Hidden Markov Models

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HMMs

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April 7th, 2020

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The model consists of a state space Q ≠ Ø (for our purposes Q is finite)

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- and a transition probability matrix p_{ij} where $i,j \in Q$
- The model has no memory, i.e. the probability of moving from state *i* to *j* depends only on the state *i*.
- multiplying the matrix *P*, we can compute the change of the probability distribution as the model "steps" forward
- We are usually interested in stationary distributions π , such that $\pi \cdot P = \pi$

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For any trajectory $\pi,$ we can calculate the probability of emiting ${\cal S}$

$$P(S,\pi) = \prod_{t=0}^{n-1} e_{\pi(t+1)}(S(t+1)) \cdot p_{\pi(t),\pi(t+1)},$$

Can we find the optimal trajectory π , given S?

$$P(S, \pi_*) = max\{P(S, \pi) \mid \pi \in Q^*, |\pi| = |S|\}.$$

Viterbi algorithm

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We can use dynamic programming, filling in the v(i, k) matrix $v(i, k) = max\{P(S[1..i], \pi) \mid \pi \in Q^i, \pi(i) = k\}.$

with the initial condition:

$$v(0,k) = \begin{cases} 1 & \text{gdy } k = k_0, \\ 0 & \text{gdy } k \neq k_0. \end{cases}$$

and step function:

$$v(i,k) = e_k(S(i)) \cdot \max_{l \in Q} [v(i-1,l) \cdot p_{l,k}].$$

To finally read out the seeked probability:

$$P(S, \pi_*) = \max_{k \in Q} [v(|S|, k)].$$

Estimating emission probabilities

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Now, we can calculate the probability of emitting S, over all possible trajectories, with the Forward-method. The initial step is as follows:

$$f(0,k) = \begin{cases} 1 & \text{gdy } k = k_0, \\ 0 & \text{gdy } k \neq k_0. \end{cases}$$

Then, we make similar steps:

$$f(i,k) = e_k(S(i)) \cdot \sum_{l \in Q} f(i-1,l) \cdot p_{l,k}.$$

and finally we can calculate the total probability at the end:

$$P(S) = \sum_{k \in Q} f(|S|, k).$$

The same works backwards:

$$b(i,k) = \sum_{l \in Q} p_{k,l} \cdot e_l(S(i+1)) \cdot b(i+1,l).$$

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Putting it together, the probability of being in state k at step i, given S is equal to:

$$P(\pi(i) = k \mid S) = \frac{P(\pi(i) = k \& S)}{P(S)} = \frac{f(i, k) \cdot b(i, k)}{P(S)}$$

emission probabilities

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Estimate of the Emission matrix:

$$e_k(x) = \frac{E_k(x)}{\sum_{y \in \Sigma} E_k(y)}.$$

Can be calculated using f and b

$$E_k(x) = \sum_{j=1}^n \sum_{i \in I_j(x)} \frac{f_{\mathcal{M}}^{(j)}(i,k) \cdot b_{\mathcal{M}}^{(j)}(i,k)}{P_{\mathcal{M}}(S_j)},$$

Calculating probabilities

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Similarly the transition matrix:

$$p_{k,l} = \frac{P_{k,l}}{\sum_{q \in Q} P_{k,q}},$$

depends on f and b

$$P_{k,l} = \sum_{j=1}^{n} \sum_{i=1}^{|S_j|} \frac{f_{\mathcal{M}}^{(j)}(i,k) \cdot p_{k,l}^{\mathcal{M}} \cdot e_l^{\mathcal{M}}(S_j(i+1)) \cdot b_{\mathcal{M}}^{(j)}(i+1,l)}{P_{\mathcal{M}}(S_j)}$$

Baum-Welch algorithm

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- Suppose, we only know the word S and the sets Q and Σ.
 Can we estimate both p_{ij} and e_{ij}?
- We can start with random p_{ij}, e_{ij} , and iteratively proceed as follows:
 - Calculate the estimates of being in each of states at each step using *f*, *b* and current estimates of *e*, *p*.
 - Find the optimal e, p, given current e, p, f, b
- This is an example of a known procedure called Expectation-Maximization
- It converges to a local optimum of the likelihood, because at every iteration, the likelihood cannot be decreased.

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- We will discuss it in more depth next week, but the HMM model is very useful in describing sequence alignments and so-called sequence profiles
- It is relatively easy to extend this model for arbitrary emissions (e.g. Gaussian or multinomial), not necessarily from a discrete space of symbols. This is frequently used for modelling functional genomics data
- It is also quite a good model for segmentation of chromosomes based on different measurements along the genome

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