

Overview of Natural Language Processing

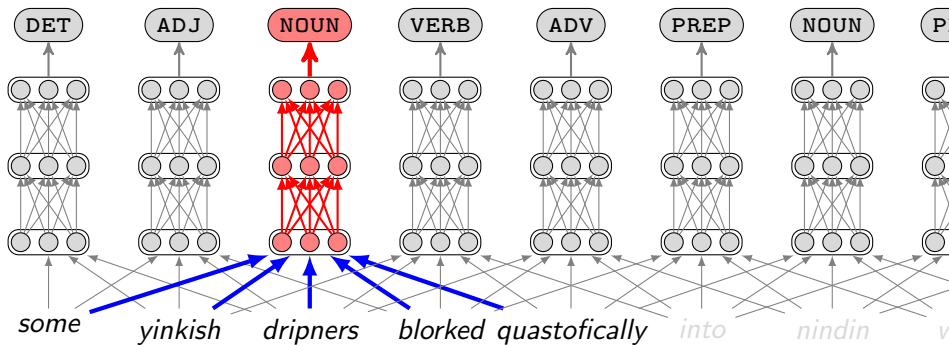
Part II & ACS L390

Gradient Descent and Neural Nets

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Gradient Descent and Neural Nets

1. Gradient Descent
2. Feedforward Neural Networks

Gradient Descent

Supervised learning

Assume there is a *good* annotated corpus

$$\mathcal{D} = \left\{ (x^{(1)}, y^{(1)}), (x^{(2)}, y^{(2)}), \dots, (x^{(l)}, y^{(l)}) \right\}$$

How can we get a *good* parameter vector?

Maximum-Likelihood Estimation

$$\hat{\theta} = \text{arg max } L(\theta)$$

where $L(\theta)$ is the log-likelihood of observing the data \mathcal{D} :

$$L(\theta) = \sum_{i=1}^l \log p(y^{(i)} | x^{(i)}; \theta)$$

Gradient Descent/Ascent

In general, finding a minimum/maximum is *hard*.

However, a simple idea that often works:

- Initialise θ with some value
- Iteratively improve θ

The derivative tells us whether to increase or decrease
(but doesn't tell us how much to increase/decrease by):

$$\theta^{[t+1]} = \theta^{[t]} + \beta \frac{dL}{d\theta}(\theta^{[t]})$$

Gradient Descent for the Log-Linear Model

Assume we have a *parameter vector* θ .

$$p(y|x; \theta) = \frac{\exp(\theta^\top f(x, y))}{\sum_{y' \in \mathcal{Y}} \exp(\theta^\top f(x, y'))}$$

i.e.

$$p(y|x; \theta) \propto \exp(\theta^\top f(x, y))$$

$$\begin{aligned} L(\theta) &= \sum_{i=1}^l \log p(y^{(i)}|x^{(i)}; \theta) \\ &= \sum_{i=1}^l \left(\theta^\top f(x^{(i)}, y^{(i)}) - \log \sum_{y' \in \mathcal{Y}} \exp(\theta^\top f(x^{(i)}, y')) \right) \end{aligned}$$

Gradient Descent for the Log-Linear Model

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Calculating gradients (chain rule)

$$\begin{aligned} \frac{dL}{d\theta_k} &= \sum_{i=1}^l \left(f_k(x^{(i)}, y^{(i)}) - \frac{\sum_{y' \in \mathcal{Y}} (\exp(\theta^\top f(x^{(i)}, y')) f_k(x^{(i)}, y'))}{\sum_{y^* \in \mathcal{Y}} \exp(\theta^\top f(x^{(i)}, y^*))} \right) \\ &= \sum_{i=1}^l f_k(x^{(i)}, y^{(i)}) - \sum_{i=1}^l \sum_{y' \in \mathcal{Y}} f_k(x^{(i)}, y') \frac{\exp(\theta^\top f(x^{(i)}, y'))}{\sum_{y^* \in \mathcal{Y}} \exp(\theta^\top f(x^{(i)}, y^*))} \end{aligned}$$

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Gradient Descent: Algorithm

Maximize $L(\theta)$ where

$$\frac{dL}{d\theta_k} = \underbrace{\sum_{i=1}^l f_k(x^{(i)}, y^{(i)})}_{\text{empirical counts}} - \underbrace{\sum_{i=1}^l \sum_{y' \in \mathcal{Y}} f_k(x^{(i)}, y') p(y'|x^{(i)}; \theta)}_{\text{expected counts}}$$

1 **Initialize** $\theta^{[0]} \leftarrow 0$

2 **for** $t = 1, \dots$

3 **calculate** $\Delta = \frac{dL(\theta^{[t-1]})}{d\theta}$

4 **calculate** $\beta_* = \arg \max_{\beta} L(\theta + \beta \Delta)$

▷ line search

5 **update** $\theta^{[t]} \leftarrow \theta^{[t-1]} + \beta_* \Delta$

Gradient Descent: Algorithm

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Challenges

- Go through every training sample to get Δ ;
- Line search is another non-trivial optimisation problem

Stochastic Gradient Descent

$$L(\theta) = \sum_{i=1}^l \log p(y^{(i)}|x^{(i)}; \theta)$$

- Randomly use one or several training samples to get a suboptimal gradient Δ' ;
- Fix β .

Feedforward Neural Networks

Recap: about linear combination

$$p(y|x; \theta) = \frac{\exp(\theta^\top f(x, y))}{\sum_{y' \in \mathcal{Y}} \exp(\theta^\top f(x, y'))}$$

... 0 0 1 0 1 0 0 1 0 1 0 0 0 0 0 0 ...

f_{1001} : if word₋₂=some and tag=N

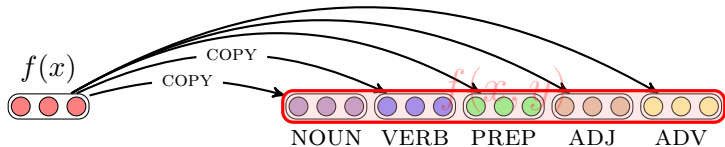
is θ_{1001} positive large?
vote for yes

$\theta^\top f(x, y)$ is a linear combination of θ . That is why a log-linear model is called a linear classifier.

Questions

Can we automate the design of features?

Is linear combination justified?

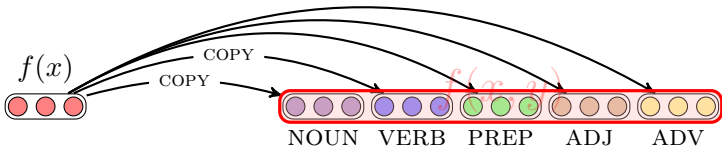


A simple way to define $f(x, y)$ based on $f(x)$ is “copy”. We assumed that θ and $f(x, y)$ have DK dimensions:

- D – number of input features
- K – number of output classes

So we can also view θ as comprising K vectors with D dimensions:

$$p(y|x; \theta) \propto \exp(\theta_y^\top f(x))$$

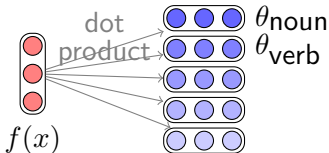


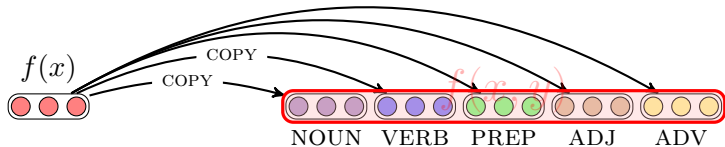
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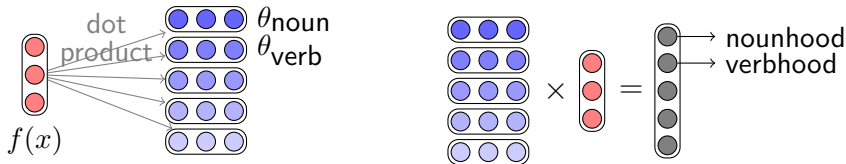


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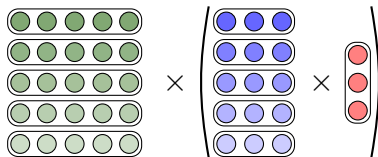
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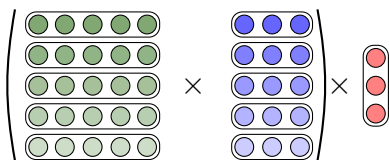
$$p(y|x; \theta) \propto \exp(\theta_y^\top f(x))$$



Now consider NER after POS tagging

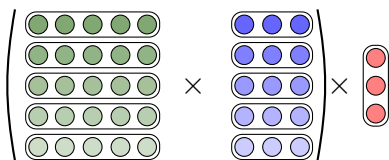


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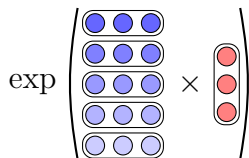
Oops! The combined transformation is linear!

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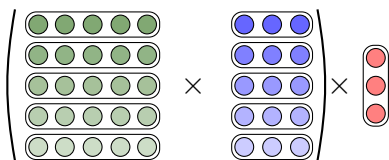
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Introduce some non-linearity



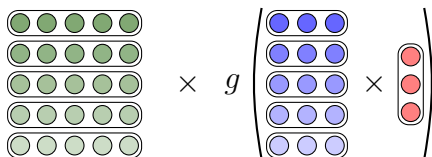
\exp / g is applied component-wise (to each dimension separately)

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Feedforward Neural Networks

Think about multi-class classification:

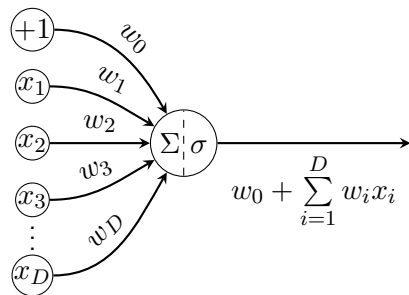
- D – number of features (input)
- K – number of classes (output)
- \mathbf{x} – the input feature vector

Think about a particular class, say y_k . We describe the “friendship” between \mathbf{x} and y_k in the following way:

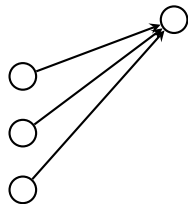
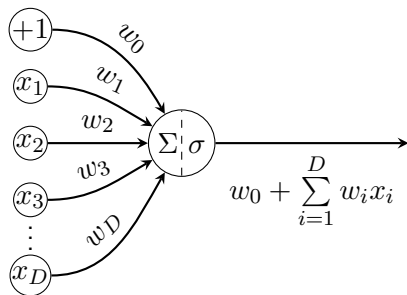
$$\text{score_function}(x, y_k) = w_0 + \sum_{i=1}^D w_i x_i$$

where \mathbf{w} measures how much each feature w_i contributes to y_k .

Feedforward Neural Networks

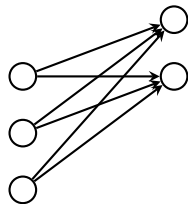
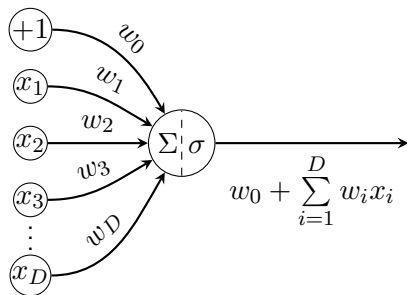


Feedforward Neural Networks



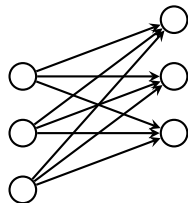
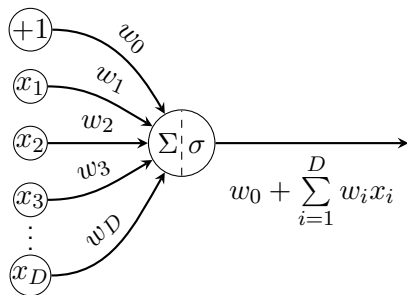
For each class y_k , we do the same thing.

Feedforward Neural Networks



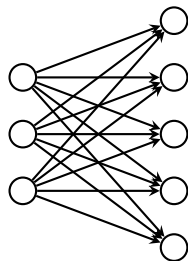
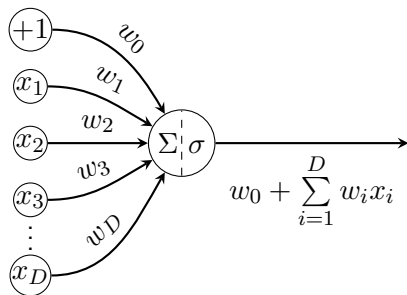
For each class y_k , we do the same thing. Again

Feedforward Neural Networks



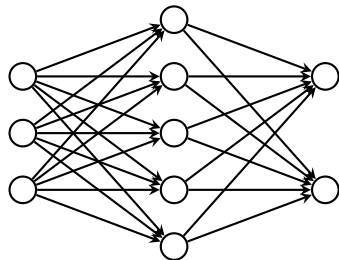
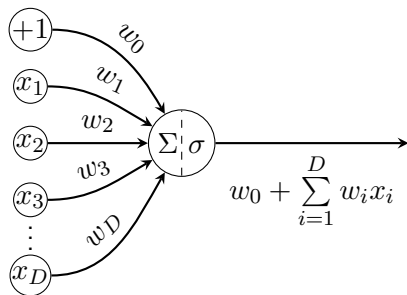
For each class y_k , we do the same thing. Again and again

Feedforward Neural Networks



For each class y_k , we do the same thing. Again and again and again. This is called **perceptron**, which was invented by Frank Rosenblatt in 1958.

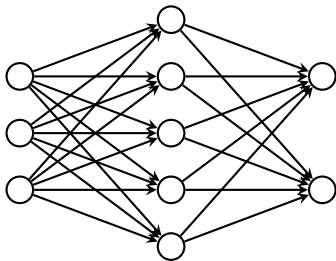
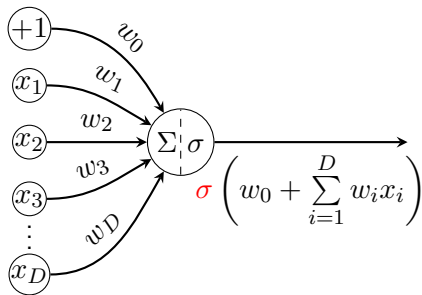
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Things will be much more fun if we have a stack of perceptrons.

Feedforward Neural Networks



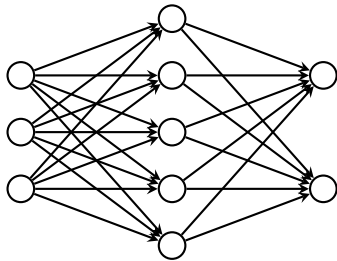
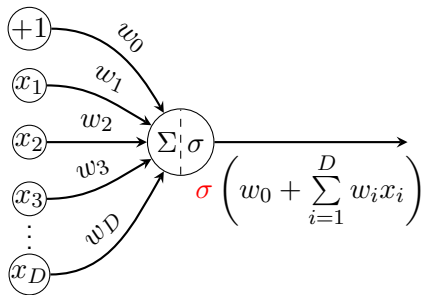
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Things will be much more fun if we have a stack of perceptrons. **Oops, must add something...**

Sigmoid $\sigma(x) = \frac{1}{1+e^{-x}}$

Otherwise, simple matrix multiplication.

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Sigmoid $\sigma(x) = \frac{1}{1+e^{-x}}$

Otherwise, simple matrix multiplication. Now you can do non-linear classification.

Feedforward Neural Networks

A feedforward neural network is a composition of simple functions:

- 1 layer: $\exp(W_1x)$ ▷ log-linear model
- 2 layers: $\exp(W_2g(W_1x))$
- 3 layers: $\exp(W_3g(W_2g(W_1x)))$
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Both \exp and g are applied component-wise (to each dimension separately)

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The activation function g should be non-linear, e.g.:

- Rectified linear unit: $\text{ReLU}(z) = \max(0, z)$
- Hyperbolic tangent: $\tanh(z) = \frac{e^{2z}-1}{e^{2z}+1}$
- Sigmoid: $\sigma(z) = \frac{1}{1+e^{-z}}$

Gradient Descent for Neural Nets

Assume there is a good annotated corpus:

$$\mathcal{D} = \left\{ (x^{(1)}, y^{(1)}), (x^{(2)}, y^{(2)}), \dots, (x^{(l)}, y^{(l)}) \right\}$$

Aim to maximise the log-likelihood (also called “cross entropy”):

$$L(\theta) = \sum_{i=1}^l \log p(y^{(i)} | x^{(i)}; \theta)$$

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θ now contains parameters from many layers,
but we can still use gradient descent:

- Backpropagation: efficient application of chain rule
- Iterate many times over training set

Reading

- D Jurafsky and J Martin. *Speech and Language Processing*
Chapter 6. web.stanford.edu/~jurafsky/slp3/6.pdf
- * Essence of linear algebra
www.youtube.com/watch?v=fNk_zzaMoSs&list=PLZHQObOWTQDPD3MizzM2xVFitgF8hE_ab