Advanced Machine Learning

Exam Training
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1)

```
library(HMM)
## Hidden variables (true positions)
states <- 1:10
0, 0.5, 0.5, 0, 0, 0, 0, 0, 0, 0,
                          0, 0, 0.5, 0.5, 0, 0, 0, 0, 0, 0,
                          0, 0, 0, 0.5, 0.5, 0, 0, 0, 0, 0,
                          0, 0, 0, 0, 0.5, 0.5, 0, 0, 0, 0,
                          0, 0, 0, 0, 0, 0.5, 0.5, 0, 0, 0,
                          0, 0, 0, 0, 0, 0.5, 0.5, 0, 0,
                          0, 0, 0, 0, 0, 0, 0.5, 0.5, 0,
                          0, 0, 0, 0, 0, 0, 0, 0.5, 0.5,
                          0.5, 0, 0, 0, 0, 0, 0, 0, 0, 0.5),
                        byrow=TRUE, nrow=length(states), ncol=length(states))
## Emission variables (observed positions)
symbols <- 1:10
0.2, 0.2, 0.2, 0.2, 0, 0, 0, 0, 0, 0.2,
                        0.2, 0.2, 0.2, 0.2, 0.2, 0, 0, 0, 0, 0,
                        0, 0.2, 0.2, 0.2, 0.2, 0.2, 0, 0, 0, 0,
                        0, 0, 0.2, 0.2, 0.2, 0.2, 0.2, 0, 0, 0,
                        0, 0, 0, 0.2, 0.2, 0.2, 0.2, 0.2, 0, 0,
                        0, 0, 0, 0, 0.2, 0.2, 0.2, 0.2, 0.2, 0,
                        0, 0, 0, 0, 0.2, 0.2, 0.2, 0.2, 0.2,
                        0.2, 0, 0, 0, 0, 0.2, 0.2, 0.2, 0.2,
                        0.2, 0.2, 0, 0, 0, 0, 0.2, 0.2, 0.2),
                      byrow=TRUE, nrow=length(states), ncol=length(states))
start_probs <- rep(1, length(states)) / length(states)</pre>
robot_hmm <- initHMM(states, symbols,</pre>
                   startProbs=start_probs,
                  transProbs=transition_probs,
                   emissionProbs=emission_probs)
```

```
set.seed(123)
samples_hmm <- simHMM(robot_hmm, 10)</pre>
```

3)

```
compute_filtered_probs <- function(hmm, observations) {</pre>
    log probs <- HMM::forward(hmm, observations)</pre>
    probs <- prop.table(exp(log_probs), 2)</pre>
    probs
}
get_most_probable_states_by_filtered <- function(hmm, observations, states) {</pre>
    probs <- compute_filtered_probs(hmm, observations)</pre>
    most_probable_states <- as.numeric(apply(probs, 2, function(x) {</pre>
        states[which.max(x)]
    }))
    most_probable_states
}
compute_smoothed_probs <- function(hmm, observations) {</pre>
    probs <- HMM::posterior(hmm, observations)</pre>
    probs
}
get_most_probable_states_by_smoothed <- function(hmm, observations, states) {</pre>
    probs <- compute_smoothed_probs(hmm, observations)</pre>
    most_probable_states <- as.numeric(apply(probs, 2, function(x) {</pre>
        states[which.max(x)]
    }))
    most_probable_states
}
get_most_probable_path_by_viterbi <- function(hmm, observations) {</pre>
    most_probable_path <- HMM::viterbi(hmm, observations)</pre>
    most_probable_path
}
get_accuracy_filtered <- function(hmm, samples, states) {</pre>
    predicted_states <- get_most_probable_states_by_filtered(hmm, samples$observation, states)</pre>
    sum(predicted_states == samples$states) / length(predicted_states)
}
get_accuracy_smoothed <- function(hmm, samples, states) {</pre>
    predicted_states <- get_most_probable_states_by_smoothed(hmm, samples$observation, states)</pre>
    sum(predicted_states == samples$states) / length(predicted_states)
}
get_accuracy_viterbi <- function(hmm, samples, states) {</pre>
    predicted_states <- get_most_probable_path_by_viterbi(hmm, samples$observation)</pre>
```

```
sum(predicted_states == samples$states) / length(predicted_states)
}
sample_states <- samples_hmm$states
sample_obs <- samples_hmm$observation

most_probable_states_filtered <- get_most_probable_states_by_filtered(robot_hmm, sample_obs, states)
most_probable_states_smoothed <- get_most_probable_states_by_smoothed(robot_hmm, sample_obs, states)
most_probable_path <- get_most_probable_path_by_viterbi(robot_hmm, sample_obs)</pre>
```

Forward-Backward

```
emission_density <- function(x, z) {</pre>
    return(emission_probs[z, x])
}
transition_density <- function(z, previous_z) {</pre>
    return(transition_probs[previous_z, z])
}
prior_density <- function(z) {</pre>
    return(1 / length(states))
myforward_backward <- function(x, states, prior_density, transition_density, emission_density) {</pre>
    alphas <- myforward(x, states, prior density, transition density, emission density)
    betas <- mybackward(x, states, prior_density, transition_density, emission_density)</pre>
    list(alpha=alphas, beta=betas)
}
myforward <- function(x, states, prior_density, transition_density, emission_density) {</pre>
    T <- length(x)
    alphas <- matrix(NA, ncol=length(states), nrow=T)</pre>
    for (state in states) {
        alphas[1, state] <- emission_density(x[1], state) * prior_density(state)
    }
    for (t in 2:T) {
        inner_sum <- 0
        for (state in states) {
             inner_sum <- 0
             for (previous_state in states) {
                 inner_sum <- inner_sum + alphas[t - 1, previous_state] *</pre>
                     transition_density(state, previous_state)
             alphas[t, state] <- emission_density(x[t], state) * inner_sum</pre>
        }
    }
    alphas
}
mybackward <- function(x, states, prior_density, transition_density, emission_density) {</pre>
    T <- length(x)
    betas <- matrix(NA, ncol=length(states), nrow=T)</pre>
    betas[T, ] <- 1
   for (t in (T - 1):1) {
```

```
for (state in states) {
            inner_sum <- 0
            for (next_state in states) {
                inner_sum <- inner_sum + betas[t + 1, next_state] *</pre>
                    emission_density(x[t + 1], next_state) * transition_density(next_state, state)
            }
           betas[t, state] <- inner_sum</pre>
       }
   }
   betas
alphabeta <- myforward_backward(x=sample_obs,</pre>
                                states = states,
                                prior_density=prior_density,
                                emission_density = emission_density,
                                transition_density = transition_density)
a <- alphabeta$alpha
b <- alphabeta$beta
filtering <- a / rowSums(a)
filtering
#>
         [,1] [,2] [,3]
                                       [,5]
                                                 [,6]
                                                            [,7]
                                                                      [,8]
                             [,4]
#> [1,]
                      0 0.2000000 0.2000000 0.2000000 0.20000000 0.2000000
#> [2,]
                      0 0.1428571 0.2857143 0.2857143 0.28571429 0.0000000
                     0 0.1250000 0.3750000 0.5000000 0.00000000 0.0000000
#> [3,]
#> [4,]
                     0 0.0625000 0.2500000 0.4375000 0.25000000 0.0000000
           0
                 0
         0
#> [5,]
                 0
                    0 0.0000000 0.0000000 0.4230769 0.42307692 0.1538462
#> [6,]
                    0 0.0000000 0.0000000 0.2291667 0.45833333 0.3125000
         0
#> [7,]
          0
                 0
                     0 0.0000000 0.0000000 0.1145833 0.34375000 0.3854167
#> [8,]
                     0 0.0000000 0.0000000 0.0000000 0.24309392 0.3867403
           0
#> [9,]
            0
                     0 0.0000000 0.0000000 0.0000000 0.15714286 0.4071429
#> [10,]
                      0 0.0000000 0.0000000 0.0000000 0.07857143 0.2821429
#>
              [,9]
                        [,10]
#> [1,] 0.0000000 0.00000000
#> [2,] 0.0000000 0.00000000
#> [3,] 0.0000000 0.00000000
#> [4,] 0.0000000 0.00000000
#> [5,] 0.0000000 0.00000000
#> [6,] 0.0000000 0.00000000
#> [7,] 0.1562500 0.00000000
#> [8,] 0.2872928 0.08287293
#> [9,] 0.4357143 0.00000000
#> [10,] 0.4214286 0.21785714
t(compute_filtered_probs(robot_hmm, sample_obs))
#>
        states
#> index 1 2 3
                                           6
```

```
1 0 0 0 0.2000000 0.2000000 0.2000000 0.20000000 0.20000000 0.0000000
     2 0 0 0 0.1428571 0.2857143 0.2857143 0.28571429 0.0000000 0.0000000
#>
     3 0 0 0 0 0.1250000 0.3750000 0.5000000 0.00000000 0.0000000 0.0000000
     4 0 0 0 0 0.0625000 0.2500000 0.4375000 0.25000000 0.0000000 0.0000000
#>
     5 0 0 0 0 0.0000000 0.0000000 0.4230769 0.42307692 0.1538462 0.0000000
#>
#>
     6 0 0 0 0.0000000 0.0000000 0.2291667 0.45833333 0.3125000 0.0000000
#>
     7 0 0 0 0.0000000 0.0000000 0.1145833 0.34375000 0.3854167 0.1562500
     8 0 0 0 0.0000000 0.0000000 0.0000000 0.24309392 0.3867403 0.2872928
#>
     9 0 0 0 0 0.0000000 0.0000000 0.0000000 0.15714286 0.4071429 0.4357143
#>
     10 0 0 0 0.0000000 0.0000000 0.0000000 0.07857143 0.2821429 0.4214286
#>
       states
#>
                10
     1 0.00000000
#>
#>
     2 0.00000000
     3 0.00000000
#>
     4 0.00000000
#>
#>
     5 0.00000000
     6 0.00000000
     7 0.00000000
#>
#>
     8 0.08287293
#>
     9 0.00000000
     10 0.21785714
smoothing <- a * b / rowSums(a * b)</pre>
smoothing
        [,1] [,2] [,3]
                             [,4]
                                      [, 5]
                                                 [,6]
                                                            [,7]
   [1,]
                0 0.45000000 0.4107143 0.13928571 0.00000000 0.00000000
#> [2,]
                     0 0.17857143 0.5428571 0.27857143 0.00000000 0.00000000
#> [3,]
         0
                   0 0.04642857 0.3964286 0.55714286 0.00000000 0.00000000
                0
#> [4,] 0 0 0.00000000 0.1857143 0.60000000 0.21428571 0.00000000
#> [5,] 0 0 0.00000000 0.0000000 0.51071429 0.43214286 0.05714286
#> [6,] 0 0 0.00000000 0.0000000 0.23571429 0.55000000 0.21428571
         0 0
#> [7,]
                   0 0.00000000 0.0000000 0.07857143 0.47142857 0.39642857
         0 0 0 0.00000000 0.0000000 0.00000000 0.31428571 0.50000000
#> [8,]
#> [9,]
                   0 0.00000000 0.0000000 0.00000000 0.15714286 0.40714286
                     0 0.00000000 0.0000000 0.00000000 0.07857143 0.28214286
#> [10,]
           0 0
#>
              [,9]
                       [,10]
#> [1,] 0.00000000 0.0000000
#> [2,] 0.00000000 0.0000000
#> [3,] 0.00000000 0.0000000
#> [4,] 0.00000000 0.0000000
#> [5,] 0.00000000 0.0000000
#> [6,] 0.00000000 0.0000000
#> [7,] 0.05357143 0.0000000
#> [8,] 0.18571429 0.0000000
#> [9,] 0.43571429 0.0000000
#> [10,] 0.42142857 0.2178571
t(compute_smoothed_probs(robot_hmm, sample_obs))
       states
#> index 1 2 3
                                5
                                           6
   1 0 0 0 0.45000000 0.4107143 0.13928571 0.00000000 0.00000000
     2 0 0 0 0.17857143 0.5428571 0.27857143 0.00000000 0.00000000
```

```
3 0 0 0 0.04642857 0.3964286 0.55714286 0.00000000 0.00000000
    4 0 0 0 0.00000000 0.1857143 0.60000000 0.21428571 0.00000000
#>
   5 0 0 0 0.00000000 0.0000000 0.51071429 0.43214286 0.05714286
#> 6 0 0 0 0.00000000 0.0000000 0.23571429 0.55000000 0.21428571
#> 7 0 0 0 0.00000000 0.0000000 0.07857143 0.47142857 0.39642857
   8 0 0 0 0.00000000 0.0000000 0.00000000 0.31428571 0.50000000
#>
#>
   9 0 0 0 0.00000000 0.0000000 0.00000000 0.15714286 0.40714286
   #>
     states
#> index
   1 0.00000000 0.0000000
#>
#>
   2 0.00000000 0.0000000
#>
   3 0.00000000 0.0000000
   4 0.00000000 0.0000000
#>
#>
   5 0.00000000 0.0000000
#>
   6 0.00000000 0.0000000
   7 0.05357143 0.0000000
#>
#> 8 0.18571429 0.0000000
#> 9 0.43571429 0.0000000
#> 10 0.42142857 0.2178571
```

Viterbi

```
myviterbi <- function(x, states, prior_density, transition_density, emission_density) {</pre>
    T <- length(x)
    weights <- matrix(NA, ncol=length(states), nrow=T)</pre>
    weights[1, ] <- log(prior_density(states)) + log(emission_density(x[1], states))</pre>
    paths <- matrix(NA, ncol=length(states), nrow=T)</pre>
    for (t in 1:(T-1)) {
        for (state in states) {
            weights[t + 1, state] <- log(emission_density(x[t + 1], state)) +</pre>
                 max(log(transition_density(state, states)) + weights[t, states])
            paths[t + 1, state] <- which.max(log(transition_density(state, states)) +</pre>
                                               weights[t, states])
        }
    }
    path <- rep(NA, T)
    path[T] <- which.max(weights[T,])</pre>
    for (t in (T - 1):1) {
        path[t] <- paths[t + 1, path[t + 1]]</pre>
    }
    path
}
myviterbi(x=sample_obs,
          states=states,
          prior_density=prior_density,
          emission_density = emission_density,
          transition_density = transition_density)
#> [1] 4 4 4 5 6 6 6 7 7 7
HMM::viterbi(robot_hmm, sample_obs)
#> [1] 4 4 4 5 6 6 6 7 7 7
```