

Outline

Hidden Markov models: shortcomings

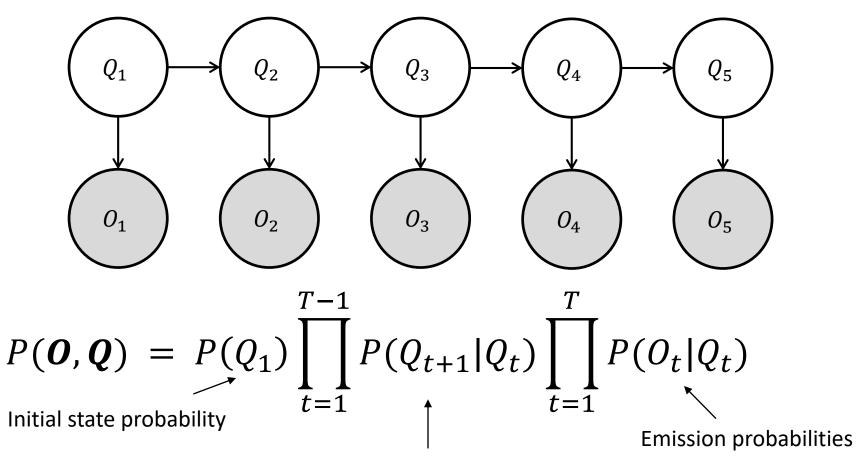
Generative vs. discriminative models

Linear-chain CRFs

Inference and learning algorithms with linear-chain CRFs

Hidden Markov Model

Graph specifies how join probability decomposes



State transition probabilities

Shortcomings of Standard HMMs

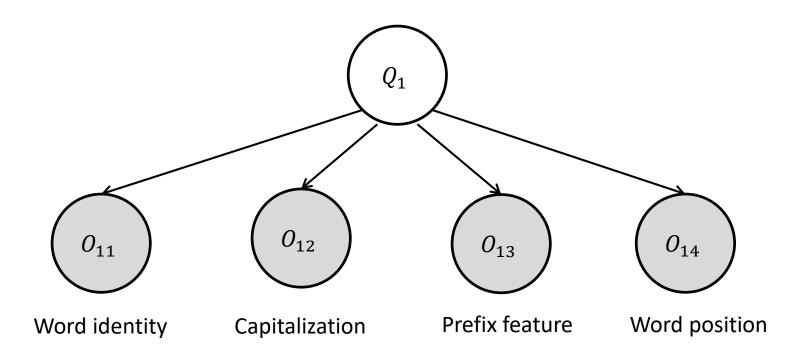
How do we add more features to HMMs?

Might be useful for POS tagging:

- Word position within sentence (1st, 2nd, last...)
- Capitalization
- Word prefix and suffixes (-tion, -ed, -ly, -er, re-, de-)
- Features that depend on more than the current word or the previous words.

Possible to Do with HMMs

Add more emissions at each timestep



Clunky

Is there a better way to do this?

Discriminative Models

HMMs are generative models

- Build a model of the joint probability distribution $P(\mathbf{0}, \mathbf{Q})$,
- Let's rename the variables
- Generative models specify $P(X,Y;\theta^{\text{gen}})$

If we are only interested in POS tagging, we can instead train a **discriminative model**

- Model specifies $P(Y|X; \theta^{\text{disc}})$
- Now a task-specific model for sequence labelling; cannot use it for generating new samples of word and POS sequences

Discriminative Sequence Model

The parallel to an HMM in the discriminative case: linear-chain conditional random fields (linear-chain CRFs) (Lafferty et al., 2001)

$$P(Y|X) = \frac{1}{Z(X)} \exp \sum_{t} \sum_{k} \theta_k f_k(y_t, y_{t-1}, x_t)$$
sum over all features
sum over all time-steps

Z(X) is a normalization constant:

$$Z(X) = \sum_{y} \exp \sum_{t} \sum_{k} \theta_{k} f_{k}(y_{t}, y_{t-1}, x_{t})$$

sum over all possible sequences of hidden states

Intuition

Standard HMM: product of probabilities; these probabilities are defined over the identity of the states and words

- Transition from state DT to NN: $P(y_{t+1} = NN | y_t = DT)$
- Emit word the from state DT: $P(x_t = the | y_t = DT)$

Linear-chain CRF: replace the products by numbers that are NOT probabilities, but linear combinations of weights and feature values.

Features in CRFs

Standard HMM probabilities as CRF features:

Transition from state DT to state NN

$$f_{DT\to NN}(y_t, y_{t-1}, x_t) = \mathbf{1}(y_{t-1} = DT) \mathbf{1}(y_t = NN)$$

Emit the from state DT

$$f_{DT \to the}(y_t, y_{t-1}, x_t) = \mathbf{1}(y_t = DT) \mathbf{1}(x_t = the)$$

Initial state is DT

$$f_{DT}(y_1, x_1) = \mathbf{1}(y_1 = DT)$$

Indicator function:

Let
$$\mathbf{1}(condition) = \begin{cases} 1 & \text{if } condition \text{ is true} \\ 0 & \text{otherwise} \end{cases}$$

Features in CRFs

More examples of features that may be useful

Word is capitalized

$$f_{cap}(y_t, y_{t-1}, x_t) = \mathbf{1}(y_t = ?)\mathbf{1}(x_t \text{ is capitalized})$$

• Word ends in –ed

$$f_{-ed}(y_t, y_{t-1}, x_t) = \mathbf{1}(y_t = ?)\mathbf{1}(x_t \text{ ends with } ed)$$

Inference with LC-CRFs

Dynamic programming still works – modify the forward and the Viterbi algorithms to work with the weight-feature products.

	нмм	LC-CRF
Forward algorithm	$P(X \theta)$	Z(X)
Viterbi algorithm	$\underset{Y}{\operatorname{argmax}} P(X,Y \theta)$	$\operatorname*{argmax}_{Y} P(Y X,\theta)$

Forward Algorithm for HMMs

Create trellis
$$\alpha_i(t)$$
 for $i=1\dots N, t=1\dots T$ $\alpha_j(1)=\pi_j b_j(O_1)$ for $j=1\dots N$ for $t=2\dots T$: for $j=1\dots N$:
$$\alpha_j(t)=\sum_{i=1}^N \alpha_i(t-1)a_{ij}b_j(O_t)$$

$$P(\boldsymbol{O}|\theta)=\sum_{j=1}^N \alpha_j(T)$$

Runtime: $O(N^2T)$

Forward Algorithm for LC-CRFs

Create trellis
$$\alpha_i(t)$$
 for $i=1\dots N, t=1\dots T$
$$\alpha_j(1) = \exp\sum_k \theta_k^{init} f_k^{init} (y_1=j,x_1) \text{ for } j=1\dots N$$
 for $t=2\dots T$: for $j=1\dots N$:
$$\alpha_j(t) = \sum_{i=1}^N \alpha_i(t-1) \exp\sum_k \theta_k f_k (y_t=j,y_{t-1},x_t)$$

$$Z(X) = \sum_{j=1}^N \alpha_j(T)$$
 Transition and emission probabilities replaced by exponent of weighted sums of features.

Runtime: $O(N^2T)$

Having Z(X) allows us to compute P(Y|X)

Viterbi Algorithm for HMMs

```
Create trellis \delta_i(t) for i=1\dots N, t=1\dots T \delta_j(1)=\pi_jb_j(O_1) for j=1\dots N for t=2\dots T: for j=1\dots N: \delta_j(t)=\max_i\delta_i(t-1)a_{ij}b_j(O_t) Take \max_i\delta_i(T)
```

Runtime: $O(N^2T)$

Viterbi Algorithm for LC-CRFs

Create trellis
$$\delta_i(t)$$
 for $i=1\dots N, t=1\dots T$
$$\delta_j(1)=\exp\sum_k \theta_k^{init} f_k^{init} (y_1=j,x_1) \text{for } j=1\dots N$$
 for $t=2\dots T$: for $j=1\dots N$:
$$\delta_j(t)=\max_i \delta_i(t-1)\exp\sum_k \theta_k f_k (y_t=j,y_{t-1},x_t)$$
 Take $\max_i \delta_i(T)$

Runtime: $O(N^2T)$

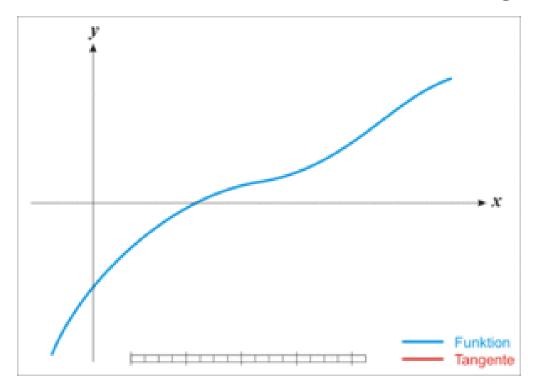
Remember that we need backpointers.

Training LC-CRFs

Unlike for HMMs, no analytic MLE solution
Use iterative method to improve data likelihood

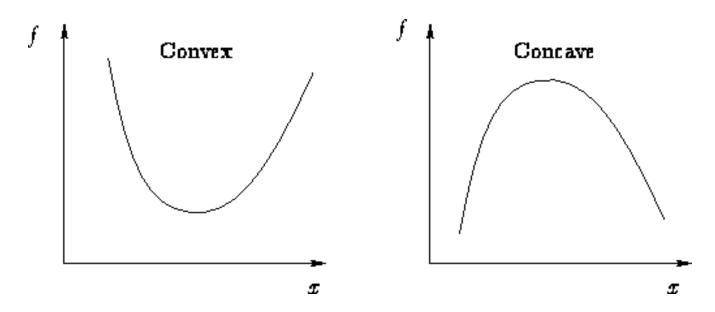
Gradient descent

A version of Newton's method to find where the gradient is 0



Convexity

Fortunately, $l(\theta)$ is a **concave** function (equivalently, its negation is a **convex** function). That means that we will find the global maximum of $l(\theta)$ with gradient ascent (equivalently, the global minimum of $-l(\theta)$ with gradient descent).



Gradient Ascent

Walk in the direction of the gradient to maximize $l(\theta)$

a.k.a., gradient descent on a loss function

$$\theta^{\text{new}} = \theta^{\text{old}} + \gamma \nabla l(\theta)$$

 γ is a learning rate that specifies how large a step to take.

There are more sophisticated ways to do this update:

- Conjugate gradient
- L-BFGS (approximates using second derivative)

Gradient Descent Summary

Descent vs ascent

Convention: think about the problem as a minimization problem

Minimize the negative log likelihood

•
$$\theta \leftarrow \theta - \gamma(-\nabla l(\theta))$$

Initialize $\theta = \{\theta_1, \theta_2, ..., \theta_k\}$ randomly

Do for a while:

Compute $\nabla l(\theta)$, which will require dynamic programming (i.e., forward algorithm)

$$\theta \leftarrow \theta + \gamma \nabla l(\theta)$$

Gradient of Log-Likelihood

Find the gradient of the log likelihood of the training corpus:

$$l(\theta) = \log \prod_{i} P(Y^{(i)}|X^{(i)})$$

Interpretation of Gradient

Overall gradient is the difference between:

$$\sum_{i} \sum_{t} f_{k}(y_{t}^{(i)}, y_{t-1}^{(i)}, x_{t}^{(i)})$$

the empirical distribution of feature f_k in the training corpus and:

$$\sum_{i} \sum_{t} \sum_{y,y'} f_k(y,y',x_t^{(i)}) P(y,y'|X^{(i)})$$

the expected distribution of f_k as predicted by the current model

Interpretation of Gradient

When the corpus likelihood is maximized, the gradient is zero, so the difference is zero.

Intuitively, this means that finding parameter estimate by gradient descent is equivalent to telling our model to predict the features in such a way that they are found in the same distribution as in the gold standard.

Regularization

To avoid overfitting, we can encourage the weights to be close to zero.

Add term to corpus log likelihood:

$$l^*(\theta) = \log \prod_{i} P(Y^{(i)}|X^{(i)}) - \sum_{k} \frac{\theta_k^2}{2\sigma^2}$$

 σ controls the amount of **regularization**

Results in extra term in gradient:

$$-\frac{\theta_k}{\sigma^2}$$

Stochastic Gradient Descent (SGD)

In the standard version of the algorithm, the gradient is computed over the entire training corpus.

Weight update only once per iteration through training corpus.

Alternative: calculate gradient over a small mini-batch of the training corpus and update weights

SGD is when mini-batch size is one.

- Many weight updates per iteration through training corpus
- Usually results in much faster convergence to final solution, without loss in performance

Stochastic Gradient Descent

Goal: Minimize $-l^*(\theta)$

Initialize $\theta = \{\theta_1, \theta_2, ..., \theta_k\}$ randomly

Do for a while:

Randomize order of samples in training corpus

For each mini-batch (of size one) in the training corpus:

Compute $\nabla l^*(\theta)$ over this mini-batch $\theta \leftarrow \theta + \gamma \nabla l^*(\theta)$