *λ* = 0*.*3583 and a log-likelihood of -174.26.

x <- lamb\_data n <- **length**(x)

lambda\_leroux <- 0.3583

- n \* lambda\_leroux + **log**(lambda\_leroux) \* **sum**(x) - **sum**(**log**(**factorial**(x)))

## [1] -201.0436344

We can see that the log-likelihood is different, but when we remove the constant − *n*

*i*=1

log(*xi*!), it

becomes

10 First Author*et al.* and dd: Running title

- n \* lambda\_leroux + **log**(lambda\_leroux) \* **sum**(x)

## [1] -174.2610803

In our estimation procedures, we use the complete formula for the log-likelihood. In addition we don’t use the Expectation-Maximization (EM) algorithm like they do. Therefore our likelihood differs sligthly from the paper.

## Computing the stationary distribution

In the objective function above in [subsubsection 3.3.1](#_bookmark3), we calculated the stationary distribution of the *m*

state HMM’s Markov chain with transition probability matrix **Γ**, but we skipped over the code to do that.

To calculate the stationary distribution, we find the solution ***δ*** of the system:

***δ***(**I***m* − **Γ** + **U**) = **1**

where **I***m* is the *m* x *m* identity matrix, **U** is a *m* x *m* matrix of ones, and **1** is a row vector of ones. An implementation of this in R is

*# Compute the stationary distribution of a Markov chain*

*# with transition probability gamma*

stat.dist <- **function**(gamma) {

m <- **dim**(gamma)[1]

**return**(**solve**(**t**(**diag**(m) - gamma + 1), **rep**(1, m)))

}

More details are available in [Zucchini et al.](#_bookmark38) ([2016](#_bookmark38)).

In order to use it in TMB, we show a way to implement it in C++ in the file utils.cpp, see Appendix A.2 [item (iv)](#_bookmark20).

## Nested model specification

Nested models can be useful for multiple reasons. For example, there can be a need to fix some parameters because of some biological or physical constraints.

In our situation, using a nested model solves this issue at the cost of having a worse fit. The reason is that some parameters are fixed and the others’ estimates have different values than in the original model. Since the estimates can be different, the fit worsens.

To showcase the advantage of nested models, we will reuse the 2 state model.

TMB can be instructed to treat some parameters as constants. We will arbitrarily fix *λ*1 to 1.

*# Get the previous values, and fix some* fixed\_par\_lambda <- lambda fixed\_par\_lambda[1] <- 1 fixed\_par\_gamma <- gamma

Since we can only fix working parameters, we chose to fix a natural parameter for which we can eas- ily identify the corresponding working parameter to fix, for simplicity. Fixing a value in the transition probability matrix would be tricky, so we chose a Poisson mean instead.

Now that the nested model parameters have been made, we need to transform them into a set of working parameters.

Biometrical Journal **52** (2020) 61 11

*# Transform them into working parameters*

new\_parameters <- **pois.HMM.pn2pw**(m = m,

lambda = fixed\_par\_lambda, gamma = fixed\_par\_gamma)

In order to have TMB treat parameters as constants, the map argument of the MakeADFun function must be a list. This list must contain named vectors filled with NA if the parameter is to be treated as a constant, or factor levels.

Equal factor level is collected to a common value, so we want 1 different factor level for each non-fixed parameter, and NA for each fixed parameter. We take increasing numbers to make factors, and replace with NA where necessary.

map <- **list**(tlambda = **as.factor**(**c**(NA, 2)), tgamma = **as.factor**(**c**(1, 3)))

*# The map is fed to the MakeADFun function*

fixed\_par\_obj\_tmb <- **MakeADFun**(TMB\_data, new\_parameters,

DLL = "poi\_hmm", silent = TRUE, map = map)

fixed\_par\_mod\_tmb <- **nlminb**(start = fixed\_par\_obj\_tmb$par,

objective = fixed\_par\_obj\_tmb$fn, gradient = fixed\_par\_obj\_tmb$gr, hessian = fixed\_par\_obj\_tmb$he)

Old estimates and standard errors, and the new ones

Before After

|  |  |  |  |  |
| --- | --- | --- | --- | --- |
|  | Estimate | Std. Error | Estimate | Std. Error |
| *λ*1 | 0.26 | 0.04 | 1.00 | 0.00 |
| *λ*2 | 3.11 | 1.02 | 3.44 | 0.86 |
| *γ*1*,*1 | 0.99 | 0.01 | 1.00 | 0.00 |
| *γ*2*,*1 | 0.31 | 0.18 | 0.19 | 0.18 |
| *γ*1*,*2 | 0.01 | 0.01 | 0.00 | 0.00 |
| *γ*2*,*2 | 0.69 | 0.18 | 0.81 | 0.18 |
| *δ*1 | 0.96 | 0.03 | 0.98 | 0.03 |

*δ*2 0.04 0.03 0.02 0.03

**Table 1** Before and after fixing low parameters to 0 and the likelihood got worse as expected.

*# The negative log-likelihood is accessed by the attribute*

*# "objective" of the optimized object*

mod\_tmb$objective

## [1] 177.5188369

fixed\_par\_mod\_tmb$objective

## [1] 264.1635573

12 First Author*et al.* and dd: Running title

It should be noted that some inconsistencies can happen. For example, the stationary distribution is a vector of probabilities and should sum to 1. However, we can see that it doesn’t exactly sum to 1.

adrep <- **summary**(**sdreport**(obj\_tmb), "report")

estimate\_delta <- adrep[**rownames**(adrep) == "delta", "Estimate"]

**sum**(estimate\_delta)

## [1] 1

**sum**(estimate\_delta) == 1

## [1] FALSE

This is likely due to machine approximations when making numbers far apart from each other interact together. For example, in R a small number is not 0 but is treated as 0 when added to a much larger number.

1e-100 == 0

## [1] FALSE

(1 + 1e-100) == 1

## [1] TRUE

This can result in inconsistencies when numbers far apart from each other are involved. Fortunately, this doesn’t seem to cause any issues.

# Confidence intervals

## Hessian based confidence intervals

First, let us consider evaluation of parameter uncertainty via the Hessian. As the Hessian ∇2 log *L*(*ψ*ˆ) refers to the transformed parameters *ψ* the delta-method must be used to obtain an estimate of the covari- ance matrix of *θ*ˆ:

Σ*θ*ˆ = −∇*g*(*ψ*ˆ) ∇

( 2

log *L*(*ψ*ˆ)

'−1

∇*g*(*ψ*ˆ)! (2)

With minimal effort from the users perspective, TMB can be instructed to calculate Σ*θ*ˆ (by automatic differentiation). Standard errors of derived parameters, such as the stationary distribution ***δ*** which is a function of **Γ**, can be calculated by the delta-method similarly.

TMB provides an easy way to retrieve these. Following the example above, as we saw in [subsubsec-](#_bookmark4) [tion 3.3.2](#_bookmark4),

*# Get all standard errors*

adrep <- **summary**(**sdreport**(obj\_tmb), "report") adrep

##

Estimate

Std. Error

## lambda 0.25636540802 0.04016444898

## lambda 3.11475432301 1.02131181176

Biometrical Journal **52** (2020) 61 13

|  |  |  |  |
| --- | --- | --- | --- |
| ## | gamma | 0.98872127904 | 0.01063570936 |
| ## | gamma | 0.31033852657 | 0.18468647942 |
| ## | gamma | 0.01127872096 | 0.01063570936 |
| ## | gamma | 0.68966147343 | 0.18468647942 |
| ## | delta | 0.96493123100 | 0.03181444960 |
| ## | delta | 0.03506876900 | 0.03181444960 |

It should be noted that although the estimates can be found in the optimization variable opt in their working form, they are also sent to the MakeADFun object obj and can be retrieved as shown above. sdreport shows information about the working parameters’ estimates, whereas the summary of that report will also show information about the variables reported by the function ADREPORT.

*# More help with ?summary.sdreport*

We will now see how to get Wald confidence intervals using TMB. NEED CITATION With our 2 state Poisson HMM, the 100(1 − *α*)% confidence interval for *a* is *a* ± *z*1−*α/*2 ∗ *σa* where *zx* is the *x*-percentile of the standard normal distribution, and *σa* is the standard error of *a*.

|  |  |  |  |
| --- | --- | --- | --- |
| ## |  | lower\_bound | upper\_bound |
| ## | lambda | 0.177644534554 | 0.33508628149 |
| ## | lambda | 1.113019954968 | 5.11648869105 |
| ## | gamma | 0.967875671737 | 1.00956688634 |
| ## | gamma | -0.051640321537 | 0.67231737467 |
| ## | gamma | -0.009566886344 | 0.03212432826 |
| ## | gamma | 0.327682625328 | 1.05164032154 |
| ## | delta | 0.902576055600 | 1.02728640640 |
| ## | delta | -0.027286406403 | 0.09742394440 |

Estimates of ***θ*** and ***δ***, with accompanying confidence intervals are displayed in [Table 11](#_bookmark15) and [Table 12](#_bookmark16).

*# sdreport gives the estimates of the parameters and their standard errors*

adrep <- **summary**(**sdreport**(obj\_tmb), "report")

*# Get the 97.5 percentile of the standard normal distribution*

q95\_norm <- **qnorm**(1 - 0.05 / 2)

*# Create the CI*

*# Extract the values*

estimates <- adrep[, "Estimate"] std\_errors <- adrep[, "Std. Error"]

*# Create the bounds*

lower\_bound <- estimates - q95\_norm upper\_bound <- estimates + q95\_norm

*# Show the CI*

**cbind**(lower\_bound, upper\_bound)

\*

\*

std\_errors std\_errors

For comparison, we have included the corresponding standard deviations resulting from replacing ∇2 log *L*(*ψ*ˆ) in [(Equation 2](#_bookmark6)) with traditional numerical approximations using the R functions nlm (nlmb) and fdHess (nlme). The difference when using these approximations are of order 1*e* − 05 or less, however one would still have to calculate ∇*g*(*ψ*ˆ) for these approximations to be useful.

As mentioned earlier, for a larger amount of hidden states, TMB may be unable to give some or any

standard standard errors because some variables are close to their boundaries.

14 First Author*et al.* and dd: Running title

In that situation, using a nested model might solve this issue at the cost of the model usually having a worse fit.

Refer to [Section 3.4](#_bookmark5) for information on how to make a nested model using TMB.

## Likelihood profile based confidence intervals

Next, we consider evaluating uncertainty using likelihood-profiles. Assuming normality of the maximum likelihood (ML) estimators, confidence intervals for the parameters can be obtained using the above esti- mates of standard deviations. For a fixed sample size, the validity of the normality assumption can often be violated and in these cases a Likelihood-based CI is more robust (reference). Let *η* be single parameter in a model with parameters (*η, θ*) and likelihood *L*(*η, θ*), and let *Lp*(*η*) be the profile likelihood defined by *Lp*(*η*) = max*θ L*(*η, θ*). Then a likelihood-based CI for *η* is given by

(

*η* : 2 log(

*L*(*η*ˆ*, θ*ˆ) *Lp*(*η*)

2

*< χ*

1*,*(1−*α*)

(3)

where *χ*2

1*,*(1−*α*)

is the 1 − *α* quantile of a *χ*2 distribution with 1 degree of freedom. TMB allows for

very efficient computation of both the profile likelihood *L*(*η*) and the CI given by [(Equation 3](#_bookmark7)), and

this has been used to produce [Table 11](#_bookmark15) and [Table 12](#_bookmark16) which display the profile log-likelihood and the corresponding likelihood-based CI’s in the model for the lamb and the simulated dataset. Note that the problem of transformed parameters is easier to deal with when making likelihood-based CI’s, since

{*g*(*η*)*, Lp*(*g*(*η*))} = {*g*(*η*)*, Lp*(*η*)} for any one-to-one function *g* (invariance principle).

Again, TMB provides an easy way to obtain those intervals. The name argument allows to profile pa- rameter. The trace argument indicates how much information on the optimization you want the function to show. Following the HMM example, if we wish to profile the 2*nd* working parameter log lambda2, we have to feed its position into the name argument:

profile <- **tmbprofile**(obj = obj\_tmb,

name = 2, trace = FALSE)

*# Confidence interval of lambda*

**exp**(**confint**(profile))

##

lower

upper

## tlambda 1.265338844 4.947732633

It should be noted that the argument lincomb allows to profile any linear combination of parameters by using a vector of the linear coefficients in the same order as the parameters.

This method can sometimes give NA values, usually when a ML estimate is close to a boundary, but not only. For example, the first working parameter log lambda is pretty low (−4*.*47). It becomes too difficult to profile the likelihood so the function fails to provide a meaningful confidence interval.

Another important issue is about profiling with a univariate model. With only 1 (hidden) state, a Poisson HMM becomes a univariate Poisson regression model. Therefore, the profile becomes a plot of the likeli- hood when the Poisson mean varies. However, tmbprofile fails to provide a confidence interval. The reason is likely that it tries to optimize the likelihood despite the lack of parameters to change, and there- fore fails. [Visser et al.](#_bookmark37) ([2000](#_bookmark37)), [Meeker and Escobar](#_bookmark34) ([1995](#_bookmark34)) and [Venzon and Moolgavkar](#_bookmark36) ([1988](#_bookmark36)) provide more details on profiling likelihoods.

Biometrical Journal **52** (2020) 61 15

## Bootstrap based confidence intervals

Finally, we consider evaluating uncertainty using bootstrap. There are many ways to bootstrap. Non- parametric bootstraping of time series is possible, as [Ha¨rdle et al.](#_bookmark28) ([2003](#_bookmark28)) has shown. But since our focus is not on bootstrapping techniques, we chose a parametric approach. See **?** for more details on bootstrapping. From the parameters’ ML estimates, we generate new data and re-estimate the parameters 500 times.

From that list of new estimates we can get the 2.5th and 97.5th percentiles and get 95% confidence intervals for the parameters.

We show below how we get confidence intervals using bootstrap, based on the 3 state Poisson HMM estimates from above.

1. First, we need a function to generate random data from a HMM.

*# Generate a random sample from a HMM*

pois.HMM.generate\_sample <- **function**(ns, mod) {

mvect <- 1:mod$m state <- **numeric**(ns)

state[1] <- **sample**(mvect, 1, prob = mod$delta)

**for** (i **in** 2:ns) {

state[i] <- **sample**(mvect, 1, prob = mod$gamma[state[i - 1], ])

}

x <- **rpois**(ns,lambda = mod$lambda[state])

**return**(x)

}

1. Then, when the model is estimated each time, we don’t impose an order for the states. This can lead to the label switching problem, where states aren’t ordered the same way in each model. To address this, we re-ordered the states by ascending Poisson means.

Sorting the means is pretty straightforward. Re-ordering the TPM is a little trickier. To do so, we took the permutations of the states given by the sorted Poisson means, and permuted each row index and column index to its new value.

The function we used is

*# Relabel states by increasing Poisson means*

pois.HMM.label.order <- **function**(m, lambda, gamma, delta = **NULL**) {

*# Get the indexes of the sorted states according to ascending lambda*

sorted\_lambda <- **sort**(lambda, index.return = TRUE)$ix

*# Re-order the TPM according to the switched states in the sorted lambda*

ordered\_gamma <- **matrix**(0, nrow = m, ncol = m)

**for** (col **in** 1:m) {

new\_col <- **which**(sorted\_lambda == col)

**for** (row **in** 1:m) {

new\_row <- **which**(sorted\_lambda == row) ordered\_gamma[row, col] <- gamma[new\_row, new\_col]

}

}

*# Re-order the stationary distribution if it was provided*

*# Generate it otherwise*

**if** (**is.null**(delta)) {

16 First Author*et al.* and dd: Running title

delta <- **stat.dist**(ordered\_gamma)

} **else** {

delta <- delta[sorted\_lambda]

}

**return**(**list**(lambda = **sort**(lambda),

gamma = ordered\_gamma, delta = delta))

}

Let’s show an example to understand the process. For readability, the TPM is filled with row and column indexes instead of probabilities.

lambda <- **c**(30, 10, 20)

gamma <- **matrix**(**c**(11, 12, 13,

21, 22, 23,

31, 32, 33), byrow = TRUE, ncol = 3)

**pois.HMM.label.order**(m = 3, lambda, gamma)

## $lambda

## [1] 10 20 30

##

## $gamma

##

## [1,]

## [2,]

## [3,]

##

[,1] [,2] [,3]

33

13

23

31

11

21

32

12

22

## $delta

## [1] -0.032786885246 0.016393442623 -0.008196721311

State 1 has been relabeled state 3, state 2 became state 1, and state 3 became state 2.

In short, 1 → 3*,* 2 → 1*,*and3 → 2.

1. Bootstrap code

bootstrap\_estimates <- **data.frame**() DATA\_SIZE <- **length**(lamb\_data)

*# Set how many parametric bootstrap samples we create*

BOOTSTRAP\_SAMPLES <- 10

*# ML parameters*

ML\_working\_estimates <- obj\_tmb$env$last.par.best ML\_natural\_estimates <- obj\_tmb$**report**(ML\_working\_estimates) gamma <- ML\_natural\_estimates$gamma

lambda <- ML\_natural\_estimates$lambda delta <- ML\_natural\_estimates$delta

*# Parameters for TMB*

cols <- **names**(ML\_working\_estimates)

Biometrical Journal **52** (2020) 61 17

tgamma <- ML\_working\_estimates[cols == "tgamma"] tlambda <- ML\_working\_estimates[cols == "tlambda"] ML\_TMB\_parameters <- **list**(tlambda = tlambda,

tgamma = tgamma) params\_names <- **c**("lambda", "gamma", "delta")

**for** (idx\_sample **in** 1:BOOTSTRAP\_SAMPLES) {

mod\_temp <- **NULL**

natural\_parameters <- **NULL**

*# This while loop is in case tmb doesn't converge for some reason*

*# And in case some parameter estimates are NA values*

**while** (**is.null**(mod\_temp) |

**anyNA**(natural\_parameters[params\_names], recursive = TRUE)) {

*#simulate the data*

bootstrap\_data <- **pois.HMM.generate\_sample**(ns = DATA\_SIZE,

mod = **list**(m = m,

lambda = lambda, gamma = gamma, delta = delta))

*# Parameters for TMB*

TMB\_data\_bootstrap <- **list**(x = bootstrap\_data, m = m)

*# Estimate the parameters*

obj <- **MakeADFun**(TMB\_data\_bootstrap,

ML\_TMB\_parameters, DLL = "poi\_hmm", silent = TRUE)

*# Sometimes TMB doesn't converge, maybe the sample*

*# has too many extremes*

*# In those cases, the data is regenerated through the "while"*

*# The error messages are useless in those cases*

**try**(mod\_temp <- **nlminb**(start = obj$par, objective = obj$fn,

gradient = obj$gr, hessian = obj$he),

silent = TRUE)

**if** (!**is.null**(mod\_temp)) {

working\_parameters <- obj$env$last.par.best natural\_parameters <- obj$**report**(working\_parameters)

*# Label switching*

natural\_parameters <- **pois.HMM.label.order**(m,

natural\_parameters$lambda, natural\_parameters$gamma,

18 First Author*et al.* and dd: Running title

natural\_parameters$delta)

}

}

*# The values from gamma are taken columnwise* natural\_parameters <- **unlist**(natural\_parameters[params\_names]) len\_par <- **length**(natural\_parameters)

bootstrap\_estimates[idx\_sample, 1:len\_par] <- natural\_parameters

}

q <- **apply**(bootstrap\_estimates, 2, **function**(par\_estimate) {

**quantile**(par\_estimate, probs = **c**(0.025, 0.975))

})

*# Get row column indexes for gamma instead of the default column*

params\_names <- **paste0**(**rep**("lambda", m), 1:m)

**for** (gamma\_idx **in** 1:m ˆ 2) {

row\_col\_idx <- **matrix.col.idx.to.rowcol**(gamma\_idx, m) params\_names <- **c**(params\_names,

**paste0**("gamma",

**paste0**(row\_col\_idx, collapse = "")))

}

params\_names <- **c**(params\_names,

**paste0**(**rep**("delta", m), 1:m))

bootstrap\_CI <- **data.frame**("Parameter" = params\_names,

"Estimate" = **c**(lambda, gamma, delta), "Lower bound" = q[1, ],

"Upper bound" = q[2, ])

**print**(bootstrap\_CI, right = FALSE, row.names = FALSE)

|  |  |  |  |  |
| --- | --- | --- | --- | --- |
| ## | Parameter | Estimate | Lower.bound | Upper.bound |
| ## | lambda1 | 0.25636540802 | 7.117338887e-09 | 0.2954041084 |
| ## | lambda2 | 3.11475432301 | 3.565352892e-01 | 3.4379317191 |
| ## | gamma11 | 0.98872127904 | 4.444232052e-01 | 0.9947809939 |
| ## | gamma21 | 0.31033852657 | 9.937752479e-02 | 0.5693025943 |
| ## | gamma12 | 0.01127872096 | 5.219006082e-03 | 0.5555767948 |
| ## | gamma22 | 0.68966147343 | 4.306974057e-01 | 0.9006224752 |
| ## | delta1 | 0.96493123100 | 2.635813800e-01 | 0.9830019462 |
| ## | delta2 | 0.03506876900 | 1.699805382e-02 | 0.7364186200 |

1. It should be noted that some estimates can be very large or small. This can happen when the randomly generated bootstrap sample contains long chains of the same values. However, a large number of bootstrap samples lowers that risk since we ”leave out” 5% of the extreme values in the CI. It could be useful to skip the bootstrap samples where the estimates are too far out of expectations.

Biometrical Journal **52** (2020) 61 19

# State inference

## Setup

Given an optimized MakeADFun object obj, we need to setup some variables the probabilities detailed below.

*# Retrieve the objects at ML value*

adrep <- obj\_tmb$**report**(obj\_tmb$env$last.par.best) delta <- adrep$delta

gamma <- adrep$gamma

emission\_probs <- adrep$emission\_probs n <- adrep$n

m <- **length**(delta) mllk <- adrep$mllk

## Log-forward probabilities

The forward probabilities have been detailed in [Section 3.2](#_bookmark2). We show here a way to compute the log of the forward probabilities, using a scaling scheme defined by [Zucchini et al.](#_bookmark38) ([2016](#_bookmark38)).

*# Compute log-forward probabilities (scaling used)*

lalpha <- **matrix**(NA, m, n)

foo <- delta \* emission\_probs[1, ] sumfoo <- **sum**(foo)

lscale <- **log**(sumfoo) foo <- foo / sumfoo

lalpha[, 1] <- **log**(foo) + lscale

**for** (i **in** 2:n) {

foo <- foo %\*% gamma \* emission\_probs[i, ] sumfoo <- **sum**(foo)

lscale <- lscale + **log**(sumfoo) foo <- foo / sumfoo

lalpha[, i] <- **log**(foo) + lscale

}

*# Since lalpha contains n=240 columns, we only display 5 for readability*

lalpha[, 1:5]

## [,1] [,2] [,3] [,4] [,5]

## [1,] -0.2920638514 -0.5591179415 -0.8265947717 -1.094088418 -1.361582733

## [2,] -6.4651986397 -7.7716768804 -8.1146145025 -8.385231634 -8.652850430

## Log-backward probabilities

The backward probabilities have been defined in the same section.

*# Compute log-backwards probabilities (scaling used)*

lbeta <- **matrix**(NA, m, n) lbeta[, n] <- **rep**(0, m) foo <- **rep** (1 / m, m)

20 First Author*et al.* and dd: Running title

lscale <- **log**(m)

**for** (i **in** (n - 1):1) {

foo <- gamma %\*% (emission\_probs[i + 1, ] \* foo) lbeta[, i] <- **log**(foo) + lscale

sumfoo <- **sum**(foo) foo <- foo / sumfoo

lscale <- lscale + **log**(sumfoo)

}

*# Since lbeta contains n=240 columns, we only display 5 for readability*

lbeta[, 1:5]

## [,1] [,2] [,3] [,4] [,5]

## [1,] -177.2274543 -176.9599600 -176.6924658 -176.4249753 -176.1575774

## [2,] -178.3455913 -178.0780696 -177.8098901 -177.5253277 -176.9034311

## Smoothing probabilities

The smoothing probabilities are defined in [Zucchini et al.](#_bookmark38) ([2016](#_bookmark38)) as *P* (*Ct* = *i*|*X*(*N* ) = *x*(*N* )) = *αt* (*i*)*βt* (*i*) .

*# Compute conditional state probabilities, smoothing probabilities*

stateprobs <- **matrix**(NA, ncol = n, nrow = m) llk <- - mllk

**for**(i **in** 1:n) {

stateprobs[, i] <- **exp**(lalpha[, i] + lbeta[, i] - llk)

}

*# Most probable states*

ldecode <- **rep**(NA, n)

**for** (i **in** 1:n) {

ldecode[i] <- **which.max**(stateprobs[, i])

}

ldecode

*LT*

|  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- |
| ## | [1] | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 |
| ## | [38] | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 |
| ## | [75] | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 2 | 2 | 2 | 2 | 2 | 2 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 |
| ## | [112] | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 |
| ## | [149] | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 |
| ## | [186] | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 2 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 |
| ## | [223] | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |

## Forecast, h-step-ahead-probabilities

The forecast distribution or h-step-ahead-probabilities as well as its implementation in R is detailed in [Zucchini et al.](#_bookmark38) ([2016](#_bookmark38)).

Let ***ϕ****N* = ***α****N* .

***α****N* **1***!*

Biometrical Journal **52** (2020) 61 21

Then,

*P* (*XN* +*h* = *x*|*X*(*N* ) = *x*(*N* )) =

***α****N* **Γ***h***P**(*x*)**1**!

***α****N* **1**!

= ***ϕ****N* **Γ***h***P**(*x*)**1**!

An implementation of this, using a scaling scheme is

*# Number of steps*

h <- 1

*# Values for which we want the forecast probabilities*

xf <- 0:50

nxf <- **length**(xf)

dxf <- **matrix**(0, nrow = h, ncol = nxf) foo <- delta \* emission\_probs[1, ] sumfoo <- **sum**(foo)

lscale <- **log**(sumfoo) foo <- foo / sumfoo

**for** (i **in** 2:n) {

foo <- foo %\*% gamma \* emission\_probs[i, ] sumfoo <- **sum**( foo)

lscale <- lscale + **log**(sumfoo) foo <- foo / sumfoo

}

emission\_probs\_xf <- **get.emission.probs**(xf, lambda)

**for** (i **in** 1:h) {

foo <- foo %\*% gamma

**for** (j **in** 1:m) {

dxf[i, ] <- dxf[i, ] + foo[j] \* emission\_probs\_xf[, j]

}

}

*# Since lbeta contains n=240 columns, we only display 5 for readability*

dxf[, 5]

## [1] 0.002181533178

## Global decoding using the Viterbi algorithm

The Viterbi algorithm is detailed in [Zucchini et al.](#_bookmark38) ([2016](#_bookmark38)). It calculates the sequence of states (*i*∗*, . . . , i*∗ )

1 *N*

which maximizes the conditional probability of all states simultaneously, i.e.

(*i*∗*, . . . , i*∗ ) = arg max

*P* (*C*1 = *i*1*, . . . , CN* = *iN* |*X*(*N* ) = *x*(*N* ))

1 *N*

*i*1 *,...,iN* ∈{1*,...,m*}

An implementation of it is

xi <- **matrix**(0, n, m)

foo <- delta \* emission\_probs[1, ] xi[1, ] <- foo / **sum**(foo)

**for** (i **in** 2:n) {

22 First Author*et al.* and dd: Running title

foo <- **apply**(xi[i - 1, ] \* gamma, 2, max) \* emission\_probs[i, ] xi[i, ] <- foo / **sum**(foo)

}

iv <- **numeric**(n)

iv[n] <- **which.max**(xi[n, ])

**for** (i **in** (n - 1):1){

iv[i] <- **which.max**(gamma[, iv[i + 1]] \* xi[i, ])

|  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- |
| }  iv |  | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| ## | [1] | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 |
| ## | [38] | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 |
| ## | [75] | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 2 | 2 | 2 | 2 | 2 | 2 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 |
| ## | [112] | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 |
| ## | [149] | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 |
| ## | [186] | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 2 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 |
| ## | [223] | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 | 1 |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |

# Application to hospital data

## Data description

We timed different parameters on a dataset provided by [Leroux and Puterman](#_bookmark31) ([1992](#_bookmark31)), on a simulated dataset, and on a larger dataset provided by a hospital from Assistance publique – Hoˆpitaux de Paris, a french hospital trust.

The first dataset is the number of movements by a fetal lamb in 240 consecutive 5-second intervals, and

is:

0 0 0 0 0 1 0 1 0 0 0 0 0 0 1 0 1 0 0 0 0 2 2 0 0 0 0 1 0 0 1 1 0 0 1 1 1 0 0 1 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0

2 0 1 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 1 0 0 0 0 0 7 3 2 3 2 4 0 0 0 0 1 0 0 0 0 0 0 0 1 0 2 0 0 0 0 1 0 0 0 0 1 0

0 0 0 0 0 0 0 0 0 0 0 0 0 1 0 0 0 0 0 2 1 0 0 1 0 0 0 1 0 1 1 0 0 0 1 0 0 1 0 0 0 1 2 0 0 0 1 0 1 1 0 1 0 0 2 0 1 2

1 1 2 1 0 1 1 0 0 1 1 0 0 0 1 1 1 0 4 0 0 2 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0

0 0 0 0 0 0 0 0

**Table 2** Numbers of movements by a fetal lamb in 240 consecutive 5-second intervals

The second is a sequence of random numbers generated by a *m*-state Poisson HMM. The code we used

is

DATA\_SIZE\_SIMULATION <- 2000

m <- 2

*# Generate parameters*

lambda <- **seq**(1, 7, length.out = m)

*# Create the transition probability matrix with 0.8 on its diagonal*

**if** (m == 1) {

gamma <- **diag**(1, m)

} **else** {

gamma <- **matrix**(0.2 / (m - 1), nrow = m, ncol = m)

**diag**(gamma) <- 0.8

}

Biometrical Journal **52** (2020) 61 23

delta <- **stat.dist**(gamma)

*#simulate the data*

simul\_data <- **pois.HMM.generate\_sample**(ns = DATA\_SIZE\_SIMULATION,

mod = **list**(m = m,

lambda = lambda, gamma = gamma, delta = delta))

The last is the number of patients who entered a hospital each hour between the first hour of 2010 and the last hour of 2019. The dataset is available at GITHUB REPO DATASET WEBSITE. The zeroes in the amount are not missing data but indicate that no one arrived at the hospital. The data was grouped using the package dplyr. The column DATE is a datetime item useful for sorting and plotting the data. The columns YEAR is an integer taking values 2010, 2011, . . . , 2019. The column PATIENTS represents the number of people arriving at the hospital between the hour mentioned and the next. For example, the dataset starts on the hour 0 with an amount of 6, and indicates that 6 people arrived between 00:00 and 00:59. The column WDAY represents the number of the day in the week: 1 is Monday and 7 is Sunday.

## Speed comparison

Different numbers of hidden states *m* were used: (1*,* 2 for the lamb dataset, and 1*,* 2*,* 3 for the simulated dataset) to compare speeds.

Overall, the acceleration grows with the amount of hidden states *m* and the size of the dataset. For readability, DM denotes the estimation time without using TMB.

TMB1, ..., TMB4 will be used depending on whether TMB gives the gradient and hessian. The following table will summarize the notation

**Table 3** Naming of TMB parameters

|  |  |  |  |  |
| --- | --- | --- | --- | --- |
|  | TMB1 | TMB2 | TMB3 | TMB4 |
| Exact gradient is used  Exact hessian is used | No  No | No  Yes | Yes  No | Yes  Yes |

The following [Figure 1](#_bookmark9) below shows the time acceleration when using TMB on the hospital dataset through boxplots, for multiple amounts of hidden states *m*.

24 First Author*et al.* and dd: Running title

Parameter estimation time

**Hospital data, size = 87648**

100.0

1

2

3

4

10.0

Time (seconds)

1.0

0.1

DM TMB1 TMB2 TMB3 TMB4 DM TMB1 TMB2 TMB3 TMB4 DM TMB1 TMB2 TMB3 TMB4 DM TMB1 TMB2 TMB3 TMB4

Exact/inexact gradient and hessian

**Figure 1** Time w/o using TMB with hospital data

Biometrical Journal **52** (2020) 61 25

The average speed increases from using TMB on the hospital dataset are summarized here:

|  |  |  |  |  |
| --- | --- | --- | --- | --- |
| m | TMB1 | TMB2 | TMB3 | TMB4 |
| 1 | 1420.29 | 1424.77 | 1184.13 | 1187.47 |
| 2 | 1629.22 | 1616.27 | 2860.31 | 2290.07 |
| 3 | 1042.35 | 1016.19 | 3595.52 | 1392.61 |
| 4 | 760.66 | 758.42 | 4634.25 | 1147.34 |

**Table 4** Speed percentage increase of TMB estimation on the hospital dataset

## Interpretation

Here are the boxplots of hourly arrivals.

**Hourly arrivals**

40



20

PATIENTS

0

0 1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23

HOUR

As could have been guessed, more patients arrive in the day rather than during the night, in the afternoon rather than in the evening. Also, most patients show up in the early part of the day and around noon. From this, we can see different distributions in the hourly arrivals of patients, justifying 3 or 4 hidden states.

Here we can see the distribution of the states of a 4 state Poisson HMM fitted on the hourly arrivals.

26 First Author*et al.* and dd: Running title

3000

1

2

3

4

5

7

8

9

10

11 12 13

14

15

16

17 18 19 20

21

22

2000

amount

state

1

2

3

4

1000

0

0 6 23

1234 1234 1234 1234 1234 1234 123 1234 1234 1234 1234 1234 1234 1234 1234 1234 1234 1234 1234 1234 1234 1234 1234 123

hour

This graph shows the amount of states per hour across the entire dataset. The states (1, 2, 3, 4) have means (3.816, 8.64, 13.186, 22.503) respectively. This model seems to agree with our guess.

Here is a table of the negative log-likelihoods corresponding to different values of *m*.

|  |  |  |  |
| --- | --- | --- | --- |
| m | nll | AIC | BIC |
| 1 | 324288.01 | 648578.03 | 648587.41 |
| 2 | 259295.22 | 518598.43 | 518635.96 |
| 3 | 247402.86 | 494823.72 | 494908.15 |
| 4 | 242586.76 | 485205.52 | 485355.62 |

**Table 5** Negative log-likelihood values corresponding to different amounts of Poisson HMM states The 4 state model has the best likelihood, the best AIC, and the best BIC. It seems it has the best fit.

# 7 Speed evaluation

## 7.1 Speed comparison

1. Speed of TMB

To compare speeds with and without using TMB, it is important to test multiple sets of parameters. To do so, we used the R package microbenchmark.

The following [Figure 2](#_bookmark10) below shows the time acceleration when using TMB on the lamb dataset from [Leroux and Puterman](#_bookmark31) ([1992](#_bookmark31)) through boxplots.

Biometrical Journal **52** (2020) 61 27

Parameter estimation time

**Lamb data, size = 240m = 2**

0.030

0.010

Time (seconds)

0.003

DM TMB1 TMB2 TMB3 TMB4

Exact/inexact gradient and hessian

**Figure 2** Time w/o using TMB with lamb data

28 First Author*et al.* and dd: Running title

The average speed increases using the lamb dataset are summarized here:

|  |  |  |  |  |
| --- | --- | --- | --- | --- |
| m | TMB1 | TMB2 | TMB3 | TMB4 |
| 1 | 2036.39 | 1860.74 | 2092.40 | 1809.81 |
| 2 | 2271.13 | 2219.18 | 4459.56 | 1617.31 |

**Table 6** Speed percentage increase of TMB estimation on the lamb dataset

The following [Figure 3](#_bookmark11) below shows the time acceleration when using TMB on a simulated dataset through boxplots.

Parameter estimation time

**Simulated data, size = 2000**

1.00

0.10

Time (seconds)

0.01

|  |  |  |  |
| --- | --- | --- | --- |
| 1 |  | 2 | 3 |
| DM TMB1 TMB2 TMB3 | TMB4 | DM TMB1 TMB2 TMB3 TMB4 | DM TMB1 TMB2 TMB3 TMB4 |
|  |  | Exact/inexact gradient and hessian |  |

**Figure 3** Time w/o using TMB with simulated data

Biometrical Journal **52** (2020) 61 29

The average speed increases using a simulated dataset are summarized here:

|  |  |  |  |  |
| --- | --- | --- | --- | --- |
| m | TMB1 | TMB2 | TMB3 | TMB4 |
| 1 | 2072.39 | 2020.43 | 1929.44 | 1705.73 |
| 2 | 1959.55 | 1763.48 | 3589.45 | 3175.87 |
| 3 | 1102.32 | 1095.07 | 4168.62 | 3263.75 |

**Table 7** Speed percentage increase of TMB estimation on the simulated dataset

30 First Author*et al.* and dd: Running title

1. Log-likelihoods

TMB also allows for a better fit of a HMM. As we can see in the following [Table 8](#_bookmark12), the likelihood gets better when TMB can use the exact gradient.

|  |  |  |  |  |
| --- | --- | --- | --- | --- |
| m | procedure | nll | AIC | BIC |
| 2 | DM | 177.52 | 363.04 | 376.96 |
| 2 | TMB1 | 177.52 | 363.04 | 376.96 |
| 2 | TMB2 | 177.52 | 363.04 | 376.96 |
| 2 | TMB3 | 177.52 | 363.04 | 376.96 |
| 2 | TMB4 | 177.52 | 363.04 | 376.96 |

**Table 8** Negative log-likelihood values in lamb dataset

Moreover, the computation time of the likelihood is better with TMB although all of the times are relatively low, as we can see in the following tables.

|  |  |  |  |  |
| --- | --- | --- | --- | --- |
| m | TMB1 | TMB2 | TMB3 | TMB4 |
| 1 | 2001.31 | 2091.77 | 2091.37 | 2029.49 |
| 2 | 1994.02 | 2073.93 | 2095.03 | 2066.85 |

**Table 9** Speed percentage increase of TMB likelihood calculation on the lamb dataset

|  |  |  |  |  |
| --- | --- | --- | --- | --- |
| m | TMB1 | TMB2 | TMB3 | TMB4 |
| 1 | 2085.74 | 2079.52 | 2066.96 | 2086.84 |
| 2 | 1784.69 | 1780.48 | 1780.22 | 1779.60 |
| 3 | 864.12 | 887.50 | 895.06 | 863.82 |

**Table 10** Speed percentage increase of TMB likelihood calculation on the simulated dataset

Biometrical Journal **52** (2020) 61 31

1. All optimization methods

Finally, we compared different optimization methods. The ones we retained were BFGS, Nelder- Mead, L-BFGS-B, nlm, nlminb, and hjn, because the others didn’t converge. We also added the algorithm for least-squares curve fitting from the package marqLevAlg. Exact gradients and hes- sians are provided by TMB and fed to each algorithm. The speed comparisons are in the following [Figure 4](#_bookmark14).

**Lamb data, size = 240, m = 2**

Parameter estimation time

**Simulated data, size = 2000, m = 3**

Parameter estimation time

5

0.5

3

Time (milliseconds)

Time (milliseconds)

0.3

2

0.2

BFGS L−BFGS−B nlm nlminb hjn marqLevAlg

Method

BFGS L−BFGS−B nlm nlminb hjn marqLevAlg

Exact/inexact gradient and hessian

**Figure 4** Log estimation time vs optimization method

32 First Author*et al.* and dd: Running title

The following tables summarize the estimates and their confidence intervals, for the lamb dataset, and for the simulated dataset. Instead of showing the standard error, we show the lower and upper bounds of the confidence intervals. The reason is that sometimes, only one bound of the interval is known and doesn’t let one have a typical standard error.

|  |  |  |  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- | --- | --- | --- |
| m | Parameter | Parameter.estimate | Profile.L | Profile.U | Bootstrap.L | Bootstrap.U | TMB.L | TMB.U |
| 1 | *λ*1 | 0.36 |  |  | 0.29 | 0.44 | 0.28 | 0.43 |
| 2 | *λ*1 | 0.26 | 0.15 | 0.33 | 0.00 | 0.33 | 0.18 | 0.34 |
| 2 | *λ*2 | 3.11 | 1.27 | 4.95 | 0.27 | 5.19 | 1.11 | 5.12 |
| 2 | *γ*1*,*1 | 0.99 | 0.93 | 1.00 | 0.48 | 1.00 | 0.97 | 1.01 |
| 2 | *γ*2*,*1 | 0.31 | 0.04 | 0.68 | 0.03 | 1.00 | -0.05 | 0.67 |
| 2 | *γ*1*,*2 | 0.01 | 0.00 | 0.07 | 0.00 | 0.52 | -0.01 | 0.03 |
| 2 | *γ*2*,*2 | 0.69 | 0.32 | 0.96 | 0.00 | 0.97 | 0.33 | 1.05 |
| 2 | *δ*1 | 0.96 |  |  | 0.17 | 0.99 | 0.90 | 1.03 |
| 2 | *δ*2 | 0.04 |  |  | 0.01 | 0.83 | -0.03 | 0.10 |

**Table 11** Estimates and standard errors on the lamb dataset

|  |  |  |  |  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- | --- | --- | --- | --- |
| m | Parameter | Parameter.value | Parameter.estimate | Profile.L | Profile.U | Bootstrap.L | Bootstrap.U | TMB.L | TMB.U |
| 1 | *λ*1 | 1.00 | 1.01 |  |  | 0.96 | 1.05 | 0.96 | 1.05 |
| 2 | *λ*1 | 1.00 | 0.98 |  |  | 0.91 | 1.04 | 0.92 | 1.04 |
| 2 | *λ*2 | 20.00 | 20.04 |  |  | 19.77 | 20.30 | 19.76 | 20.32 |
| 2 | *γ*1*,*1 | 0.80 | 0.79 |  |  | 0.76 | 0.82 | 0.77 | 0.82 |
| 2 | *γ*2*,*1 | 0.20 | 0.22 |  |  | 0.19 | 0.25 | 0.19 | 0.24 |
| 2 | *γ*1*,*2 | 0.20 | 0.21 |  |  | 0.18 | 0.24 | 0.18 | 0.23 |
| 2 | *γ*2*,*2 | 0.80 | 0.78 |  |  | 0.75 | 0.81 | 0.76 | 0.81 |
| 2 | *δ*1 | 0.50 | 0.51 |  |  | 0.47 | 0.55 | 0.47 | 0.55 |
| 2 | *δ*2 | 0.50 | 0.49 |  |  | 0.45 | 0.53 | 0.45 | 0.53 |
| 3 | *λ*1 | 1.00 | 0.97 | 0.89 | 1.05 | 0.90 | 1.04 | 0.89 | 1.04 |
| 3 | *λ*2 | 10.50 | 10.57 | 10.29 | 10.85 | 10.28 | 10.86 | 10.29 | 10.85 |
| 3 | *λ*3 | 20.00 | 20.10 | 19.70 | 20.51 | 19.69 | 20.53 | 19.70 | 20.51 |
| 3 | *γ*1*,*1 | 0.80 | 0.80 | 0.76 | 0.84 | 0.77 | 0.83 | 0.77 | 0.83 |
| 3 | *γ*2*,*1 | 0.10 | 0.10 | 0.08 | 0.12 | 0.08 | 0.12 | 0.07 | 0.12 |
| 3 | *γ*3*,*1 | 0.10 | 0.10 | 0.08 | 0.12 | 0.07 | 0.12 | 0.07 | 0.12 |
| 3 | *γ*1*,*2 | 0.10 | 0.09 | 0.07 | 0.11 | 0.07 | 0.12 | 0.07 | 0.11 |
| 3 | *γ*2*,*2 | 0.80 | 0.82 | 0.77 | 0.86 | 0.78 | 0.84 | 0.78 | 0.85 |
| 3 | *γ*3*,*2 | 0.10 | 0.11 | 0.09 | 0.14 | 0.09 | 0.14 | 0.08 | 0.14 |
| 3 | *γ*1*,*3 | 0.10 | 0.11 | 0.09 | 0.13 | 0.09 | 0.14 | 0.08 | 0.13 |
| 3 | *γ*2*,*3 | 0.10 | 0.09 | 0.06 | 0.11 | 0.06 | 0.11 | 0.06 | 0.11 |
| 3 | *γ*3*,*3 | 0.80 | 0.79 | 0.74 | 0.83 | 0.75 | 0.82 | 0.75 | 0.82 |
| 3 | *δ*1 | 0.33 | 0.33 |  |  | 0.28 | 0.38 | 0.28 | 0.38 |
| 3 | *δ*2 | 0.33 | 0.36 |  |  | 0.30 | 0.41 | 0.30 | 0.41 |
| 3 | *δ*3 | 0.33 | 0.31 |  |  | 0.27 | 0.36 | 0.27 | 0.36 |

**Table 12** Estimates and standard errors on the simulated dataset

All the code used to produce those results is available in the GitHub repository GITHUB REPO. The files executing the code are poi hmm lamb.R, poi hmm simul.R, and poi hmm hosp.R.

# 8 Discussion

1. Throughout this paper, when we actually used hidden states (*m >* 1), we found that TMB3 yields the greatest speed increase (when TMB’s exact gradient is provided to the optimizer but not the hessian). This also stayed true when using other optimizers such as nlm. The reason is unclear. Although adding the exact hessian on top of the gradient to the optimizer seems to slow down the computation, it might help the optimizer converge in some cases, and could be worth investigating.
2. When providing only TMB’s hessian, optimizers sometimes fail to converge.

Biometrical Journal **52** (2020) 61 33

1. Mention the prospect of using TMB for panel data with random effects, Laplace approximation.

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## Conflict of Interest

*The authors have declared no conflict of interest.*

# Appendix

## R Code

1. Code to setup global parameters and declare functions used
2. Packages used
3. Functions used in TMB related code
4. Functions used in non TMB related code
5. Code to run estimations and comparisons using the lamb dataset
6. Code to run estimations and comparisons using a simulated dataset
7. Code to run estimations and comparisons using the hospital dataset

## C++ Code

1. Poisson HMM negative log-likelihood calculation This is the file poi hmm.cpp which contains the negative likelihood function.

It should be noted that in the C++ file, testing if the value is missing (i.e. testing for NaN (Not A Number) values) requires a little trick. The reason is that the standard test function std::isnan() doesn’t currently work on a single data value inside TMB.

Cuurently in C++, comparisons involving NaN values are always false, except when testing inequality between 2 NaN values. In other words, for a float f, the expression f != f will be true if and only if f is a NaN value. Similarly, f == f returns false if and only if f is a NaN value.

**# i n c l u d e** *<*TMB. hpp*>*

**# i n c l u d e** ” . . / f u n c t i o n s / u t i l s . cpp ”

*/ / L i k e l i h o o d f o r a poisson h i dden markov model .*

**template** *<***c l a s s** Type*>*

Type o b j e c t i v e f u n c t i o n *<*Type *>* : : **operator** ( ) ( )

*{*

*/ / Data*

DATA VECTOR( x ) ; DATA INTEGER(m) ;

*/ / t i m e s e r i e s v e c t o r*

*/ / Number o f s t a t e s m*

*/ / Parameters*

PARAMETER VECTOR( t l ambda ) ;

*/ / c o n d i t i o n a l l o g s d ' s*

34 First Author*et al.* and dd: Running title

PARAMETER VECTOR( tgamma ) ; */ / m( m−1) working parameters o f TPM*

*/ / Uncomment only u s i ng a non s t a t i o n a r y d i s t r i b u t i o n*

*/ / PARAMETER VECTOR( t d e l t a ) ; / / t r ansformed s t a t i o n a r y d i s t r i b u t i o n ,*

*/ / Transform working parameters t o n a t u r a l parameters :*

v e c t o r *<*Type*>* lambda = t l ambda . exp ( ) ;

matrix *<*Type*>* gamma = Gamma w2n (m, tgamma ) ;

*/ / C o n s t r u c t s t a t i o n a r y d i s t r i b u t i o n*

v e c t o r *<*Type*>* d e l t a = S t a t d i s t (m, gamma ) ;

*/ / I f u s i ng a non s t a t i o n a r y d i s t r i b u t i o n , use t h i s i n s t e a d*

*/ / v ec t o r <Type> d e l t a = Delta w 2 n ( m, t d e l t a ) ;*

*/ / Get number o f t i m e s t e p s ( n )*

**i n t** n = x . s i z e ( ) ;

*/ / Evaluate c o n d i t i o n a l d i s t r i b u t i o n : Put c o n d i t i o n a l*

*/ / p r o b a b i l i t i e s o f observed x i n n t i m e s m matrix*

*/ / ( one column f o r each s t a t e , one row f o r each d a t a p o i n t ) :*

matrix *<*Type*>* e m i s s i o n p r o b s ( n , m) ; matrix *<*Type*>* row 1 vec ( 1 , m) ;

row 1 vec . s e t O n e s ( ) ;

**f o r** ( **i n t** i = 0 ; i *<* n ; i ++) *{*

**i f** ( x [ i ] != x [ i ] ) *{ / / f != f r e t u r n s t r u e i f and only i f f i s NaN .*

*/ / Replace missing v a l u e s ( NA i n R , NaN i n C++) with 1*

e m i s s i o n p r o b s . row ( i ) = row 1 vec ;

*}*

**e l s e** *{*

e m i s s i o n p r o b s . row ( i ) = d p o i s ( x [ i ] , lambda , **f a l s e** ) ;

*}*

*}*

*/ / Corresponds t o t h e book page 333*

matrix *<*Type*>* foo , P ;

Type mllk , sumfoo , l s c a l e ;

**i f** (m == 1 ) *{*

mllk = *−* e m i s s i o n p r o b s . c o l ( 0 ) . a r r a y ( ) . l o g ( ) . sum ( ) ;

*/ / Use ad r e por t on v a r i a b l e s we are i n t e r e s t e d i n :*

ADREPORT( lambda ) ; ADREPORT( gamma ) ; ADREPORT( d e l t a ) ;

Biometrical Journal **52** (2020) 61 35

REPORT( lambda ) ; REPORT( gamma ) ; REPORT( d e l t a ) ;

**return** mllk ;

*}*

foo = ( d e l t a \* v e c t o r *<*Type *>*( e m i s s i o n p r o b s . row ( 0 ) ) ) . m a t r i x ( ) ; sumfoo = foo . sum ( ) ;

l s c a l e = l o g ( sumfoo ) ;

foo . t r a n s p o s e I n P l a c e ( ) ; foo /= sumfoo ;

**f o r** ( **i n t** i = 2 ; i *<*= n ; i ++) *{*

P = e m i s s i o n p r o b s . row ( i *−* 1 ) ;

foo = ( ( foo \* gamma ) . a r r a y ( ) \* P . a r r a y ( ) ) . m a t r i x ( ) ; sumfoo = foo . sum ( ) ;

l s c a l e += l o g ( sumfoo ) ; foo /= sumfoo ;

*}*

mllk = *−* l s c a l e ;

*/ / Use ad r e po r t on v a r i a b l e s f o r which we want s t andard e r r o r s*

ADREPORT( lambda ) ; ADREPORT( gamma ) ; ADREPORT( d e l t a ) ;

*/ / V a r i a b l e s we need f o r l o c a l decoding and conveniency*

REPORT( lambda ) ; REPORT( gamma ) ; REPORT( d e l t a ) ; REPORT( n ) ;

REPORT( e m i s s i o n p r o b s ) ; REPORT( mllk ) ;

**return** mllk ;

*}*

1. This optional function is in the file utils.cpp. It is only necessary if a stationary distribution is not assumed, and ***δ*** is used as a parameter instead of being derived from the transistion probability matrix **Γ**. It codes the function to convert the working parameter tdelta into its natural format.

*/ / Function t r a n s f o r m i n g working parameters i n i n i t i a l d i s t r i b u t i o n*

*/ / t o n a t u r a l parameters*

**template** *<***c l a s s** Type*>*

v e c t o r *<*Type*>* Delta w 2 n ( **i n t** m, v e c t o r *<*Type*>* t d e l t a ) *{*

v e c t o r *<*Type*>* d e l t a (m) ;

36 First Author*et al.* and dd: Running title

**i f** (m == 1 )

**return** Type ( 1 ) ;

*/ / s e t f i r s t e l ement t o one .*

*/ / F i l l i n t h e l a s t m − 1 e l em en t s with working parameters*

*/ / and t a k e e x p o n e n t i a l*

foo *<<* Type ( 1 ) , t d e l t a . exp ( ) ;

*/ / normalize*

d e l t a = foo / foo . sum ( ) ;

**return** d e l t a ;

*}*

1. This function is in the file utils.cpp. It transforms the transition probability matrix **Γ** from its working format to its natural format.

*/ / Function t r a n s f o r m i n g t h e working parameters i n TPM t o*

*/ / n a t u r a l parameters ( w2n )*

**template** *<***c l a s s** Type*>*

matrix *<*Type*>* Gamma w2n ( **i n t** m, v e c t o r *<*Type*>* tgamma ) *{*

*/ / C o n s t r u c t m x m i d e n t i t y matrix*

matrix *<*Type*>* gamma (m, m) ; gamma . s e t I d e n t i t y ( ) ;

**i f** (m == 1 )

**return** gamma ;

*/ / F i l l o f f d i a g o n a l e l e me n t s with working parameters column−wise :*

**i n t** i d x = 0 ;

**f o r** ( **i n t** i = 0 ; i *<* m; i ++)*{*

**f o r** ( **i n t** j = 0 ; j *<* m; j ++)*{*

**i f** ( j != i ) *{*

*/ / F i l l gamma according t o mapping and t a k e e x p o n e n t i a l*

gamma ( j , i ) = tgamma . exp ( ) ( i d x ) ; i d x ++;

*}*

*}*

*}*

*/ / Normalize each row :*

v e c t o r *<*Type*>* cs = gamma . rowwise ( ) . sum ( ) ;

**f o r** ( **i n t** i = 0 ; i *<* m; i ++) gamma . row ( i ) /= cs [ i ] ;

**return** gamma ;

*}*

Biometrical Journal **52** (2020) 61 37

1. This function is in the file utils.cpp. It derives the stationary distribution from the transition probability matrix **Γ**.

*/ / Function computing t h e s t a t i o n a r y d i s t r i b u t i o n o f a Markov chain*

**template** *<***c l a s s** Type*>*

v e c t o r *<*Type*>* S t a t d i s t ( **i n t** m, matrix *<*Type*>* gamma ) *{*

*/ / C o n s t r u c t s t a t i o n a r y d i s t r i b u t i o n*

matrix *<*Type*>* I (m, m) ; matrix *<*Type*>* U(m, m) ;

matrix *<*Type*>* row 1 vec ( 1 , m) ; U = U. s e t O n e s ( ) ;

I = I . s e t I d e n t i t y ( ) ; row 1 vec . s e t O n e s ( ) ;

matrix *<*Type*>* A = I *−* gamma + U;

matrix *<*Type*>*

matrix *<*Type*>*

v e c t o r *<*Type*>*

Ainv = A. i n v e r s e ( ) ;

d e l t a m a t = row 1 vec

\*

Ainv ;

d e l t a = d e l t a m a t . row ( 0 ) ;

**return** d e l t a ;

*}*

1. Functions used in C++ Poisson HMM code
2. Linear model negative log-likelihood calculation
3. Functions used in C++ linear model code
4. simple C acceleration of Zucchini scripts p. 333, A 1.7, A 1.8 with conditional probabilities outside of the forward / backward loop
5. Use same order as in Zucchini
6. the .cpp file with transformation code
7. the .cpp file with likelihood
8. an .R file showcasing the use

38 First Author*et al.* and dd: Running title

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