

# Hidden Markov models

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# Part of speech tagging

Given input:

this is a simple sentence

the goal is to identify the part of speech (syntactic category) for each word:

this/DET is/VERB a/DET simple/ADJ sentence/NOUN

I use this task as a running example to introduce HMMs.

The set of part of speech (POS) categories can differ based on the application, corpus annotators and language.

One universal tag set used by Google (Petrov et al., 2011):

| Tag  | Description                                  | Example                  |
|------|--|--------------------------|
| VERB | Verbs (all tenses and modes)                 | eat, ate, eats           |
| NOUN | Nouns (common and proper)                    | home, Micah              |
| PRON | Pronouns                                     | I, you, your, he         |
| ADJ  | Adjectives                                   | yellow, bigger, wildest  |
| ADV  | Adverbs                                      | quickly, faster, fastest |
| ADP  | Adpositions (prepositions and postpositions) | of, in, by, under        |
| CONJ | Conjunctions                                 | and, or, but             |
| DET  | Determiners                                  | a, an, the, this         |
| NUM  | Cardinal numbers                             | one, two, first, second  |
| PRT  | Particles, other function words              | up, down, on, off        |
| X    | Other: foreign words, typos, abbreviations   | brasserie, abcense, HMM  |
| .    | Punctuation                                  | ?, !, .                  |

## POS tagging is hard

Ambiguity:

|                     |    |                       |
|---------------------|----|-----------------------|
| glass of water/NOUN | vs | water/VERB the plants |
| lie/VERB down       | vs | tell a lie/NOUN       |
| wind/VERB down      | vs | a mighty wind/NOUN    |

Sparsity:

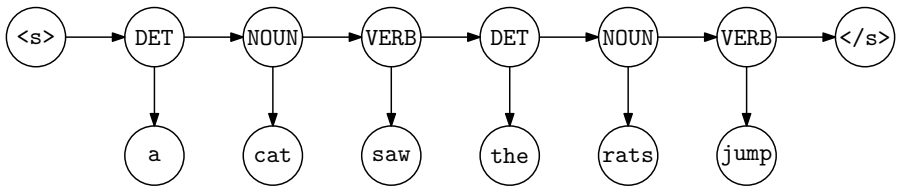
- Words we never see
- Word-tag pairs we never see

## A probabilistic model for tagging

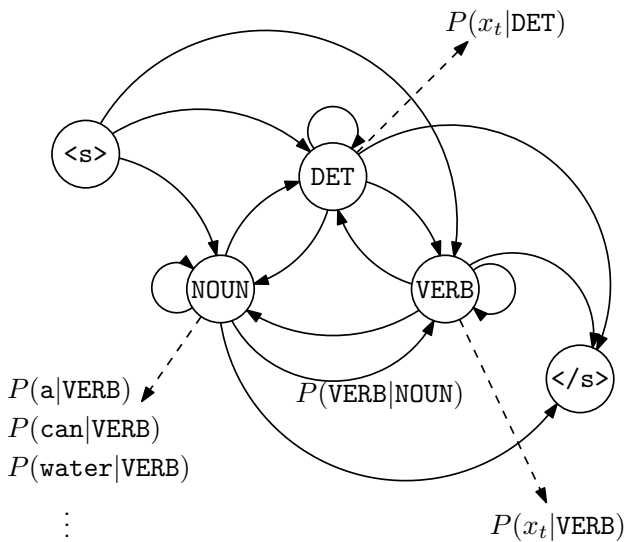
Let  $x_t$  denote the word and  $z_t$  denote the tag at time step  $t$ .

- Initialization:  $z_0 = \langle s \rangle$
- Repeat:
  - Choose a tag based on the previous tag:  $P(z_t | z_{t-1})$
  - If  $z_t = \langle /s \rangle$ : Break
  - Choose a word conditioned on its tag:  $P(x_t | z_t)$

If we knew all the transition probabilities  $P(z_t | z_{t-1})$  and all the emission probabilities  $P(x_t | z_t)$ , we could generate a sentence:



Could also represent this with a state diagram:

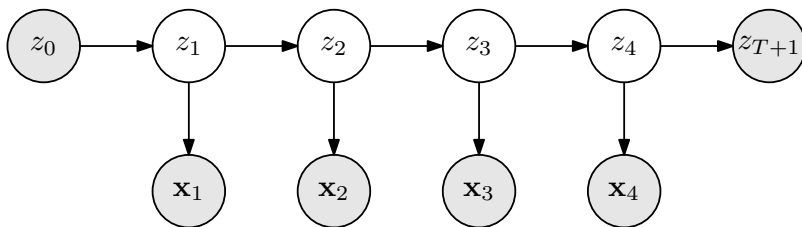


# HMM definition and notation

A hidden Markov model (HMM) defines a probability distribution over a sequence of states and output observations:

- Output sequence:  $\mathbf{x}_{1:T} = \mathbf{x}_1, \mathbf{x}_2, \dots, \mathbf{x}_T$ . We denote these as vectors, but they can also be scalars or discrete observations.
- State sequence:  $z_{0:T+1} = z_0, z_1, \dots, z_T, z_{T+1}$ . The integer variable  $z_t \in \{0, \dots, K + 1\}$  represents the state at  $t$ .

Often  $\mathbf{x}_{1:T}$  is observed and  $z_{0:T+1}$  is hidden:



This is a directed graphical model representation. Contrast this with the state diagram drawn earlier.

Why “hidden”? Why “Markov”?

An HMM is specified by:

- A set of states:  $\{0, 1, \dots, K + 1\}$
- Transition probabilities:  $\mathbf{A}$  with  $A_{i,j} = P_{\mathbf{A}}(z_t = j | z_{t-1} = i)$
- Emission distribution for each state:  $p_{\phi}(\mathbf{x}_t | z_t)$ . We denote the emission distribution as continuous, but can also be discrete.
- Group the parameters together:  $\theta = \{\mathbf{A}, \phi\}$

The start and end states are special:<sup>1</sup>

- Always start in  $z_0 = 0$ .  
Transitioning out of this start state is captured by  $A_{0,j}$ .
- Always end in  $z_{T+1} = K + 1$ .  
Transitioning into this final state is captured by  $A_{j,K+1}$ .
- States 0 and  $K + 1$  are non-emitting: They don't have a corresponding  $\mathbf{x}$  when we move in or out of them.

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<sup>1</sup>The convention in HMM definitions is to explicitly define separate start and end probabilities, but we don't need to do this if we use these non-emitting states.

# The three HMM problems

## Problem 1: The marginal probability

Given an observed sequence  $\mathbf{x}_{1:T}$  and a trained HMM with parameters  $\theta$ , what is the probability of the observed sequence  $p_{\theta}(\mathbf{x}_{1:T})$ ?

## Problem 2: The most likely state sequence

Given an observed sequence  $\mathbf{x}_{1:T}$  and a trained HMM with parameters  $\theta$ , what is the most likely state sequence through the HMM?

$$\arg \max_{z_{0:T+1}} P_{\theta}(z_{0:T+1} | \mathbf{x}_{1:T})$$

## Problem 3: Learning

Given training data  $\mathbf{x}_{1:T}$ , how do we choose the HMM parameters  $\theta$  to maximize  $p_{\theta}(\mathbf{x}_{1:T})$ ?<sup>2</sup>

This is the conventional ordering of the problems (Rabiner and Juang, 1993), but we will do it slightly differently by starting with problem 2.

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<sup>2</sup>In this note I will sometimes be a bit sloppy and drop the dependence of the distribution on  $\theta$  (or  $\mathbf{A}$  or  $\phi$ ), but just remember that they are there.



## 2. Most likely state sequence: Viterbi

We want to find the most likely state sequence given the observed sequence:

$$\begin{aligned}\arg \max_{z_{0:T+1}} P(z_{0:T+1} | \mathbf{x}_{1:T}) &= \arg \max_{z_{0:T+1}} \frac{p(\mathbf{x}_{1:T}, z_{0:T+1})}{p(\mathbf{x}_{1:T})} \\ &= \arg \max_{z_{0:T+1}} p(\mathbf{x}_{1:T}, z_{0:T+1})\end{aligned}$$

We could just calculate  $p(\mathbf{x}_{1:T}, z_{0:T+1})$  for every possible state sequence  $z_{0:T+1}$ . But there are  $K^T$  possible sequences. Instead we use dynamic programming. (Of course.)

Define

$$\delta_t(j) \triangleq \max_{z_{0:t-1}} p(\mathbf{x}_{1:t}, z_{0:t-1}, z_t = j)$$

These variables can be calculated recursively:

$$\delta_t(j) = \max_{z_{0:t-1}} p(\mathbf{x}_{1:t}, z_{0:t-1}, z_t = j)$$

$$= \max_{i=1}^K p(\mathbf{x}_t | z_t = j) A_{i,j} \delta_{t-1}(i)$$

In a similar way you can show that

$$\hat{P} = \max_{z_{0:T+1}} p(\mathbf{x}_{1:T}, z_{0:T+1}) = \max_{j=1}^K A_{j,K+1} \delta_T(j)$$

To find the best state sequence, we start with this last equation and backtrack our steps:

$$\hat{z}_T = \arg \max_{j=1}^K A_{j,K+1} \delta_T(j)$$

The next state is determined by looking at the  $\arg \max$  that gave us  $\delta_T(\hat{z}_T)$ , i.e.

$$\hat{z}_{T-1} = \arg \max_{j=1}^K p(\mathbf{x}_T | z_T = \hat{z}_T) A_{j,\hat{z}_T} \delta_{T-1}(j)$$

and so on.

When calculating  $\delta_t(j)$ , we normally just store the backtrace (the  $\arg \max$ ) denoted as  $\psi_t(j)$  to use when we backtrack.

## Viterbi algorithm

- Initialization:

$$\begin{aligned}\delta_1(j) &= p(\mathbf{x}_1 | z_1 = j) A_{0,j} \\ &\text{for } j = 1, 2, \dots, K\end{aligned}$$

- Recursion:

$$\begin{aligned}\delta_t(j) &= \max_{i=1}^K p(\mathbf{x}_t | z_t = j) A_{i,j} \delta_{t-1}(i) \\ \psi_t(j) &= \arg \max_{i=1}^K p(\mathbf{x}_t | z_t = j) A_{i,j} \delta_{t-1}(i) \\ &\text{for } t = 2, 3, \dots, T \text{ and } j = 1, 2, \dots, K\end{aligned}$$

- Termination:

$$\begin{aligned}\hat{P} &= \max_{j=1}^K A_{j,K+1} \delta_T(j) \\ \hat{z}_T &= \arg \max_{j=1}^K A_{j,K+1} \delta_T(j)\end{aligned}$$

- Backtracking:

$$\begin{aligned}\hat{z}_t &= \arg \max_{j=1}^K p(\mathbf{x}_{t+1} | z_{t+1} = \hat{z}_{t+1}) A_{j,\hat{z}_{t+1}} \delta_t(j) \\ &= \psi_{t+1}(\hat{z}_{t+1}) \\ &\text{for } t = T-1, T-2, \dots, 1\end{aligned}$$

## Example: Viterbi

Transition and emission probabilities:

$$\mathbf{A} = \begin{array}{c} \begin{array}{ccc} \text{c} & \text{v} & \text{</s>} \\ \begin{bmatrix} 1.0 & 0.0 & 0.0 \\ 0.2 & 0.4 & 0.4 \\ 0.7 & 0.1 & 0.2 \end{bmatrix} & \begin{array}{l} \text{<s>} \\ \text{c} \\ \text{v} \end{array} \end{array} \end{array}$$

$$\phi = \begin{array}{c} \begin{array}{ccc} \text{m} & \text{h} & \text{o} \\ \begin{bmatrix} 0.6 & 0.2 & 0.2 \\ 0.1 & 0.3 & 0.6 \end{bmatrix} & \begin{array}{l} \text{c} \\ \text{v} \end{array} \end{array} \end{array}$$

Observed sequence:  $x_{1:3} = \text{m o h}$

Calculate the Viterbi variables  $\delta_t(j)$ :

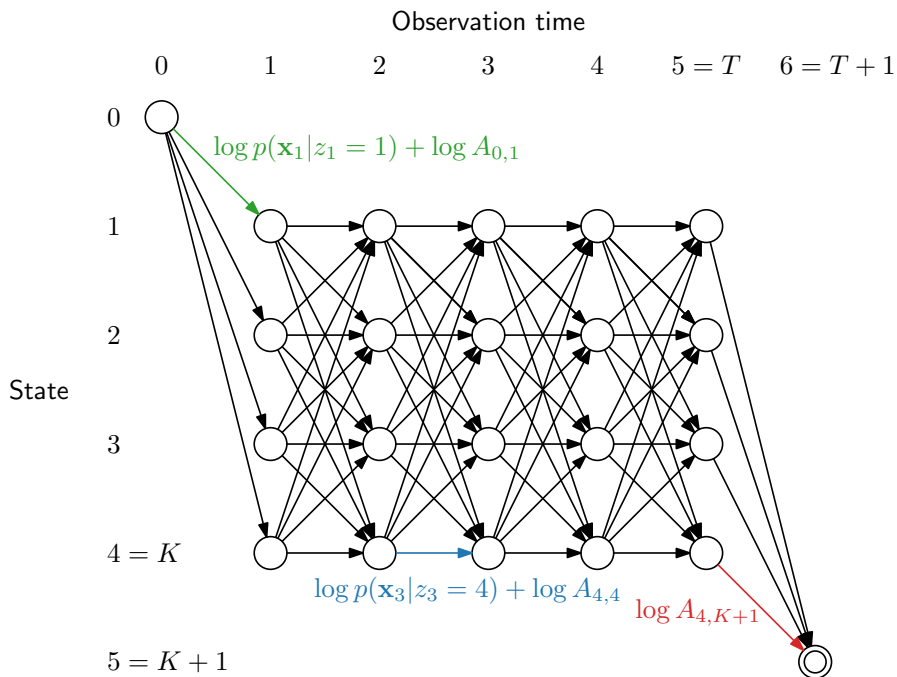
|          |  | $x_1 = \mathfrak{m}$ | $x_2 = \mathfrak{o}$ | $x_3 = \mathfrak{h}$ |  |
|----------|--|----------------------|----------------------|----------------------|--|
| 0 : <s>  |  |                      |                      |                      |  |
| 1 : C    |  |                      |                      |                      |  |
| 2 : V    |  |                      |                      |                      |  |
| 3 : </s> |  |                      |                      |                      |  |

Calculated Viterbi variables  $\delta_t(j)$  with backtracking:

|          |   | $x_1 = \mathfrak{m}$ | $x_2 = \mathfrak{o}$ | $x_3 = \mathfrak{h}$ |          |
|----------|---|----------------------|----------------------|----------------------|----------|
| 0 : <s>  | 1 |                      |                      |                      |          |
| 1 : C    |   | $\nwarrow 0.6$       | $\leftarrow 0.024$   | $\swarrow 0.02016$   |          |
| 2 : V    |   | $\nwarrow 0$         | $\nwarrow 0.144$     | $\leftarrow 0.00432$ |          |
| 3 : </s> |   |                      |                      |                      | 0.008064 |

Does this example seem familiar?

## Just finding the shortest path through a DAG



See my separate note on [dynamic programming](#) for more detail.

Time complexity: In the Viterbi recursion, for each node in the graph, we need to consider the max over  $K$  terms. There are roughly  $K \cdot T$  nodes, which gives  $\mathcal{O}(K^2T)$  terms in the max operations. This is much better than the  $\mathcal{O}(K^T)$  possible paths.

# 1. Marginal probability of sequence

We want the probability of a sequence  $p(\mathbf{x}_{1:T})$ . We will also need this for solving problem 3.

Marginalizing we have:

$$p(\mathbf{x}_{1:T}) = \sum_{z_{0:T+1}} p(\mathbf{x}_{1:T}, z_{0:T+1})$$

We are summing over all possible state sequences. One of these correspond exactly to the Viterbi path above. In the summation, this path will be the one with the highest value  $\hat{P}$ .

Again we could in principle enumerate over all possible sequences, but in practice there are too many:  $K^T$ . Again we turn to dynamic programming. (Of course.)

Define forward variables:

$$\alpha_t(j) \triangleq p(\mathbf{x}_{1:t}, z_t = j)$$

These variables can be calculated recursively:

$$\begin{aligned} \alpha_t(j) &= p(\mathbf{x}_{1:t}, z_t = j) \\ &= p(\mathbf{x}_t | \mathbf{x}_{1:t-1}, z_t = j) p(\mathbf{x}_{1:t-1}, z_t = j) \\ &= p(\mathbf{x}_t | z_t = j) \sum_{i=1}^K p(\mathbf{x}_{1:t-1}, z_t = j, z_{t-1} = i) \\ &= p(\mathbf{x}_t | z_t = j) \sum_{i=1}^K P(z_t = j | z_{t-1} = i) p(\mathbf{x}_{1:t-1}, z_{t-1} = i) \\ &= p(\mathbf{x}_t | z_t = j) \sum_{i=1}^K A_{i,j} \alpha_{t-1}(i) \end{aligned}$$

The marginal probability can then be calculated as:

$$p(\mathbf{x}_{1:T}) = \sum_{j=1}^K A_{j,K+1} \alpha_T(j)$$



## Forward algorithm

- Initialization:

$$\alpha_0(j) = \begin{cases} 1 & \text{if } j = 0 \\ 0 & \text{for } j = 1, 2, \dots, K \end{cases}$$
$$\alpha_1(j) = p(\mathbf{x}_1 | z_1 = j) A_{0,j}$$

for  $j = 1, 2, \dots, K$

- Recursion:

$$\alpha_t(j) = p(\mathbf{x}_t | z_t = j) \sum_{i=1}^K A_{i,j} \alpha_{t-1}(i)$$

for  $t = 2, 3, \dots, T$  and  $j = 1, 2, \dots, K$

- Termination:

$$p(\mathbf{x}_{1:T}) = \sum_{j=1}^K A_{j,K+1} \alpha_T(j)$$

Compare this algorithm to the Viterbi algorithm: The max operations have just been changed to sums.

The  $\alpha_0(j)$  initialization is only used later for solving problem 3 and not necessary for calculating  $p(\mathbf{x}_{1:T})$ .

## Example: Forward algorithm

Calculate the forward variables  $\alpha_t(j)$ :

|          |  | $x_1 = \text{m}$ | $x_2 = \text{o}$ | $x_3 = \text{h}$ |  |
|----------|--|------------------|------------------|------------------|--|
| 0 : <s>  |  |                  |                  |                  |  |
| 1 : C    |  |                  |                  |                  |  |
| 2 : V    |  |                  |                  |                  |  |
| 3 : </s> |  |                  |                  |                  |  |

Calculated forward variables  $\alpha_t(j)$  and marginal probability  $P(x_{1:3})$ :

|          |   | $x_1 = \text{m}$ | $x_2 = \text{o}$ | $x_3 = \text{h}$ |          |
|----------|---|------------------|------------------|------------------|----------|
| 0 : <s>  | 1 |                  |                  |                  |          |
| 1 : C    |   | 0.6              | 0.024            | 0.02112          |          |
| 2 : V    |   | 0                | 0.144            | 0.0072           |          |
| 3 : </s> |   |                  |                  |                  | 0.009888 |

## Backward algorithm

We could have alternatively gone in the reverse direction by defining backwards variables:

$$\beta_t(j) \triangleq p(\mathbf{x}_{t+1:T} | z_t = j)$$

which can also be calculated recursively.

- Initialization:

$$\beta_{T+1}(j) = \begin{cases} 1 & \text{for } j = K + 1 \\ 0 & \text{for } j = 1, 2, \dots, K \end{cases}$$
$$\beta_T(j) = A_{j,K+1}$$
$$\text{for } j = 1, 2, \dots, K$$

- Recursion:

$$\beta_t(j) = \sum_{i=1}^K p(\mathbf{x}_{t+1} | z_{t+1} = i) A_{j,i} \beta_{t+1}(i)$$
$$\text{for } t = T - 1, T - 2, \dots, 0 \text{ and } j = 1, 2, \dots, K$$

- Termination:

$$p(\mathbf{x}_{1:T}) = \sum_{j=1}^K A_{0,j} \beta_0(j)$$

Again the  $\beta_{T+1}(j)$  initialization isn't required for  $p(\mathbf{x}_{1:T})$  and only becomes relevant for solving problem 3.

### 3. Learning HMM parameters

Goal: We want to learn the HMM parameters  $\theta = \{\mathbf{A}, \phi\}$  from training sequences  $\mathbf{x}_{1:T}$ .

Chicken and egg problem:

- If we were in a supervised setting with a known label sequence  $z_{0:T+1}$  that goes with training sequence  $\mathbf{x}_{1:T}$ , then we could just use maximum likelihood estimation (MLE), e.g.

$$\hat{A}_{i,j} = \frac{C(z_t = i, z_{t+1} = j)}{C(z_t = i)}$$

For POS tagging, this would be like knowing the tags for the words in your training data.

- If we weren't given the labelled state sequences but we knew the HMM parameters  $\theta$ , then we could find the best state sequence using Viterbi.

But we have neither labels nor the parameters!

#### A sketch of expectation-maximisation (EM)

- Initialization: Choose initial  $\theta^{(0)}$ .
- Recursion: For each iteration  $m$ :
  - **E-step:** Compute expected counts using  $\theta^{(m-1)}$ .
  - **M-step:** Set  $\theta^{(m)}$  by using MLE with the expected counts from the E-step.
- Termination: Repeat until  $\theta$  converges (or some other criterion).

We consider two flavours of EM for HMMs.

## Hard EM: Viterbi re-estimation

If we had the hard assignments of observations  $\mathbf{x}_t$  to states  $z_t$ , then the MLE estimates for the HMM parameters  $\theta$  would be as follows.

- Transition probabilities:

$$\hat{A}_{i,j} = \frac{C(z_t = i, z_{t+1} = j)}{C(z_t = i)}$$

- Emission distributions:

- We collect all the  $\mathbf{x}$ 's assigned to state  $k$  and then set  $\phi_k$  to the MLE.
- If  $p_{\phi_k}(\mathbf{x}_t | z_t = k)$  is a continuous density, this will be the MLE for that density.
- If  $P_{\phi_k}(x_t | z_t = k)$  is a discrete probability mass function for a categorical distribution, then the MLE is

$$\begin{aligned}\hat{\phi}_{k,c} &= P(x_t = c | z_t = k) \\ &= \frac{C(z_t = k, x_t = c)}{C(z_t = k)}\end{aligned}$$

## The hard EM algorithm

- Initialization:
  - For each training sequence  $\mathbf{x}_{1:T}$ : Assign a sensible  $z_{0:T+1}$  (what is sensible?)
  - Estimate  $\theta^{(0)}$  using the MLE equations (previous page).
- Recursion: For each iteration  $m$ :
  - **E-step:** For each training sequence  $\mathbf{x}_{1:T}$ :
    - \* Apply Viterbi with  $\theta^{(m-1)}$  to get the optimal state sequence  $\hat{z}_{0:T+1}$ .
    - \* Calculate:  $\hat{P} = p(\mathbf{x}_{1:T}, \hat{z}_{0:T+1})$
  - Accumulate the scores:  $J^{(m)} = \sum_{n=1}^N \log \hat{P}^{(n)}$   
(where  $n$  is the index of the training sequence)
  - **M-step:** Use all the optimal state sequences  $\hat{z}_{0:T+1}$  for all the training sequences and estimate  $\theta^{(m)}$  using the MLE equations (previous page).
- Termination: Compare score  $J^{(m)}$  to the previous one  $J^{(m-1)}$  and decide whether to stop.

Although this algorithm uses hard assignments, it is still guaranteed to converge to a local optimum. This is because every E- and M-step guarantees that  $J^{(m)} \geq J^{(m-1)}$ .

## Soft EM: Baum-Welch re-estimation

Instance of the forward-backward algorithm. Here we use soft counts instead of hard assignments.

Define the probability of being in state  $j$  at time  $t$ :

$$\gamma_t(j) \triangleq P(z_t = j | \mathbf{x}_{1:T})$$

which can be calculated as

$$\begin{aligned}\gamma_t(j) &= \frac{p(z_t = j, \mathbf{x}_{1:T})}{p(\mathbf{x}_{1:T})} \\ &= \frac{p(\mathbf{x}_{1:t}, \mathbf{x}_{t+1:T}, z_t = j)}{p(\mathbf{x}_{1:T})} \\ &= \frac{\alpha_t(j) \beta_t(j)}{p(\mathbf{x}_{1:T})}\end{aligned}$$

Define the probability of transitioning from state  $i$  at time  $t$  to state  $j$  at  $t + 1$ :

$$\xi_t(i, j) \triangleq P(z_t = i, z_{t+1} = j | \mathbf{x}_{1:T})$$

which can be calculated as

$$\begin{aligned}\xi_t(i, j) &= \frac{p(z_t = i, z_{t+1} = j, \mathbf{x}_{1:T})}{p(\mathbf{x}_{1:T})} \\ &= \frac{p(\mathbf{x}_{1:t}, \mathbf{x}_{t+1}, \mathbf{x}_{t+2:T}, z_t = i, z_{t+1} = j)}{p(\mathbf{x}_{1:T})} \\ &= \frac{\alpha_t(i) A_{i,j} p(\mathbf{x}_{t+1} | z_{t+1} = j) \beta_{t+1}(j)}{p(\mathbf{x}_{1:T})}\end{aligned}$$

The E- and M-steps at iteration  $m$  now change to:

- **E-step:**

- Forward-backward: Calculate all the forward variables  $\alpha_t(j)$  and backward variables  $\beta_t(j)$  using the current  $\theta^{(m-1)} = \{\mathbf{A}^{(m-1)}, \phi^{(m-1)}\}$ .
- Calculate the expected number of transitions out of state  $j$ :

$$\sum_{t=0}^T \gamma_t(j)$$

- Calculate the expected number of transitions from state  $i$  to  $j$ :

$$\sum_{t=0}^T \xi_t(i, j)$$

- **M-step:**

- Transition probabilities:

$$\hat{A}_{i,j} = \frac{\sum_{t=0}^T \xi_t(i, j)}{\sum_{t=0}^T \gamma_t(i)}$$

- Emission probability mass function for categorical distribution:

$$\begin{aligned} \hat{\phi}_{k,c} &= P(x_t = c | z_t = k) \\ &= \frac{\sum_{t=1}^T \mathbb{I}\{x_t = c\} \gamma_t(k)}{\sum_{t=1}^T \gamma_t(k)} \end{aligned}$$



### Side note: Start and end states

We assume that we start in state 0 and end in state  $K + 1$ . We therefore need to be a bit careful with  $\gamma_t(j)$  when  $t = 0$  and with  $\xi_t(i, j)$  when  $t = 0$  or  $t = T$ .

As one example of these edge cases, we need

$$\gamma_0(j) = \begin{cases} 1 & \text{if } j = 0 \\ 0 & \text{otherwise} \end{cases}$$

Fortunately the initializations for  $\alpha_0(j)$  and  $\beta_{T+1}(j)$  given earlier ensure that  $\gamma_t(j)$  and  $\xi_t(i, j)$  are correct at the start and end of a sequence. The one additional requirement is that we don't have a  $p(\mathbf{x}_{T+1}|z_{t+1} = j)$  term in the numerator when calculating  $\xi_T(i, j)$ . (Why not?)

## Example: Soft and hard EM

- States:  $\{0, 1, 2, 3\}$
- Output vocabulary:  $\{m, o, h\}$
- Training sequence:  $x_{1:T} = m \ o \ o \ m \ o \ h \ o \ h \ o$

**Hard EM**

m o o m o h o h o

**Soft EM**

m o o m o h o h o

## Use logs: The log-sum-exp trick

We have seen before that when multiplying many probabilities, we can get underflow. So let's work in the log domain! But just taking the log isn't enough in this case.

Consider the forward variables:

$$\alpha_t(j) = p(\mathbf{x}_t | z_t = j) \sum_{i=1}^K A_{i,j} \alpha_{t-1}(i)$$

We now switch to working with  $\log \alpha_t(j)$  throughout (and similar for the other variables).

$$\begin{aligned} \log \alpha_t(j) &= \\ &= \log p(\mathbf{x}_t | z_t = j) + \log \sum_{i=1}^K e^{\log A_{i,j} + \log \alpha_{t-1}(i)} \end{aligned}$$

But, sadly, the summation for calculating  $\log \alpha_t(j)$  can't happen in the log domain. Even with all the required probabilities in the log domain, we would first have to bring the probabilities back to the linear domain before summing.

## The log-sum-exp trick

In general we actually often end up having to calculate terms like

$$\log \sum_{k=1}^K e^{b_k}$$

where we could potentially get underflow/overflow when calculating  $e^{b_k}$ .

Luckily we have a trick that allows us to calculate the sum! The log-sum-exp trick works as follows:

$$\begin{aligned} \log \sum_{k=1}^K e^{b_k} &= \\ &= B + \log \sum_{k=1}^K e^{b_k - B} \end{aligned}$$

with  $B = \max_{k=1}^K b_k$ .

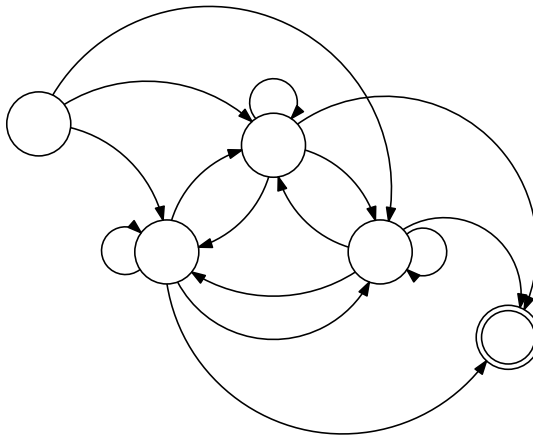
Now at least we don't get underflow/overflow from the biggest term in the summation.

# HMM topologies

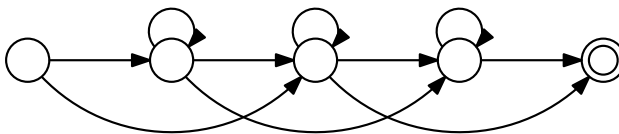
Say  $\hat{A}_{i,j} = 0$ , i.e.  $P_{\hat{A}}(z_t = j | z_{t-1} = i) = 0$ . How will the estimate  $\hat{A}_{i,j}$  change during EM training?

Turning off some transitions at initialization allows us to specify different HMM topologies.

Fully connected topology:



Left-to-right topology:



# Other aspects of HMMs

## Smoothing

Some state transitions or some state-output combinations might not occur in the training data. Similar to what we do with  $N$ -gram language models, we could use smoothing to assign some mass in  $P(x_t|z_t)$  and  $P(z_t|z_{t-1})$  to unseen events.

## Extensions of HMMs

- Higher-order HMMs

## HMMs today

- Although we looked at POS tagging as a running example in this note, HMMs are rarely used for supervised or unsupervised POS tagging today.
- But HMMs are still used a lot in speech recognition:
  - HMM-based speech recognition was the standard for a long time (up to the early 2010s).
  - HMMs are still used today in hybrid HMM-DNN speech recognition systems.



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## References

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