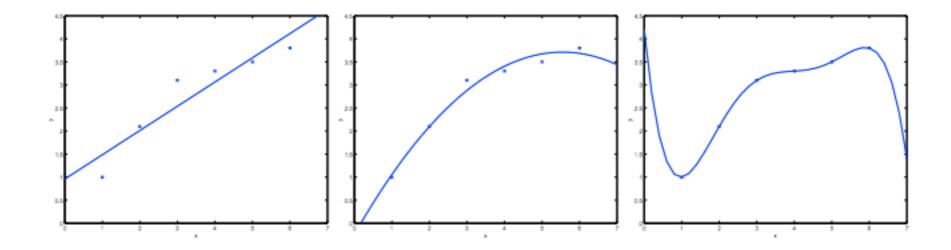
### DAT1

26/3/16

#### Agenda

- Learning Theory
- Reinforcement Learning Intro
- Spark
- (Recommender Practice)
- Random Topics

Bias / Variance: parameters, models



How much data needed?

Which models to even consider?

**Lemma.** (The union bound). Let  $A_1, A_2, \ldots, A_k$  be k different events (that may not be independent). Then

$$P(A_1 \cup \cdots \cup A_k) \leq P(A_1) + \ldots + P(A_k).$$

**Lemma.** (Hoeffding inequality) Let  $Z_1, \ldots, Z_m$  be m independent and identically distributed (iid) random variables drawn from a Bernoulli( $\phi$ ) distribution. I.e.,  $P(Z_i = 1) = \phi$ , and  $P(Z_i = 0) = 1 - \phi$ . Let  $\hat{\phi} = (1/m) \sum_{i=1}^m Z_i$  be the mean of these random variables, and let any  $\gamma > 0$  be fixed. Then

$$P(|\phi - \hat{\phi}| > \gamma) \le 2 \exp(-2\gamma^2 m)$$

How many training points (samples) to be within 0.01 of each other with probability 95%?

Binary classification empirical error

$$\hat{\varepsilon}(h) = \frac{1}{m} \sum_{i=1}^{m} 1\{h(x^{(i)}) \neq y^{(i)}\}.$$

(fraction misclassified)

Generalisation error

$$\varepsilon(h) = P_{(x,y)\sim\mathcal{D}}(h(x) \neq y).$$

PAC framework "probably approximately correct"

i.e. training and testing from same distribution

#### Minimising training error

$$\hat{ heta} = rg \min_{ heta} \hat{arepsilon}(h_{ heta}).$$

$$\mathcal{H} = \{h_{\theta} : h_{\theta}(x) = 1\{\theta^{T}x \geq 0\}, \theta \in \mathbb{R}^{n+1}\}$$

equivalent to 
$$\hat{h} = \arg \min_{h \in \mathcal{H}} \hat{\varepsilon}(h)$$

Finite 
$$\mathcal{H} = \{h_1, \dots, h_k\}$$

We want guarantees on generalisation error when we minimise empirical error

- theoretical framework
- applied cases e.g. too few data points for cross-validation

- Take fixed h\_i in H

- Let 
$$Z_j = 1\{h_i(x^{(j)}) \neq y^{(j)}\}$$

- Training error of h\_i

$$\hat{arepsilon}(h_i) = rac{1}{m} \sum_{j=1}^m Z_j.$$

- Hoeffding

$$P(|\varepsilon(h_i) - \hat{\varepsilon}(h_i)| > \gamma) \le 2\exp(-2\gamma^2 m).$$

Great. But we don't just want this for one h\_i, we want it for all of H simultaneously

$$P(|arepsilon(h_i) - \hat{arepsilon}(h_i)| > \gamma) \leq 2\exp(-2\gamma^2 m).$$
  $P(A_i)$ 

$$P(\exists h \in \mathcal{H}.|\varepsilon(h_i) - \hat{\varepsilon}(h_i)| > \gamma) = P(A_1 \cup \dots \cup A_k)$$

$$\leq \sum_{i=1}^k P(A_i)$$

$$\leq \sum_{i=1}^k 2 \exp(-2\gamma^2 m)$$

$$= 2k \exp(-2\gamma^2 m)$$

$$P(\exists h \in \mathcal{H}.|\varepsilon(h_i) - \hat{\varepsilon}(h_i)| > \gamma) = P(A_1 \cup \dots \cup A_k)$$

$$\leq \sum_{i=1}^k P(A_i)$$

$$\leq \sum_{i=1}^k 2 \exp(-2\gamma^2 m)$$

$$= 2k \exp(-2\gamma^2 m)$$

If we subtract both sides from 1, we find that

$$P(\neg \exists h \in \mathcal{H}. |\varepsilon(h_i) - \hat{\varepsilon}(h_i)| > \gamma) = P(\forall h \in \mathcal{H}. |\varepsilon(h_i) - \hat{\varepsilon}(h_i)| \le \gamma)$$
  
 
$$\ge 1 - 2k \exp(-2\gamma^2 m)$$

#### Example

For instance, we can ask the following question: Given  $\gamma$  and some  $\delta > 0$ , how large must m be before we can guarantee that with probability at least  $1 - \delta$ , training error will be within  $\gamma$  of generalization error? By setting  $\delta = 2k \exp(-2\gamma^2 m)$  and solving for m, [you should convince yourself this is the right thing to do!], we find that if

$$m \ge \frac{1}{2\gamma^2} \log \frac{2k}{\delta},$$

then with probability at least  $1 - \delta$ , we have that  $|\varepsilon(h) - \hat{\varepsilon}(h)| \leq \gamma$  for all  $h \in \mathcal{H}$ . (Equivalently, this shows that the probability that  $|\varepsilon(h) - \hat{\varepsilon}(h)| > \gamma$  for some  $h \in \mathcal{H}$  is at most  $\delta$ .) This bound tells us how many training examples we need in order make a guarantee. The training set size m that a certain method or algorithm requires in order to achieve a certain level of performance is also called the algorithm's **sample complexity**.

With just a bit more effort, we can get an explicit guarantee on generalisation error

**Theorem.** Let  $|\mathcal{H}| = k$ , and let any  $m, \delta$  be fixed. Then with probability at least  $1 - \delta$ , we have that

$$\varepsilon(\hat{h}) \le \left(\min_{h \in \mathcal{H}} \varepsilon(h)\right) + 2\sqrt{\frac{1}{2m} \log \frac{2k}{\delta}}.$$

lower bias: increasing H class

variance tradeoff of increasing H class

In summary, given a hypothesis class of complexity k, we can figure out how many training examples we need to be within a certain distance from minimising generalisation error when carrying out empirical risk minimisation

Analog for continuous case: Vapnik-Chervonenki (VC) dimension

Previously, y had a "correct" value

Now instead, what if the best we can provide is a reward function

e.g. robot motion, gaming...trading?

https://www.youtube.com/watch?v=qv6UVOQ0F44 https://www.youtube.com/watch?v=M8YjvHYbZ9w https://www.youtube.com/watch?v=rVlhMGQgDkY

A Markov decision process is a tuple  $(S, A, \{P_{sa}\}, \gamma, R)$ , where:

- S is a set of states. (For example, in autonomous helicopter flight, S
  might be the set of all possible positions and orientations of the helicopter.)
- A is a set of actions. (For example, the set of all possible directions in which you can push the helicopter's control sticks.)
- $P_{sa}$  are the state transition probabilities. For each state  $s \in S$  and action  $a \in A$ ,  $P_{sa}$  is a distribution over the state space. We'll say more about this later, but briefly,  $P_{sa}$  gives the distribution over what states we will transition to if we take action a in state s.
- $\gamma \in [0, 1)$  is called the **discount factor**.
- R: S × A → R is the reward function. (Rewards are sometimes also written as a function of a state S only, in which case we would have R: S → R).

$$s_0 \xrightarrow{a_0} s_1 \xrightarrow{a_1} s_2 \xrightarrow{a_2} s_3 \xrightarrow{a_3} \dots$$

$$s_1 \sim P_{s_0 a_0}$$

Upon visiting the sequence of states  $s_0, s_1, \ldots$  with actions  $a_0, a_1, \ldots$ , our total payoff is given by

$$R(s_0, a_0) + \gamma R(s_1, a_1) + \gamma^2 R(s_2, a_2) + \cdots$$

Or, when we are writing rewards as a function of the states only, this becomes

$$R(s_0) + \gamma R(s_1) + \gamma^2 R(s_2) + \cdots$$

We want to maximise expected discounted reward

Economic interpretation of gamma?

$$\mathrm{E}\left[R(s_0) + \gamma R(s_1) + \gamma^2 R(s_2) + \cdots\right]$$

#### **Policy** to choose actions

$$\pi: S \mapsto A$$

$$a=\pi(s)$$

#### Value function based on policy

$$V^{\pi}(s) = \mathrm{E}\left[R(s_0) + \gamma R(s_1) + \gamma^2 R(s_2) + \cdots \mid s_0 = s, \pi\right].$$

$$V^{\pi}(s) = \mathrm{E}\left[R(s_0) + \gamma R(s_1) + \gamma^2 R(s_2) + \cdots \mid s_0 = s, \pi\right].$$

#### **Bellman equation**

$$V^{\pi}(s) = R(s) + \gamma \sum_{s' \in S} P_{s\pi(s)}(s') V^{\pi}(s').$$

Can you draw this (perhaps as a decision tree?)

We want to find the optimal value

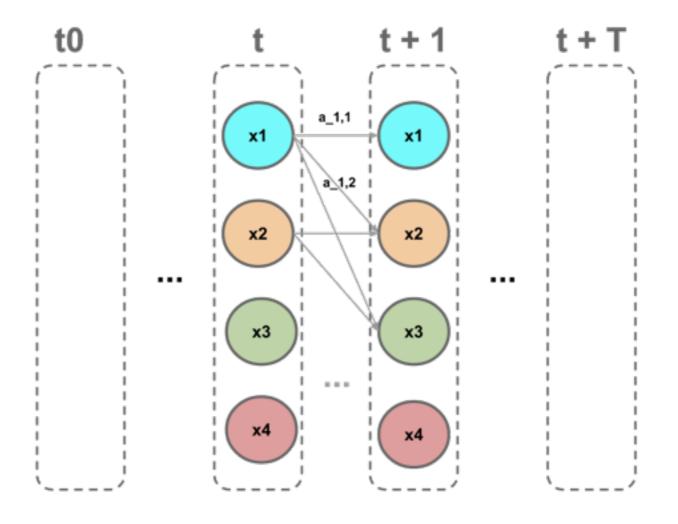
$$V^*(s) = \max_{\pi} V^{\pi}(s).$$

or

$$V^*(s) = R(s) + \max_{a \in A} \gamma \sum_{s' \in S} P_{sa}(s') V^*(s').$$

max over possible actions

Exponential paths to explore



Hansel and Grettel!

Recommended for Dynamic Programming and Value Iteration:

https://en.wikipedia.org/wiki/Memoization

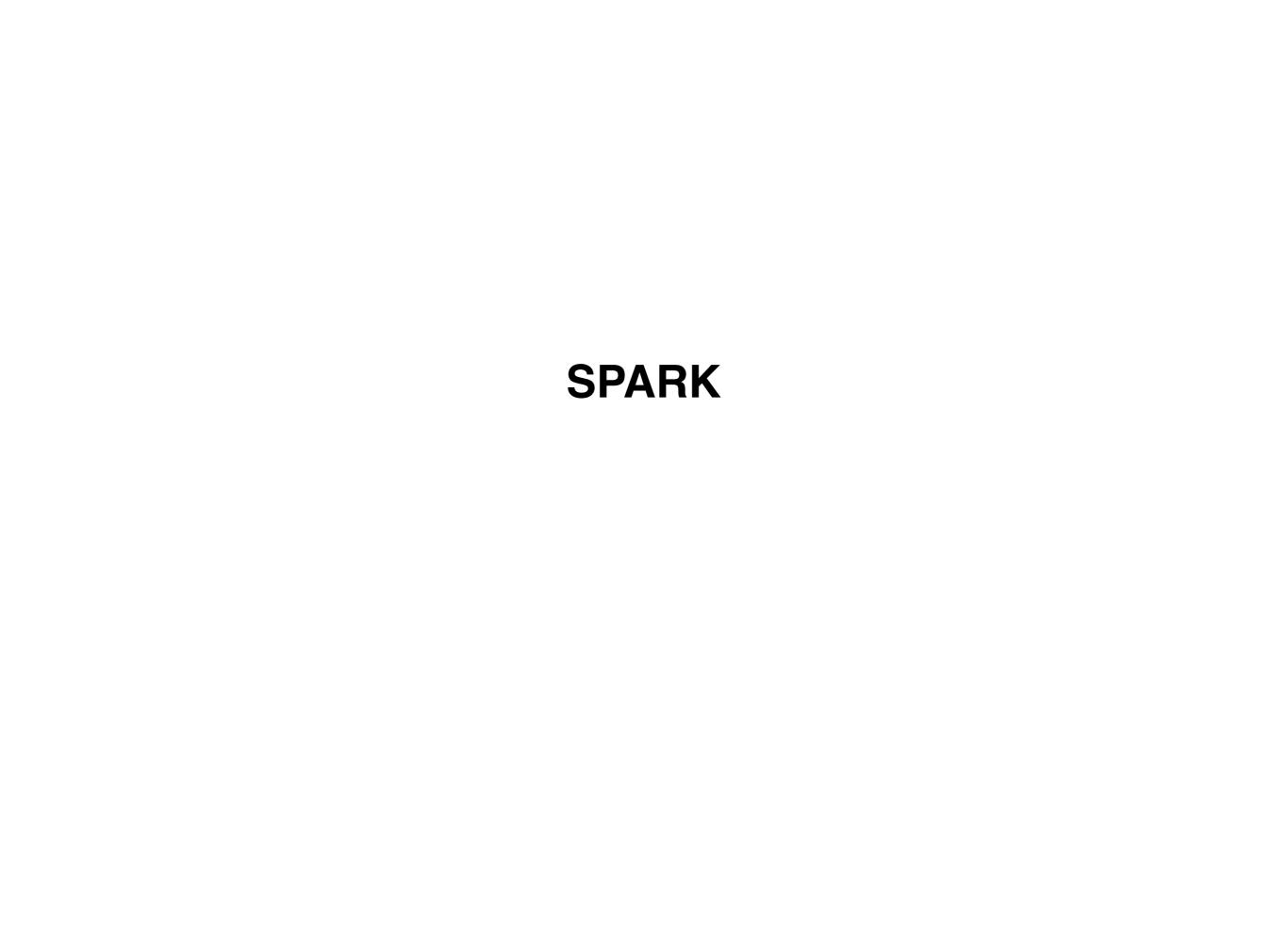
https://www.youtube.com/watch?v=oefOCk3koZo

https://www.youtube.com/watch?v=ip4iSMRW5X4

Extensions:

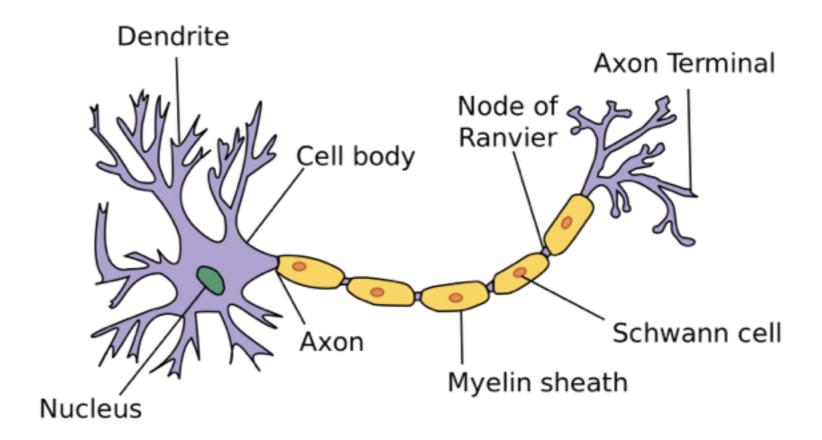
Learning P\_sa

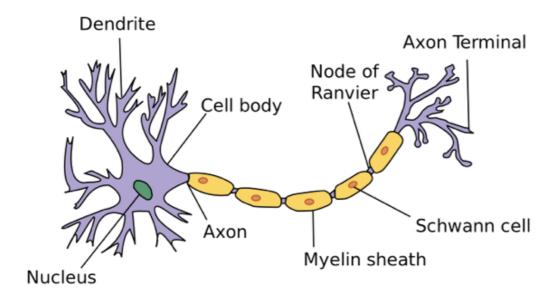
Continuous space (discretisation)

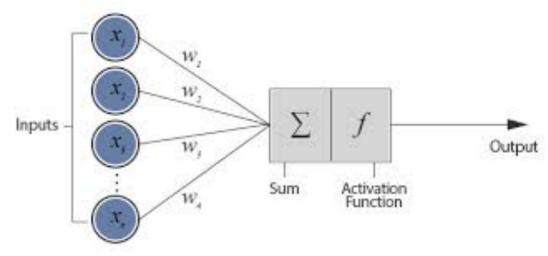


A computational system comprised of layers and each layer is built of interconnected perceptrons

Built to model the animal nervous system?



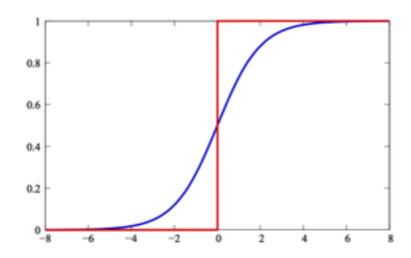




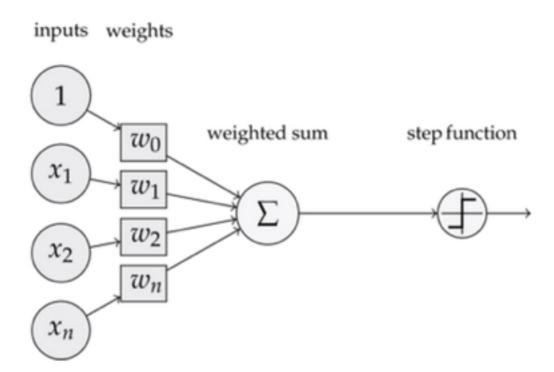
#### Single Perceptron

$$f_{log}(z) = \frac{1}{1 + e^{-z}}$$

 $f_{log}$  is called logistic function

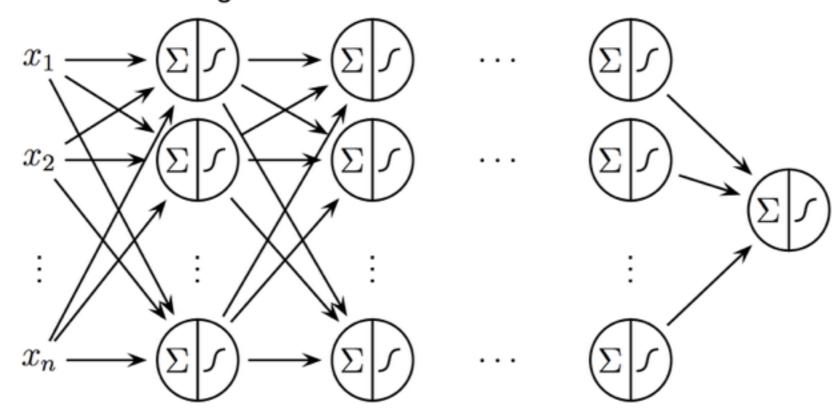


Takes in input and uses an activation function in order to output



# Artificial Neural Networks are also known as multi layer perceptrons

A multi layer perceptrons (MLP) is a finite acyclic graph. The nodes are neurons with logistic activation.



But how does it learn?!

Back-Propagation

https://en.wikipedia.org/wiki/Backpropagation

Pros Cons

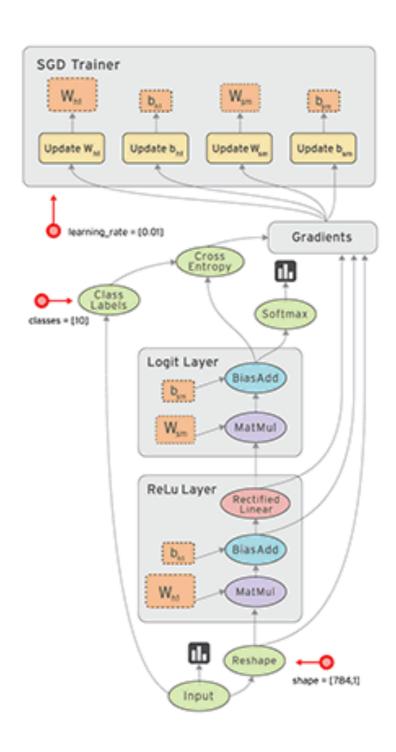
- Online model (updates as you go)
  - Doesn't need to be fit all of the time
- Very fast predictions
- Can approximate almost any type of function
- Can be used in a supervised and unsupervised manner
- Super cool

- Requires many training samples to be considered good
- Hard to describe what is happening
- Requires a lot of hardware / computation power
- Slow to train
- Can be difficult to use

https://www.tensorflow.org/

# TensorFlow is an Open Source Software Library for Machine Intelligence

**GET STARTED** 



# Random Topics Web Scraping

Python Beautiful Soup
<a href="http://www.crummy.com/software/">http://www.crummy.com/software/</a>
<a href="mailto:BeautifulSoup/bs4/doc/">BeautifulSoup/bs4/doc/</a>

Exercise: scrape this

list of buyers <a href="http://econpy.pythonanywhere.com/ex/001.html">http://econpy.pythonanywhere.com/ex/001.html</a>

finance http://stackoverflow.com/questions/10040954/ alternative-to-google-finance-api

# Random Topics Web Scraping

#### Concepts

- Robustness: changes in the web page/api
- Scale: massive web scraping fault tolerance
- Integrity: external data sources

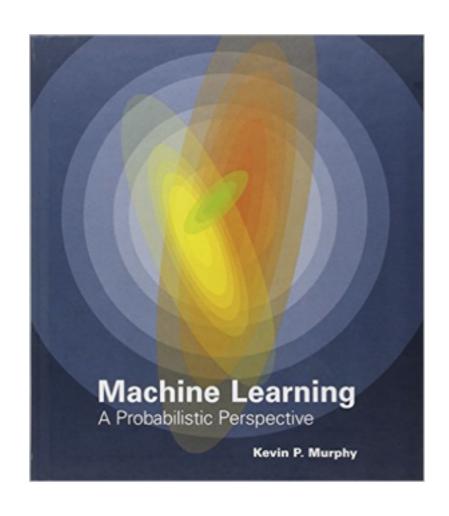
# Random Topics Setting up servers

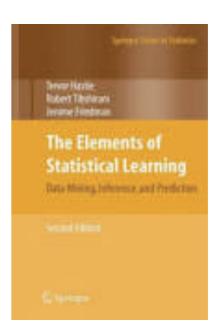




- 1. AWS account
- 2. ec2 instance
- 3. ssh into it

- 1. heroku account
- 2. python "hello" api
- 3. push to heroku

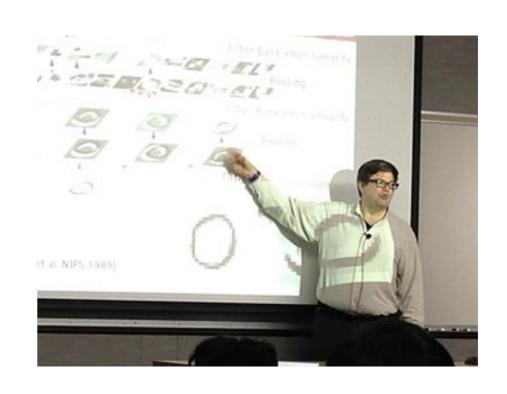




Michael Nielsen, Christopher Olah <a href="http://colah.github.io/">http://colah.github.io/</a>



Geoffrey Hinton



Yann LeCun



Andrej Karpathy



Hugo Larochelle

Linear Algebra

Probability

Parallel Programming

Computer Science

Topology

Information Theory