Smoothing discrete data (II)

– using a hidden Markov model as implemented in the mhsmm package

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We apply the mhsmm package for a simple smoothing task. The data pertain to classification of a cows eating behaviour over time. The true eating status S_t is in the vector cls0 where '1' denotes not eating and '2' denotes eating. The time resolution is one minute. The observed variables x_t in cls1 are actually not observations per se but the result of a classification obtained by a neural network (using nnet() from the nnet package).

See also the vignette "Smoothing discrete data (I)" for an alternative approach to smoothing the data.

```
> load("clsX.RData")
> length(cls0)

[1] 8640
> length(cls1)

[1] 8640
> library(mhsmm)
> plot(cls1[1:2000], type='1', ylim=c(.8,2))
> addStates(cls0[1:2000])
```

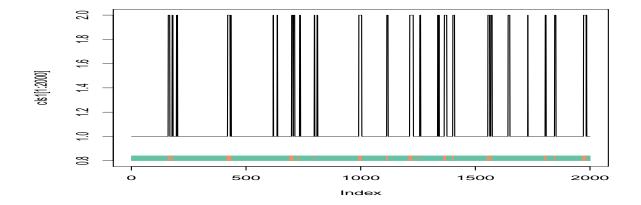


Figure 1: Observed and true eating states

A simple 'smoothing' of the observed states can be obtained as follows: The density function for the emission distribution is

```
> dpmf <- function(x,j,model) model$parms.emission$pmf[j,x]</pre>
An initial setting of the parameters is as follows:
> J <- 2
> init <- c(.5,.5)
> P \leftarrow matrix(c(.9,.1,.1,.9),nrow=J)
> B <- list(pmf=matrix(.1,ncol=J,nrow=J))</pre>
> diag(B$pmf) <- .9
> init.spec <- hmmspec(init,trans=P,parms.emission=B,dens.emission=dpmf)</pre>
> init.spec
Hidden Markov Model specification:
J (number of states):
init:
[1] 0.5 0.5
transition:
     [,1] [,2]
[1,] 0.9 0.1 [2,] 0.1 0.9
emission:
$pmf
[,1] [,2]
[1,] 0.9 0.1
[2,] 0.1 0.9
To fit the model we need to provide the function for the M-step of the EM-algorithm:
> mstep.pmf <- function(x,wt) {</pre>
    ans <- matrix(ncol=ncol(wt),nrow=ncol(wt))</pre>
    for(i in 1:ncol(wt))
      for(j in 1:ncol(wt))
        ans[i,j] \leftarrow sum(wt[which(x==j),i])/sum(wt[,i])
    list(pmf=ans)
For training the model we use the first 1000 cases
> samp <- 1:2640
> train <- list(s=cls0[samp], x=cls1[samp], N=length(cls0[samp]))
> valid <- list(x=cls1[-samp], N=length(cls1[-samp]))</pre>
We fit the model with
> hmm.obj <- hmmfit(train, init.spec,mstep=mstep.pmf)</pre>
> summary(hmm.obj)
init:
 1 0
transition:
      [,1]
            [,2]
[1,] 0.983 0.017
[2,] 0.177 0.823
emission:
$pmf
             [,1]
[1,] 0.999756954 0.0002430464
[2,] 0.008608553 0.9913914471
```

Two types of predictions can be made: Default is to use the Viterbi algorithm for producing the jointly most likely sequence of states given the observed data:

```
> vit <- predict(hmm.obj, valid)</pre>
```

Alternatively we can get the individually most likely state sequence as:

```
> smo <- predict(hmm.obj, valid, method="smoothed")
The prediction results are quite similar:
> normtab <- function(tt) round(sweep(tt,1,rowSums(tt),"/"),2)</pre>
> SS <- cls0[-samp]
> XX <- cls1[-samp]
> cls2.vit <- vit$s
> cls2.smo <- smo$s
> normtab(table(SS,XX))
SS
            2
       1
  1 0.98 0.02
  2 0.27 0.73
> normtab(table(SS,cls2.vit))
   cls2.vit
    1 2
SS
  1 0.98 0.02
  2 0.26 0.74
> normtab(table(SS,cls2.smo))
   cls2.smo
SS
            2
      1
  1 0.98 0.02
  2 0.26 0.74
> show <- 1:2000
> cls0b <- cls0[-samp]
> cls1b <- cls1[-samp]
> c(length(SS),length(cls2.vit),length(cls2.smo), length(cls0b), length(cls1b))
[1] 6000 6000 6000 6000 6000
> plot(cls1b[show], type='1', ylim=c(.8,2))
> addStates(list(cls0b[show],cls2.vit[show], cls2.smo[show]))
```

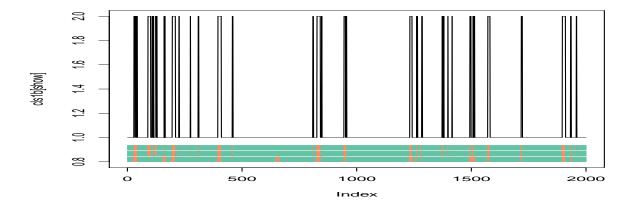


Figure 2: Observed and true eating states